

**SITE ASSESSMENT FOR PROPOSED COKE
POINT DREDGED MATERIAL CONTAINMENT
FACILITY AT SPARROWS POINT
BALTIMORE COUNTY, MARYLAND**

APPENDIX C

Mass Distribution Calculations

Prepared for:



Maryland Port Administration
2310 Broening Highway
Baltimore, Maryland 21224

Under Contract to:



Maryland Environmental Service
259 Najoles Road
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APPENDIX C. MASS DISTRIBUTION CALCULATIONS

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NAPL RECOVERY TEST CALCULATIONS

Results from LNAPL Bail Down Tests
Wells BP-MW-5, BP-MW-8, and BP-MW-10
Coke Point Peninsula, Sparrows Pt., Maryland, 24 June 2009

OBJECTIVE

The objective of the test was to remove product in the well to a minimal thickness and time the recovery of product to its original thickness.

This provides an estimate of the potential recovery rate using a skimmer or other product recovery tool.

APPROACH

Recovery rate was calculated by determining the time it takes for 80 percent of the original NAPL thickness to return to the well. This was used to calculate volume per time by converting the recovery (linear ft per time) into volume per time using a conversion factor based on the well radius.

Bail down tests were performed on wells BP-MW-5, BP-MW-8, and BP-MW-10 at the Benzol Processing area of the Coke Point Peninsula on 24 June 2009.

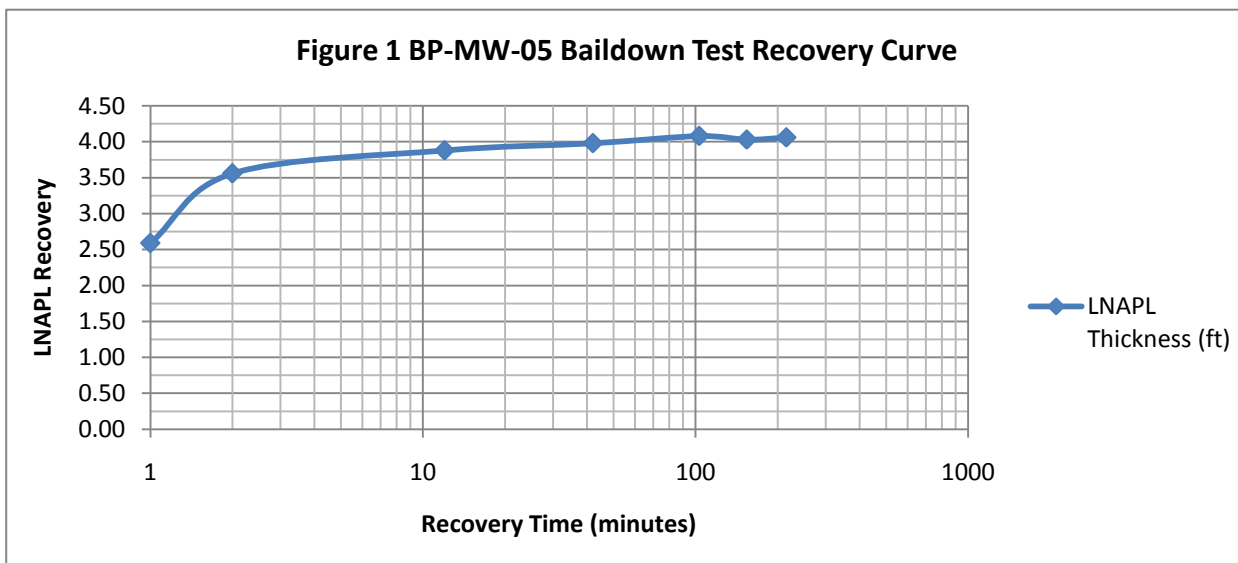
Results from LNAPL Bail Down Tests
 Wells BP-MW-5, BP-MW-8, and BP-MW-10
 Coke Point Peninsula, Sparrows Pt., Maryland, 24 June 2009

Well BP-MW-05

- Initial LNAPL thickness was 4.1 ft.
- Well was bailed for 33 minutes to a thickness of 2.59 ft.

Table 1. LNAPL Recovery Data for Well BP-MW-05

Recovery Time (minutes)	LNAPL Thickness (ft)	Percent Recovery
1	2.59	0.0%
2	3.56	64.2%
12	3.88	85.4%
42	3.98	92.1%
103	4.08	98.7%
154	4.03	95.4%
215	4.06	97.4%



Recovery rate calculation

1. Maximum baildown = 4.1 - 2.59 = 1.51 ft
2. 80% recovery = (0.8 x 1.51 ft) + 2.59 ft = 1.21 ft + 2.59 ft = 3.80 ft
3. Elapsed time to 80% recovery (interpolation from Figure 1) = 6 min
4. Compute average recovery rate (in gal/day) to 80% recovery
 $\pi(\text{well radius in ft})^2 \times (7.48 \text{ gal/ft}^3) = \text{gal per linear ft in a well}$
 $\pi(0.08 \text{ ft})^2 \times (7.48 \text{ gal/ft}^3) = 0.05 \text{ gal per linear ft (2-in. well)}$

$$0.05 \text{ gal/ft} \times (1.21 \text{ ft}) / (6 \text{ min}) = 0.01 \text{ gal/min} = 14.5 \text{ gal/day}$$

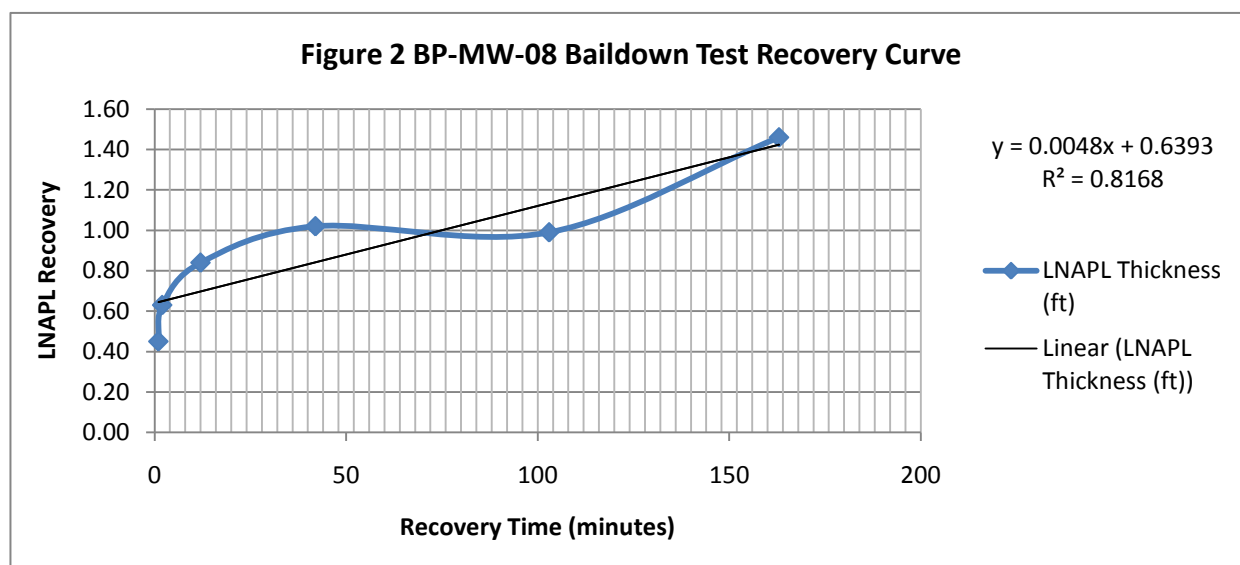
Results from LNAPL Bail Down Tests
Wells BP-MW-5, BP-MW-8, and BP-MW-10
Coke Point Peninsula, Sparrows Pt., Maryland, 24 June 2009

Well BP-MW-08

- Initial LNAPL thickness was 4.49 ft.
- Well was bailed for 8 minutes to a thickness of 0.45 ft.

Table 2. LNAPL Recovery Data for Well BP-MW-08

Recovery Time (minutes)	LNAPL Thickness (ft)	Percent Recovery
1	0.45	0.0%
2	0.63	4.5%
12	0.84	9.7%
42	1.02	14.1%
103	0.99	13.4%
163	1.46	25.0%



Recovery rate calculation

1. Maximum baildown = 4.49 - 0.45 = 4.04 ft
2. 80% recovery = 0.8 x 4.04 = 3.23 ft
3. Compute elapsed time to 80% recovery (linear extrapolation from Figure 2)
 $x = (y - 0.6393)/0.0048 = (3.23 - 0.6393)/0.0048 = 540 \text{ min}$
4. Compute average recovery rate to 80% recovery
 $0.05 \text{ gal/ft} \times (3.23 \text{ ft}) / (540 \text{ min}) = 2.99\text{E-}4 \text{ gal/min} = 0.43 \text{ gal/day}$

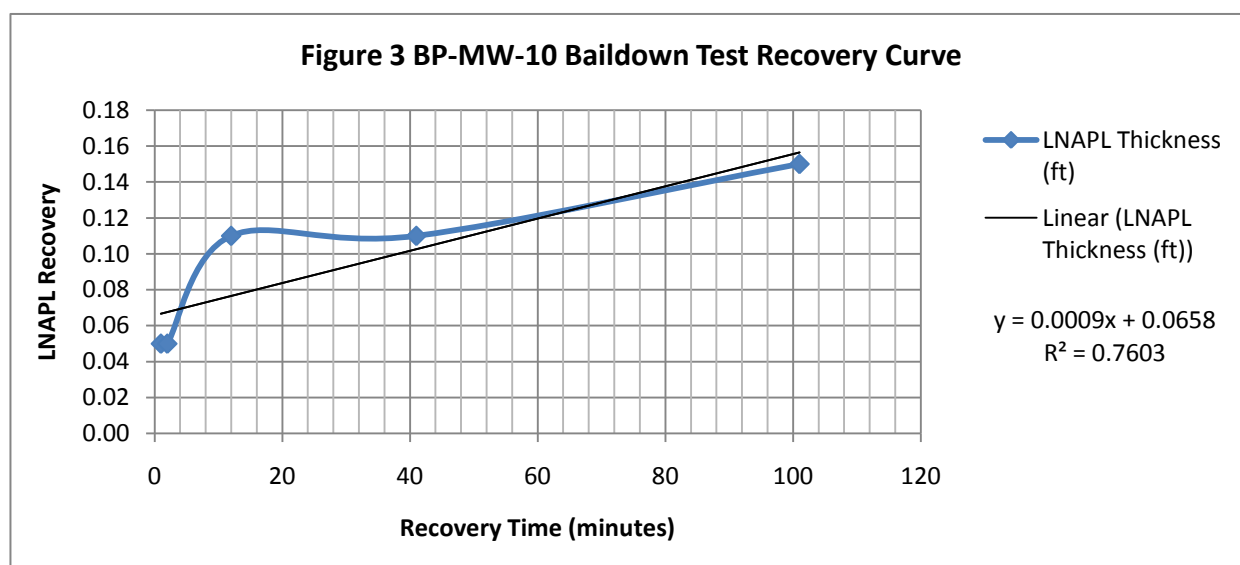
Results from LNAPL Bail Down Tests
Wells BP-MW-5, BP-MW-8, and BP-MW-10
Coke Point Peninsula, Sparrows Pt., Maryland, 24 June 2009

Well BP-MW-10

- Initial LNAPL thickness was 0.70 ft.
- Well was bailed for 4 minutes to a thickness of 0.05 ft.

Table 3. LNAPL Recovery Data for Well BP-MW-10

Recovery Time (minutes)	LNAPL Thickness (ft)	Percent Recovery
1	0.05	0.0%
2	0.05	0.0%
12	0.11	9.2%
41	0.11	9.2%
101	0.15	15.4%



Recovery rate calculation

1. Maximum baildown = 0.70 - 0.05 = 0.65 ft
2. 80% recovery = 0.8 x 0.65 = 0.52 ft
3. Compute elapsed time to 80% recovery (linear extrapolation from Figure 2)
 $x = (y - 0.0658)/0.0009 = (0.52 - 0.0658)/0.0009 = 505 \text{ min}$
4. Compute average recovery rate to 80% recovery
 $0.05 \text{ gal/ft} \times (0.52 \text{ ft}) / (505 \text{ min}) = 5.15\text{E-}5 \text{ gal/min} = 0.07 \text{ gal/day}$

NAPLANAL MODEL

NAPLANAL MODEL INFORMATION

Excerpted from NAPLANAL Version 1.0.0 Readme.TXT, December 10 1997.

Introduction

NAPLANAL is a program developed to estimate the residual NAPL saturation and individual chemical component compositional distribution in a soil sample from the results of soil chemical analysis. The model is especially useful for those who collect soil samples but have been unable to relate the total soil chemical concentrations with fundamental parameters such as NAPL saturation.

The original NAPLANAL code was written using FORTRAN and operates under DOS environment. The current version of NAPLANAL program was rewritten using Visual Basic which provides the end user with a visual interface.

Understanding the Model Output

The output file summarizes the input and output for each sample in a single page. The type of model used (saturated vs. unsaturated, porosity known vs. water content known, etc.), and the values of input parameters except for the chemical concentrations are displayed in the first several lines for verification. In the table of results, the first two columns display the chemicals detected in the sample, and the third column, "total mass (mg/kg)*", displays the input sample chemical concentrations for verification. The asterisk "*" signals that the concentration is normalized by the wet mass of the sample.

The next several columns give the calculated masses of each chemical in each phase, also normalized by the wet mass of the sample. Thus, the sum of the masses in each phase should add up to the total chemical masses displayed in the third column, "total mass (mg/kg)*". These values can be used to determine fractional distribution among phases.

After these columns are columns that give the calculated concentrations of chemicals in each phase normalized by the volume or mass of the phase. These include aqueous concentrations (mg in aqueous phase per liter of water), sorbed concentrations (mg sorbed per kg of dry soil), and concentrations in NAPL (kg in NAPL per liter of NAPL).

The last column contains calculated mole fractions of chemicals in the NAPL phase. The sum of the mole fractions equals one by definition. When no NAPL is calculated to be present in the sample, the mole fractions represent the estimated mole fractions of a NAPL that may have caused the observed chemical concentrations in the sample. Refer to the discussion of the dilution model at the end of this section.

Additional calculations follow the table displaying the calculated chemical component distributions. First, given and calculated phase volumes are presented as fractions of the total sample volume. The sum of the volume fractions of each phase should equal one. Also, the sum of the water, NAPL, and air volume fractions should equal the given or calculated porosity. Second, the calculated densities of the bulk sample and the NAPL itself (if NAPL is present in the sample) are presented. (Note that the calculated NAPL density is an approximation - it is simply the weighted average of the pure component densities.) Finally, if NAPL is present, the NAPL saturation (volume of NAPL per volume of pore space) is presented as a percentage.

If no NAPL is present in the sample, a dilution factor is given.

This dilution factor provides a measure of how much more concentrated the chemicals in the sample must be before a NAPL would be calculated to exist in the sample. A dilution factor of zero and a NAPL

saturation of zero indicates that the sample is at the solubility limit with respect to NAPL without a NAPL present. A dilution factor of 10 implies that the sample theoretically could be reproduced by mixing 10 parts clean soil sample with one part sample at the solubility limit with respect to NAPL. In other words, a dilution factor of 10 implies that the chemical concentrations need to be 11 times their measured values to reach the solubility limit with respect to a NAPL.

Reference

Mariner, P.E., M. Jin, and R.E. Jackson, "Algorithms for the Estimation of NAPL Saturation and Composition from Typical Soil Chemical Analyses." *Ground Water Monitoring and Remediation*, Spring Issue, p. 122. 1997.

Notes

Naphthalene was not detected in sample BH-SED-09-12 and the model was not run for that sample.

The model did not converge for two samples (BP-B02-14 and BP-B04-16) and therefore did not produce an output file. It is believed that the concentrations of naphthalene were too low for the model to process.

**NAPLANAL MODEL
ONSHORE INVESTIGATION OUTPUT FILES**

Sample Name Identification, BP-SO-B01-8

Model used: Liquid saturated & water moisture content known,
 Water moisture content (Volume Frac.), 0.2200,

NAPLANAL ANALYSIS RESULTS:

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| ID#, | Name,                 | Total,     | Mass,     | Mass in,   | Mass,      | Conc.,     | Sorbed,    | Conc.,   | Mole fraction, |
|------|-----------------------|------------|-----------|------------|------------|------------|------------|----------|----------------|
| ,    | ,                     | mass,      | in water, | in soil,   | in NAPL,   | in water,  | in soil,   | in NAPL, | in NAPL,       |
| ,    | ,                     | (mg/kg)*,  | (mg/kg)*, | (mg/kg)*,  | (mg/kg)*,  | (mg/L),    | (mg/kg)^,  | (kg/L),  | ,              |
| ===, | =====,                | =====,     | =====,    | =====,     | =====,     | =====,     | =====,     | =====,   | =====,         |
| 29,  | 1-methylnaphthalene,  | 3.7000,    | 0.0009,   | 0.7105,    | 2.9886,    | 0.0093,    | 0.7910,    | 0.0006,  | 0.0004,        |
| 30,  | 2-methylnaphthalene,  | 7.4000,    | 0.0018,   | 1.4032,    | 5.9951,    | 0.0184,    | 1.5622,    | 0.0011,  | 0.0007,        |
| 32,  | acenapthene,          | 0.9400,    | 0.0000,   | 0.0146,    | 0.9254,    | 0.0004,    | 0.0162,    | 0.0002,  | 0.0001,        |
| 31,  | acenapthylene,        | 4.1000,    | 0.0002,   | 0.0404,    | 4.0594,    | 0.0018,    | 0.0450,    | 0.0008,  | 0.0005,        |
| 35,  | anthracene,           | 8.4000,    | 0.0000,   | 0.0046,    | 8.3954,    | 0.0000,    | 0.0051,    | 0.0016,  | 0.0008,        |
| 42,  | benzene,              | 5700.0000, | 160.7603, | 1234.1157, | 4305.1240, | 1655.4107, | 1373.9909, | 0.8026,  | 0.9459,        |
| 39,  | benz(a)anthracene,    | 12.0000,   | 0.0000,   | 0.0634,    | 11.9366,   | 0.0000,    | 0.0706,    | 0.0022,  | 0.0009,        |
| 40,  | benzo(a)pyrene,       | 9.7000,    | 0.0000,   | 0.0390,    | 9.6610,    | 0.0000,    | 0.0434,    | 0.0018,  | 0.0007,        |
| 50,  | benzo(b)fluoranthene, | 12.0000,   | 0.0000,   | 0.0562,    | 11.9438,   | 0.0000,    | 0.0626,    | 0.0022,  | 0.0008,        |
| 51,  | benzo(k)fluoranthene, | 4.0000,    | 0.0000,   | 0.0058,    | 3.9942,    | 0.0000,    | 0.0064,    | 0.0007,  | 0.0003,        |
| 33,  | 9h-fluorene,          | 11.0000,   | 0.0002,   | 0.1244,    | 10.8754,   | 0.0019,    | 0.1385,    | 0.0020,  | 0.0011,        |
| 38,  | chrysene,             | 13.0000,   | 0.0000,   | 0.0032,    | 12.9968,   | 0.0000,    | 0.0035,    | 0.0024,  | 0.0010,        |
| 44,  | ethylbenzene,         | 140.0000,  | 0.2683,   | 27.2960,   | 112.4357,  | 2.7627,    | 30.3898,   | 0.0210,  | 0.0182,        |
| 36,  | fluoranthene,         | 18.0000,   | 0.0000,   | 0.1068,    | 17.8932,   | 0.0003,    | 0.1189,    | 0.0033,  | 0.0015,        |
| 28,  | naphthalene,          | 29.0000,   | 0.0114,   | 1.3689,    | 27.6197,   | 0.1172,    | 1.5240,    | 0.0051,  | 0.0037,        |
| 34,  | phenanthrene,         | 18.0000,   | 0.0002,   | 0.2153,    | 17.7845,   | 0.0017,    | 0.2397,    | 0.0033,  | 0.0017,        |
| 37,  | pyrene,               | 16.0000,   | 0.0000,   | 0.0609,    | 15.9390,   | 0.0002,    | 0.0678,    | 0.0030,  | 0.0014,        |
| 43,  | toluene,              | 140.0000,  | 1.0600,   | 29.4114,   | 109.5287,  | 10.9150,   | 32.7449,   | 0.0204,  | 0.0204,        |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.220000,  
 NAPL Volume Frac.(l/l), 0.012151,  
 Soil Volume Frac.(l/l), 0.767849,  
 Porosity (Volume Frac.), 0.232151,

Bulk Density (kg/l), 2.2654,  
 NAPL Density (kg/l), 0.8744,

NAPL Saturation (%), 5.2341,

Numerical Accuracy Information

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 The solution converged in 46 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
          (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene      142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene      142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene      152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene      228.3000      0.9000      0.0012      0.0057      138000.0000
  40  50-32-8      benzo(a)pyrene      252.3200      0.9000      0.0016      0.0012      550000.0000
  50  205-99-2      benzo(b)fluoranthene      252.3200      0.9000      0.0119      0.0140      55000.0000
  51  207-08-9      benzo(k)fluoranthene      252.3200      0.9000      0.0394      0.0043      55000.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      20000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
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 END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/8/2009 3:47:36 PM

Sample Name Identification, BP-SO-B01-14

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.2500,

NAPLANAL ANALYSIS RESULTS:

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| ID# | Name                  | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|-----------------------|----------------------|-------------------------|------------------------|------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene,  | 1.4000               | 0.0005                  | 0.3134                 | 1.0862                 | 0.0042                  | 0.3547                    | 0.0002                 | 0.0002                  |
| 30  | 2-methylnaphthalene,  | 2.1000               | 0.0007                  | 0.4644                 | 1.6349                 | 0.0062                  | 0.5256                    | 0.0004                 | 0.0002                  |
| 32  | acenapthene,          | 0.7100               | 0.0000                  | 0.0133                 | 0.6967                 | 0.0003                  | 0.0150                    | 0.0002                 | 0.0001                  |
| 31  | acenapthylene,        | 1.5000               | 0.0001                  | 0.0179                 | 1.4820                 | 0.0008                  | 0.0203                    | 0.0003                 | 0.0002                  |
| 35  | anthracene,           | 1.4000               | 0.0000                  | 0.0009                 | 1.3991                 | 0.0000                  | 0.0010                    | 0.0003                 | 0.0002                  |
| 42  | benzene,              | 4300.0000            | 163.9120                | 1067.4555              | 3068.6325              | 1455.6042               | 1208.1515                 | 0.6959                 | 0.8318                  |
| 39  | benz(a)anthracene,    | 2.6000               | 0.0000                  | 0.0167                 | 2.5833                 | 0.0000                  | 0.0188                    | 0.0006                 | 0.0002                  |
| 40  | benzo(a)pyrene,       | 1.8000               | 0.0000                  | 0.0088                 | 1.7912                 | 0.0000                  | 0.0099                    | 0.0004                 | 0.0002                  |
| 50  | benzo(b)fluoranthene, | 3.2000               | 0.0000                  | 0.0182                 | 3.1818                 | 0.0000                  | 0.0206                    | 0.0007                 | 0.0003                  |
| 3   | bromodichloromethane, | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 33  | 9h-fluorene,          | 8.9000               | 0.0002                  | 0.1219                 | 8.7779                 | 0.0019                  | 0.1379                    | 0.0020                 | 0.0011                  |
| 38  | chrysene,             | 2.4000               | 0.0000                  | 0.0007                 | 2.3993                 | 0.0000                  | 0.0008                    | 0.0005                 | 0.0002                  |
| 44  | ethylbenzene,         | 81.0000              | 0.2130                  | 18.3838                | 62.4032                | 1.8915                  | 20.8069                   | 0.0142                 | 0.0124                  |
| 36  | fluoranthene,         | 6.7000               | 0.0000                  | 0.0482                 | 6.6518                 | 0.0001                  | 0.0545                    | 0.0015                 | 0.0007                  |
| 28  | naphthalene,          | 67.0000              | 0.0372                  | 3.7988                 | 63.1640                | 0.3307                  | 4.2995                    | 0.0143                 | 0.0104                  |
| 34  | phenanthrene,         | 5.7000               | 0.0001                  | 0.0825                 | 5.6174                 | 0.0007                  | 0.0934                    | 0.0013                 | 0.0007                  |
| 37  | pyrene,               | 4.5000               | 0.0000                  | 0.0208                 | 4.4792                 | 0.0001                  | 0.0235                    | 0.0010                 | 0.0005                  |
| 43  | toluene,              | 820.0000             | 8.4732                  | 199.4480               | 612.0788               | 75.2454                 | 225.7362                  | 0.1388                 | 0.1406                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.250000,  
NAPL Volume Frac.(l/l), 0.009790,  
Soil Volume Frac.(l/l), 0.740210,  
Porosity (Volume Frac.), 0.259790,

Bulk Density (kg/l), 2.2201,  
NAPL Density (kg/l), 0.8726,

NAPL Saturation (%), 3.7684,

Numerical Accuracy Information

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 The solution converged in 42 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene  142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene  142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene    152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene        178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene           78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene  228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene    252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene  252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane  163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene          228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene       128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene            202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene           92.1400      0.8620      6.3700      535.0000      300.0000
=====
  
```

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/22/2009 2:39:37 PM

Sample Name Identification, BP-SO-B01-20

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.3700,

NAPLANAL ANALYSIS RESULTS:
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL,<br>, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|---------------------------------|
| 29,  | 1-methylnaphthalene,  | 3.7000,                      | 0.0091,                         | 3.4716,                           | 0.2194,                        | 0.0499,                        | 4.2420,                          | 0.0014,                       | 0.0019,                         |
| 30,  | 2-methylnaphthalene,  | 7.8000,                      | 0.0191,                         | 7.3116,                           | 0.4693,                        | 0.1051,                        | 8.9342,                          | 0.0029,                       | 0.0041,                         |
| 32,  | acenaphthene,         | 1.5000,                      | 0.0037,                         | 0.7653,                           | 0.7311,                        | 0.0203,                        | 0.9351,                          | 0.0045,                       | 0.0059,                         |
| 31,  | acenaphthylene,       | 4.2000,                      | 0.0148,                         | 1.6676,                           | 2.5176,                        | 0.0815,                        | 2.0377,                          | 0.0156,                       | 0.0207,                         |
| 35,  | anthracene,           | 11.0000,                     | 0.0006,                         | 0.3850,                           | 10.6144,                       | 0.0034,                        | 0.4704,                          | 0.0658,                       | 0.0747,                         |
| 42,  | benzene,              | 140.0000,                    | 28.3440,                        | 106.0957,                         | 5.5603,                        | 156.1928,                      | 129.6400,                        | 0.0345,                       | 0.0893,                         |
| 39,  | benz(a)anthracene,    | 14.0000,                     | 0.0001,                         | 3.6567,                           | 10.3432,                       | 0.0003,                        | 4.4682,                          | 0.0641,                       | 0.0568,                         |
| 40,  | benzo(a)pyrene,       | 9.8000,                      | 0.0000,                         | 2.0737,                           | 7.7263,                        | 0.0000,                        | 2.5339,                          | 0.0479,                       | 0.0384,                         |
| 50,  | benzo(b)fluoranthene, | 16.0000,                     | 0.0002,                         | 3.8154,                           | 12.1845,                       | 0.0008,                        | 4.6621,                          | 0.0755,                       | 0.0605,                         |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 33,  | 9h-fluorene,          | 8.9000,                      | 0.0117,                         | 3.8427,                           | 5.0457,                        | 0.0643,                        | 4.6954,                          | 0.0313,                       | 0.0381,                         |
| 38,  | chrysene,             | 14.0000,                     | 0.0000,                         | 0.2229,                           | 13.7771,                       | 0.0001,                        | 0.2724,                          | 0.0854,                       | 0.0757,                         |
| 44,  | ethylbenzene,         | 0.9200,                      | 0.0171,                         | 0.8502,                           | 0.0526,                        | 0.0944,                        | 1.0389,                          | 0.0003,                       | 0.0006,                         |
| 36,  | fluoranthene,         | 39.0000,                     | 0.0065,                         | 11.0838,                          | 27.9098,                       | 0.0356,                        | 13.5434,                         | 0.1730,                       | 0.1730,                         |
| 28,  | naphthalene,          | 48.0000,                     | 0.6202,                         | 36.3585,                          | 11.0213,                       | 3.4175,                        | 44.4270,                         | 0.0683,                       | 0.1078,                         |
| 34,  | phenanthrene,         | 36.0000,                     | 0.0254,                         | 16.0546,                          | 19.9199,                       | 0.1401,                        | 19.6174,                         | 0.1234,                       | 0.1401,                         |
| 37,  | pyrene,               | 20.0000,                     | 0.0024,                         | 4.0565,                           | 15.9411,                       | 0.0130,                        | 4.9567,                          | 0.0988,                       | 0.0988,                         |
| 43,  | toluene,              | 20.0000,                     | 1.3084,                         | 17.7013,                          | 0.9903,                        | 7.2098,                        | 21.6295,                         | 0.0061,                       | 0.0135,                         |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.370000,  
NAPL Volume Frac.(l/l), 0.000329,  
Soil Volume Frac.(l/l), 0.629671,  
Porosity (Volume Frac.), 0.370329,

Bulk Density (kg/l), 2.0389,  
NAPL Density (kg/l), 0.8987,

NAPL Saturation (%), 0.0888,

Numerical Accuracy Information

~~~~~  
 The solution converged in 7 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene  142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene  142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene    152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene  228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene    252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene  252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane  163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/28/2009 3:48:52 PM

Sample Name Identification, BP-SO-B02-08

Model used: Unsaturated sample,
Porosity (Volume Frac.), 0.3000,
Water moisture content (Volume Frac.), 0.1100,
Fraction organic carbon (foc), 0.0100,

NAPLANAL ANALYSIS RESULTS:

Table with 12 columns: ID#, Name, Total, Mass, Mass, Mass, Mass, Conc., Sorbed, Conc., Conc., Mole. Rows include various chemical compounds like 1-methylnaphthalene, 2-methylnaphthalene, acenaphthene, etc.

(mg/kg)* --- mg per kg of soil sample (wet soil)
(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.110000,
NAPL Volume Frac.(l/l), 0.000066,
Soil Volume Frac.(l/l), 0.700000,
Air Volume Frac.(l/l), 0.189934,
Porosity (Volume Frac.), 0.300000,

Bulk Density (kg/l), 1.9653,
 NAPL Density (kg/l), 0.8944,
 NAPL Saturation (%), 0.0221,

Numerical Accuracy Information

~~~~~  
 The solution converged in 12 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
29  90-12-0  1-methylnaphthalene  142.2000  0.9000  0.2600  25.8000  8500.0000
30  91-57-6  2-methylnaphthalene  142.2000  0.9000  0.5180  25.4000  8500.0000
32  83-32-9  acenaphthene  154.2100  0.9000  0.0920  3.4200  4600.0000
31  208-96-8  acenaphthylene  152.2000  0.9000  1.4800  3.9300  2500.0000
35  120-12-7  anthracene  178.2400  0.9000  1.0200  0.0450  14000.0000
42  71-43-2  benzene  78.1100  0.8740  5.5900  1750.0000  83.0000
39  56-55-3  benz(a)anthracene  228.3000  0.9000  0.0012  0.0057  138000.0000
40  50-32-8  benzo(a)pyrene  252.3200  0.9000  0.0016  0.0012  550000.0000
50  205-99-2  benzo(b)fluoranthene  252.3200  0.9000  0.0119  0.0140  550000.0000
3  75-27-4  bromodichloromethane  163.8000  1.9700  2.0600  6716.9000  61.0000
33  86-73-7  9h-fluorene  166.2200  0.9000  0.0642  1.6900  7300.0000
38  218-01-9  chrysene  228.3000  0.9000  0.0011  0.0018  200000.0000
44  100-41-4  ethylbenzene  106.1700  0.8670  6.4300  152.0000  1100.0000
36  206-44-0  fluoranthene  202.2600  0.9000  0.0065  0.2060  38000.0000
28  91-20-3  naphthalene  128.1800  0.9000  1.1500  31.7000  1300.0000
34  85-01-8  phenanthrene  178.2400  0.9000  0.1590  1.0000  14000.0000
37  129-00-0  pyrene  202.2600  0.9000  0.0050  0.1320  38000.0000
43  108-88-3  toluene  92.1400  0.8620  6.3700  535.0000  300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/15/2009 8:37:15 AM

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Sample Name Identification, BP-SO-B02-20

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.3300,

NAPLANAL ANALYSIS RESULTS:  
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ID#,	Name,	Total, mass, (mg/kg)*,	Mass, in water, (mg/kg)*,	Mass in, in soil, (mg/kg)*,	Mass, in NAPL, (mg/kg)*,	Conc., in water, (mg/L),	Sorbed, in soil, (mg/kg)^,	Conc., in NAPL, (kg/L),	Mole fraction, in NAPL, ,
==,	====,	====,	====,	====,	====,	====,	====,	====,	====,
29,	1-methylnaphthalene,	11.0000,	0.0216,	9.8942,	1.0842,	0.1381,	11.7366,	0.0051,	0.0054,
30,	2-methylnaphthalene,	23.0000,	0.0452,	20.6557,	2.2991,	0.2883,	24.5021,	0.0109,	0.0113,
32,	acenapthene,	2.0000,	0.0030,	0.7517,	1.2452,	0.0194,	0.8917,	0.0059,	0.0057,
31,	acenapthylene,	1.4000,	0.0029,	0.3862,	1.0110,	0.0183,	0.4581,	0.0048,	0.0047,
35,	anthracene,	4.4000,	0.0001,	0.0901,	4.3097,	0.0008,	0.1069,	0.0204,	0.0170,
42,	benzene,	360.0000,	61.3471,	273.7732,	24.8797,	391.2683,	324.7527,	0.1178,	0.2236,
39,	benz(a)anthracene,	6.3000,	0.0000,	1.0669,	5.2330,	0.0001,	1.2656,	0.0248,	0.0161,
40,	benzo(a)pyrene,	3.4000,	0.0000,	0.4557,	2.9443,	0.0000,	0.5406,	0.0139,	0.0082,
50,	benzo(b)fluoranthene,	6.3000,	0.0000,	0.9636,	5.3363,	0.0002,	1.1431,	0.0253,	0.0148,
3,	bromodichloromethane,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
33,	9h-fluorene,	6.1000,	0.0047,	1.8601,	4.2352,	0.0302,	2.2065,	0.0201,	0.0179,
38,	chrysene,	7.1000,	0.0000,	0.0656,	7.0344,	0.0000,	0.0779,	0.0333,	0.0216,
44,	ethylbenzene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
36,	fluoranthene,	17.0000,	0.0016,	3.1676,	13.8309,	0.0099,	3.7574,	0.0655,	0.0480,
28,	naphthalene,	260.0000,	2.4155,	168.8379,	88.7466,	15.4059,	200.2773,	0.4203,	0.4860,
34,	phenanthrene,	20.0000,	0.0084,	6.3435,	13.6481,	0.0537,	7.5247,	0.0646,	0.0537,
37,	pyrene,	9.6000,	0.0006,	1.2285,	8.3709,	0.0038,	1.4572,	0.0396,	0.0291,
43,	toluene,	58.0000,	3.1024,	50.0427,	4.8549,	19.7871,	59.3612,	0.0230,	0.0370,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.330000,
NAPL Volume Frac.(l/l), 0.000444,
Soil Volume Frac.(l/l), 0.669556,
Porosity (Volume Frac.), 0.330444,

Bulk Density (kg/l), 2.1047,
NAPL Density (kg/l), 0.8955,

NAPL Saturation (%), 0.1345,

Numerical Accuracy Information

~~~~~  
 The solution converged in 6 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene      142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene      142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene      152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene      228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene      252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene      252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane      163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/15/2009 8:43:41 AM

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Sample Name Identification, BP-SO-B02-20-2

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.3300,

NAPLANAL ANALYSIS RESULTS:  
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ID#,	Name,	Total, mass, (mg/kg)*,	Mass, in water, (mg/kg)*,	Mass in, in soil, (mg/kg)*,	Mass, in NAPL, (mg/kg)*,	Conc., in water, (mg/L),	Sorbed, in soil, (mg/kg)^,	Conc., in NAPL, (kg/L),	Mole fraction, in NAPL, ,
29,	1-methylnaphthalene,	10.0000,	0.0197,	8.9936,	0.9867,	0.1255,	10.6684,	0.0047,	0.0049,
30,	2-methylnaphthalene,	24.0000,	0.0472,	21.5512,	2.4016,	0.3008,	25.5643,	0.0114,	0.0118,
32,	acenapthene,	2.2000,	0.0033,	0.8263,	1.3704,	0.0213,	0.9801,	0.0065,	0.0062,
31,	acenapthylene,	1.7000,	0.0035,	0.4685,	1.2280,	0.0222,	0.5558,	0.0058,	0.0057,
35,	anthracene,	4.4000,	0.0001,	0.0900,	4.3098,	0.0008,	0.1068,	0.0204,	0.0170,
42,	benzene,	360.0000,	61.3421,	273.7505,	24.9074,	391.2361,	324.7260,	0.1178,	0.2236,
39,	benz(a)anthracene,	7.2000,	0.0000,	1.2181,	5.9818,	0.0001,	1.4450,	0.0283,	0.0184,
40,	benzo(a)pyrene,	3.9000,	0.0000,	0.5222,	3.3778,	0.0000,	0.6194,	0.0160,	0.0094,
50,	benzo(b)fluoranthene,	6.3000,	0.0000,	0.9627,	5.3373,	0.0002,	1.1419,	0.0252,	0.0148,
3,	bromodichloromethane,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
33,	9h-fluorene,	6.0000,	0.0047,	1.8281,	4.1673,	0.0297,	2.1685,	0.0197,	0.0176,
38,	chrysene,	7.3000,	0.0000,	0.0674,	7.2326,	0.0000,	0.0800,	0.0342,	0.0222,
44,	ethylbenzene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
36,	fluoranthene,	16.0000,	0.0015,	2.9784,	13.0202,	0.0093,	3.5330,	0.0616,	0.0451,
28,	naphthalene,	260.0000,	2.4145,	168.7692,	88.8163,	15.3997,	200.1959,	0.4199,	0.4858,
34,	phenanthrene,	20.0000,	0.0084,	6.3383,	13.6532,	0.0537,	7.5186,	0.0645,	0.0537,
37,	pyrene,	8.9000,	0.0006,	1.1377,	7.7618,	0.0036,	1.3495,	0.0367,	0.0269,
43,	toluene,	58.0000,	3.1021,	50.0377,	4.8602,	19.7851,	59.3553,	0.0230,	0.0370,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.330000,
NAPL Volume Frac.(l/l), 0.000445,
Soil Volume Frac.(l/l), 0.669555,
Porosity (Volume Frac.), 0.330445,

Bulk Density (kg/l), 2.1047,
NAPL Density (kg/l), 0.8955,

NAPL Saturation (%), 0.1347,

Numerical Accuracy Information

~~~~~  
 The solution converged in 6 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene  142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene  142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene    152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene  228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene    252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene  252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane  163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
=====
  
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END OF RECORDS



NAPLANAL Version 1.0.0

Date and Time: 7/22/2009 3:35:58 PM

-----  
Sample Name Identification, BP-SO-B03-4  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1100,  
-----

NAPLANAL ANALYSIS RESULTS:

-----  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.0170, 0.0000, 0.0170, 0.0000, 0.0002, 0.0178, 0.0006,  
30, 2-methylnaphthalene, 0.0270, 0.0000, 0.0270, 0.0000, 0.0003, 0.0282, 0.0010,  
32, acenaphthene, 0.0140, 0.0000, 0.0140, 0.0000, 0.0003, 0.0146, 0.0068,  
31, acenaphthylene, 0.0038, 0.0000, 0.0038, 0.0000, 0.0002, 0.0040, 0.0030,  
35, anthracene, 0.0056, 0.0000, 0.0056, 0.0000, 0.0000, 0.0059, 0.0682,  
42, benzene, 0.0410, 0.0022, 0.0388, 0.0000, 0.0490, 0.0406, 0.0020,  
39, benz(a)anthracene, 0.0280, 0.0000, 0.0280, 0.0000, 0.0000, 0.0293, 0.0273,  
40, benzo(a)pyrene, 0.0470, 0.0000, 0.0470, 0.0000, 0.0000, 0.0492, 0.0546,  
50, benzo(b)fluoranthene, 0.0650, 0.0000, 0.0650, 0.0000, 0.0000, 0.0680, 0.0648,  
3, bromodichloromethane, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
33, 9h-fluorene, 0.0058, 0.0000, 0.0058, 0.0000, 0.0001, 0.0061, 0.0036,  
38, chrysene, 0.0300, 0.0000, 0.0300, 0.0000, 0.0000, 0.0314, 0.6392,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 0.0290, 0.0000, 0.0290, 0.0000, 0.0001, 0.0303, 0.0284,  
28, naphthalene, 0.2200, 0.0008, 0.2192, 0.0000, 0.0176, 0.2294, 0.0408,  
34, phenanthrene, 0.0330, 0.0000, 0.0330, 0.0000, 0.0002, 0.0345, 0.0181,  
37, pyrene, 0.0270, 0.0000, 0.0270, 0.0000, 0.0001, 0.0283, 0.0413,  
43, toluene, 0.0088, 0.0001, 0.0087, 0.0000, 0.0030, 0.0091, 0.0004,  
-----

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.890000,  
Water Volume Frac.(l/l), 0.110000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.110000,

Bulk Density (kg/l), 2.4685,

Dilution Factor (Vol. fac.), 72.2865,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
              (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene      142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene      142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene      152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene      228.3000      0.9000      0.0012      0.0057      138000.0000
  40  50-32-8      benzo(a)pyrene      252.3200      0.9000      0.0016      0.0012      550000.0000
  50  205-99-2      benzo(b)fluoranthene      252.3200      0.9000      0.0119      0.0140      55000.0000
  3  75-27-4      bromodichloromethane      163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      20000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
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 END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/22/2009 3:51:40 PM

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Sample Name Identification, BP-SO-B03-12  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.0980,  
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NAPLANAL ANALYSIS RESULTS:

| ID# | Name                 | Total mass, (mg/kg)* | Mass, in water, (mg/kg)* | Mass, in soil, (mg/kg)* | Mass, in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|--------------------------|-------------------------|-------------------------|-------------------------|---------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 0.0180               | 0.0000                   | 0.0180                  | 0.0000                  | 0.0002                  | 0.0187                    | 0.0012                  |
| 30  | 2-methylnaphthalene  | 0.0380               | 0.0000                   | 0.0380                  | 0.0000                  | 0.0005                  | 0.0395                    | 0.0026                  |
| 32  | acenapthene          | 0.0072               | 0.0000                   | 0.0072                  | 0.0000                  | 0.0002                  | 0.0075                    | 0.0068                  |
| 31  | acenapthylene        | 0.0078               | 0.0000                   | 0.0078                  | 0.0000                  | 0.0003                  | 0.0081                    | 0.0118                  |
| 35  | anthracene           | 0.0120               | 0.0000                   | 0.0120                  | 0.0000                  | 0.0001                  | 0.0125                    | 0.2831                  |
| 42  | benzene              | 0.0720               | 0.0034                   | 0.0686                  | 0.0000                  | 0.0861                  | 0.0714                    | 0.0070                  |
| 39  | benz(a)anthracene    | 0.0073               | 0.0000                   | 0.0073                  | 0.0000                  | 0.0000                  | 0.0076                    | 0.0138                  |
| 40  | benzo(a)pyrene       | 0.0084               | 0.0000                   | 0.0084                  | 0.0000                  | 0.0000                  | 0.0087                    | 0.0189                  |
| 50  | benzo(b)fluoranthene | 0.0130               | 0.0000                   | 0.0130                  | 0.0000                  | 0.0000                  | 0.0135                    | 0.0251                  |
| 3   | bromodichloromethane | 0.0000               | 0.0000                   | 0.0000                  | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                  |
| 33  | 9h-fluorene          | 0.0110               | 0.0000                   | 0.0110                  | 0.0000                  | 0.0002                  | 0.0114                    | 0.0132                  |
| 38  | chrysene             | 0.0078               | 0.0000                   | 0.0078                  | 0.0000                  | 0.0000                  | 0.0081                    | 0.3221                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                   | 0.0000                  | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                  |
| 36  | fluoranthene         | 0.0180               | 0.0000                   | 0.0180                  | 0.0000                  | 0.0000                  | 0.0187                    | 0.0342                  |
| 28  | naphthalene          | 0.4600               | 0.0014                   | 0.4586                  | 0.0000                  | 0.0367                  | 0.4774                    | 0.1654                  |
| 34  | phenanthrene         | 0.0510               | 0.0000                   | 0.0510                  | 0.0000                  | 0.0004                  | 0.0531                    | 0.0541                  |
| 37  | pyrene               | 0.0130               | 0.0000                   | 0.0130                  | 0.0000                  | 0.0000                  | 0.0135                    | 0.0385                  |
| 43  | toluene              | 0.0230               | 0.0003                   | 0.0227                  | 0.0000                  | 0.0079                  | 0.0236                    | 0.0021                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.902000,  
Water Volume Frac.(l/l), 0.098000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.098000,

Bulk Density (kg/l), 2.4883,

Dilution Factor (Vol. fac.), 141.8007,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
          (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene      142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene      142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene      152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene      228.3000      0.9000      0.0012      0.0057      138000.0000
  40  50-32-8      benzo(a)pyrene      252.3200      0.9000      0.0016      0.0012      550000.0000
  50  205-99-2      benzo(b)fluoranthene      252.3200      0.9000      0.0119      0.0140      55000.0000
  3  75-27-4      bromodichloromethane      163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      20000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
=====
  
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 END OF RECORDS

Sample Name Identification, BP-SO-B03-32

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.3000,

NAPLANAL ANALYSIS RESULTS:

Table with 10 columns: ID#, Name, Total mass, Mass in water, Mass in soil, Mass in NAPL, Conc. in water, Sorbed, Conc. in NAPL, Mole fraction in NAPL. Rows list various hydrocarbons like 1-methylnaphthalene, 2-methylnaphthalene, acenaphthene, etc.

(mg/kg)\* --- mg per kg of soil sample (wet soil)
(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.300000,
NAPL Volume Frac.(l/l), 0.000005,
Soil Volume Frac.(l/l), 0.699995,
Porosity (Volume Frac.), 0.300005,

Bulk Density (kg/l), 2.1550,
NAPL Density (kg/l), 0.8994,

NAPL Saturation (%),

0.0017,

Numerical Accuracy Information

The solution converged in 7 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenapthene          | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenapthylene        | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/15/2009 1:24:42 PM

-----  
Sample Name Identification, BP-SO-B04-10  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1800,  
-----

NAPLANAL ANALYSIS RESULTS:

| ID# | Name                 | Total,<br>mass,<br>(mg/kg)* | Mass,<br>in water,<br>(mg/kg)* | Mass,<br>in soil,<br>(mg/kg)* | Mass,<br>in NAPL,<br>(mg/kg)* | Conc.,<br>in water,<br>(mg/L) | Sorbed,<br>in soil,<br>(mg/kg)^ | Mole fraction,<br>in NAPL,<br>, |
|-----|----------------------|-----------------------------|--------------------------------|-------------------------------|-------------------------------|-------------------------------|---------------------------------|---------------------------------|
| 29  | 1-methylnaphthalene  | 0.3300                      | 0.0003                         | 0.3297                        | 0.0000                        | 0.0042                        | 0.3570                          | 0.0013                          |
| 30  | 2-methylnaphthalene  | 0.6400                      | 0.0006                         | 0.6394                        | 0.0000                        | 0.0081                        | 0.6923                          | 0.0026                          |
| 32  | acenapthene          | 0.1600                      | 0.0003                         | 0.1597                        | 0.0000                        | 0.0038                        | 0.1729                          | 0.0090                          |
| 31  | acenapthylene        | 0.0290                      | 0.0001                         | 0.0289                        | 0.0000                        | 0.0013                        | 0.0313                          | 0.0026                          |
| 35  | anthracene           | 0.0610                      | 0.0000                         | 0.0610                        | 0.0000                        | 0.0005                        | 0.0660                          | 0.0858                          |
| 42  | benzene              | 79.0000                     | 7.1688                         | 71.8312                       | 0.0000                        | 93.7124                       | 77.7813                         | 0.4386                          |
| 39  | benz(a)anthracene    | 0.0870                      | 0.0000                         | 0.0870                        | 0.0000                        | 0.0000                        | 0.0942                          | 0.0098                          |
| 40  | benzo(a)pyrene       | 0.0600                      | 0.0000                         | 0.0600                        | 0.0000                        | 0.0000                        | 0.0650                          | 0.0081                          |
| 50  | benzo(b)fluoranthene | 0.1000                      | 0.0000                         | 0.1000                        | 0.0000                        | 0.0000                        | 0.1083                          | 0.0115                          |
| 3   | bromodichloromethane | 0.0000                      | 0.0000                         | 0.0000                        | 0.0000                        | 0.0000                        | 0.0000                          | 0.0000                          |
| 33  | 9h-fluorene          | 0.0000                      | 0.0000                         | 0.0000                        | 0.0000                        | 0.0000                        | 0.0000                          | 0.0000                          |
| 38  | chrysene             | 0.0750                      | 0.0000                         | 0.0750                        | 0.0000                        | 0.0000                        | 0.0812                          | 0.1848                          |
| 44  | ethylbenzene         | 0.0000                      | 0.0000                         | 0.0000                        | 0.0000                        | 0.0000                        | 0.0000                          | 0.0000                          |
| 36  | fluoranthene         | 0.2000                      | 0.0000                         | 0.2000                        | 0.0000                        | 0.0006                        | 0.2165                          | 0.0227                          |
| 28  | naphthalene          | 7.3000                      | 0.0462                         | 7.2538                        | 0.0000                        | 0.6042                        | 7.8546                          | 0.1561                          |
| 34  | phenanthrene         | 0.3100                      | 0.0002                         | 0.3098                        | 0.0000                        | 0.0024                        | 0.3355                          | 0.0196                          |
| 37  | pyrene               | 0.1500                      | 0.0000                         | 0.1500                        | 0.0000                        | 0.0004                        | 0.1624                          | 0.0265                          |
| 43  | toluene              | 3.9000                      | 0.1048                         | 3.7952                        | 0.0000                        | 1.3699                        | 4.1096                          | 0.0210                          |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.820000,  
Water Volume Frac.(l/l), 0.180000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.180000,

Bulk Density (kg/l), 2.3530,

Dilution Factor (Vol. fac.), 7.1903,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
29  90-12-0      1-methylnaphthalene      142.2000      0.9000      0.2600      25.8000      8500.0000
30  91-57-6      2-methylnaphthalene      142.2000      0.9000      0.5180      25.4000      8500.0000
32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
31  208-96-8      acenaphthylene      152.2000      0.9000      1.4800      3.9300      2500.0000
35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
39  56-55-3      benz(a)anthracene      228.3000      0.9000      0.0012      0.0057      138000.0000
40  50-32-8      benzo(a)pyrene      252.3200      0.9000      0.0016      0.0012      550000.0000
50  205-99-2      benzo(b)fluoranthene      252.3200      0.9000      0.0119      0.0140      55000.0000
3  75-27-4      bromodichloromethane      163.8000      1.9700      2.0600      6716.9000      61.0000
33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      20000.0000
44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
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 END OF RECORDS



NAPLANAL Version 1.0.0

Date and Time: 7/15/2009 9:05:43 AM

-----  
Sample Name Identification, BP-SO-B04-24

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.2500,

NAPLANAL ANALYSIS RESULTS:  
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL,<br>, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|---------------------------------|
| ==,  | ====,                 | ====,                        | ====,                           | ====,                             | ====,                          | ====,                          | ====,                            | ====,                         | ====,                           |
| 29,  | 1-methylnaphthalene,  | 1.1000,                      | 0.0016,                         | 1.0930,                           | 0.0054,                        | 0.0145,                        | 1.2305,                          | 0.0004,                       | 0.0006,                         |
| 30,  | 2-methylnaphthalene,  | 3.0000,                      | 0.0044,                         | 2.9806,                           | 0.0150,                        | 0.0395,                        | 3.3556,                          | 0.0010,                       | 0.0016,                         |
| 32,  | acenapthene,          | 1.7000,                      | 0.0043,                         | 1.5776,                           | 0.1181,                        | 0.0386,                        | 1.7761,                          | 0.0079,                       | 0.0113,                         |
| 31,  | acenapthylene,        | 1.6000,                      | 0.0072,                         | 1.4244,                           | 0.1685,                        | 0.0641,                        | 1.6036,                          | 0.0112,                       | 0.0163,                         |
| 35,  | anthracene,           | 3.2000,                      | 0.0009,                         | 1.0124,                           | 2.1867,                        | 0.0081,                        | 1.1397,                          | 0.1457,                       | 0.1809,                         |
| 42,  | benzene,              | 38.0000,                     | 4.9834,                         | 32.8816,                          | 0.1350,                        | 44.6004,                       | 37.0183,                         | 0.0090,                       | 0.0255,                         |
| 39,  | benz(a)anthracene,    | 4.2000,                      | 0.0000,                         | 3.4381,                           | 0.7618,                        | 0.0003,                        | 3.8707,                          | 0.0508,                       | 0.0492,                         |
| 40,  | benzo(a)pyrene,       | 3.6000,                      | 0.0000,                         | 2.7866,                           | 0.8134,                        | 0.0001,                        | 3.1372,                          | 0.0542,                       | 0.0475,                         |
| 50,  | benzo(b)fluoranthene, | 5.6000,                      | 0.0001,                         | 4.4793,                           | 1.1206,                        | 0.0009,                        | 5.0428,                          | 0.0747,                       | 0.0655,                         |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 33,  | 9h-fluorene,          | 3.0000,                      | 0.0047,                         | 2.7159,                           | 0.2794,                        | 0.0419,                        | 3.0576,                          | 0.0186,                       | 0.0248,                         |
| 38,  | chrysene,             | 4.0000,                      | 0.0000,                         | 0.6847,                           | 3.3152,                        | 0.0004,                        | 0.7709,                          | 0.2210,                       | 0.2141,                         |
| 44,  | ethylbenzene,         | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 36,  | fluoranthene,         | 9.9000,                      | 0.0027,                         | 8.2666,                           | 1.6307,                        | 0.0245,                        | 9.3066,                          | 0.1087,                       | 0.1189,                         |
| 28,  | naphthalene,          | 19.0000,                     | 0.1779,                         | 18.3855,                          | 0.4366,                        | 1.5922,                        | 20.6985,                         | 0.0291,                       | 0.0502,                         |
| 34,  | phenanthrene,         | 11.0000,                     | 0.0090,                         | 10.0173,                          | 0.9737,                        | 0.0806,                        | 11.2776,                         | 0.0649,                       | 0.0806,                         |
| 37,  | pyrene,               | 6.5000,                      | 0.0016,                         | 4.9687,                           | 1.5296,                        | 0.0147,                        | 5.5938,                          | 0.1019,                       | 0.1115,                         |
| 43,  | toluene,              | 2.3000,                      | 0.0922,                         | 2.1982,                           | 0.0096,                        | 0.8249,                        | 2.4747,                          | 0.0006,                       | 0.0015,                         |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.250000,  
NAPL Volume Frac.(l/l), 0.000034,  
Soil Volume Frac.(l/l), 0.749966,  
Porosity (Volume Frac.), 0.250034,

Bulk Density (kg/l), 2.2374,  
NAPL Density (kg/l), 0.8997,

NAPL Saturation (%), 0.0134,

Numerical Accuracy Information

~~~~~  
 The solution converged in 6 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene  142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene  142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene    152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene  228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene    252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene  252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane  163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
=====
  
```

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/14/2009 12:39:19 PM

Sample Name Identification, BP-SO-B05-8

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.1000,

NAPLANAL ANALYSIS RESULTS:
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|----------------------------|
| ==,  | ====,                 | ====,                        | ====,                           | ====,                             | ====,                          | ====,                          | ====,                            | ====,                         | ====,                      |
| 29,  | 1-methylnaphthalene,  | 24.0000,                     | 0.0071,                         | 14.2432,                          | 9.7498,                        | 0.1748,                        | 14.8598,                         | 0.0075,                       | 0.0068,                    |
| 30,  | 2-methylnaphthalene,  | 63.0000,                     | 0.0184,                         | 37.1506,                          | 25.8310,                       | 0.4560,                        | 38.7589,                         | 0.0198,                       | 0.0180,                    |
| 32,  | acenapthene,          | 4.6000,                      | 0.0004,                         | 0.4053,                           | 4.1943,                        | 0.0092,                        | 0.4229,                          | 0.0032,                       | 0.0027,                    |
| 31,  | acenapthylene,        | 22.0000,                     | 0.0021,                         | 1.2676,                           | 20.7302,                       | 0.0529,                        | 1.3225,                          | 0.0159,                       | 0.0135,                    |
| 35,  | anthracene,           | 2.2000,                      | 0.0000,                         | 0.0073,                           | 2.1927,                        | 0.0001,                        | 0.0077,                          | 0.0017,                       | 0.0012,                    |
| 42,  | benzene,              | 470.0000,                    | 14.7233,                        | 290.4106,                         | 164.8660,                      | 365.0394,                      | 302.9827,                        | 0.1267,                       | 0.2086,                    |
| 39,  | benz(a)anthracene,    | 1.1000,                      | 0.0000,                         | 0.0348,                           | 1.0652,                        | 0.0000,                        | 0.0363,                          | 0.0008,                       | 0.0005,                    |
| 40,  | benzo(a)pyrene,       | 0.7800,                      | 0.0000,                         | 0.0189,                           | 0.7611,                        | 0.0000,                        | 0.0197,                          | 0.0006,                       | 0.0003,                    |
| 50,  | benzo(b)fluoranthene, | 1.2000,                      | 0.0000,                         | 0.0337,                           | 1.1663,                        | 0.0000,                        | 0.0352,                          | 0.0009,                       | 0.0005,                    |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 33,  | 9h-fluorene,          | 14.0000,                     | 0.0005,                         | 0.9196,                           | 13.0799,                       | 0.0131,                        | 0.9594,                          | 0.0100,                       | 0.0078,                    |
| 38,  | chrysene,             | 1.3000,                      | 0.0000,                         | 0.0019,                           | 1.2981,                        | 0.0000,                        | 0.0020,                          | 0.0010,                       | 0.0006,                    |
| 44,  | ethylbenzene,         | 17.0000,                     | 0.0388,                         | 10.1543,                          | 6.8068,                        | 0.9631,                        | 10.5939,                         | 0.0052,                       | 0.0063,                    |
| 36,  | fluoranthene,         | 3.7000,                      | 0.0000,                         | 0.1309,                           | 3.5691,                        | 0.0004,                        | 0.1365,                          | 0.0027,                       | 0.0017,                    |
| 28,  | naphthalene,          | 1000.0000,                   | 0.7551,                         | 233.2740,                         | 765.9709,                      | 18.7210,                       | 243.3726,                        | 0.5885,                       | 0.5906,                    |
| 34,  | phenanthrene,         | 14.0000,                     | 0.0003,                         | 0.9695,                           | 13.0302,                       | 0.0072,                        | 1.0115,                          | 0.0100,                       | 0.0072,                    |
| 37,  | pyrene,               | 3.1000,                      | 0.0000,                         | 0.0712,                           | 3.0288,                        | 0.0002,                        | 0.0742,                          | 0.0023,                       | 0.0015,                    |
| 43,  | toluene,              | 330.0000,                    | 2.8571,                         | 203.6953,                         | 123.4475,                      | 70.8378,                       | 212.5134,                        | 0.0948,                       | 0.1324,                    |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.100000,  
NAPL Volume Frac.(l/l), 0.003227,  
Soil Volume Frac.(l/l), 0.896773,  
Porosity (Volume Frac.), 0.103227,

Bulk Density (kg/l), 2.4793,  
NAPL Density (kg/l), 0.8919,

NAPL Saturation (%), 3.1261,

Numerical Accuracy Information

~~~~~  
 The solution converged in 168 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)  (kg/l)  (atm-L/mol)  (mg/L)  (ml/g)
=====
  29  90-12-0  1-methylnaphthalene  142.2000  0.9000  0.2600  25.8000  8500.0000
  30  91-57-6  2-methylnaphthalene  142.2000  0.9000  0.5180  25.4000  8500.0000
  32  83-32-9  acenaphthene  154.2100  0.9000  0.0920  3.4200  4600.0000
  31  208-96-8  acenaphthylene  152.2000  0.9000  1.4800  3.9300  2500.0000
  35  120-12-7  anthracene  178.2400  0.9000  1.0200  0.0450  14000.0000
  42  71-43-2  benzene  78.1100  0.8740  5.5900  1750.0000  83.0000
  39  56-55-3  benz(a)anthracene  228.3000  0.9000  0.0012  0.0057  1380000.0000
  40  50-32-8  benzo(a)pyrene  252.3200  0.9000  0.0016  0.0012  5500000.0000
  50  205-99-2  benzo(b)fluoranthene  252.3200  0.9000  0.0119  0.0140  550000.0000
  3  75-27-4  bromodichloromethane  163.8000  1.9700  2.0600  6716.9000  61.0000
  33  86-73-7  9h-fluorene  166.2200  0.9000  0.0642  1.6900  7300.0000
  38  218-01-9  chrysene  228.3000  0.9000  0.0011  0.0018  200000.0000
  44  100-41-4  ethylbenzene  106.1700  0.8670  6.4300  152.0000  1100.0000
  36  206-44-0  fluoranthene  202.2600  0.9000  0.0065  0.2060  38000.0000
  28  91-20-3  naphthalene  128.1800  0.9000  1.1500  31.7000  1300.0000
  34  85-01-8  phenanthrene  178.2400  0.9000  0.1590  1.0000  14000.0000
  37  129-00-0  pyrene  202.2600  0.9000  0.0050  0.1320  38000.0000
  43  108-88-3  toluene  92.1400  0.8620  6.3700  535.0000  300.0000
=====
  
```

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/8/2009 2:44:13 PM

Sample Name Identification, BP-SO-B05-14

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.3500,

NAPLANAL ANALYSIS RESULTS:
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL,<br>, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|---------------------------------|
| ==,  | ====,                 | ====,                        | ====,                           | ====,                             | ====,                          | ====,                          | ====,                            | ====,                         | ====,                           |
| 29,  | 1-methylnaphthalene,  | 38.0000,                     | 0.0493,                         | 20.5223,                          | 17.4284,                       | 0.2911,                        | 24.7457,                         | 0.0116,                       | 0.0113,                         |
| 30,  | 2-methylnaphthalene,  | 98.0000,                     | 0.1262,                         | 52.5463,                          | 45.3275,                       | 0.7454,                        | 63.3602,                         | 0.0301,                       | 0.0293,                         |
| 32,  | acenapthene,          | 7.5000,                      | 0.0024,                         | 0.5418,                           | 6.9558,                        | 0.0142,                        | 0.6533,                          | 0.0046,                       | 0.0042,                         |
| 31,  | acenapthylene,        | 31.0000,                     | 0.0119,                         | 1.4556,                           | 29.5325,                       | 0.0702,                        | 1.7552,                          | 0.0196,                       | 0.0179,                         |
| 35,  | anthracene,           | 4.4000,                      | 0.0000,                         | 0.0118,                           | 4.3881,                        | 0.0001,                        | 0.0143,                          | 0.0029,                       | 0.0023,                         |
| 42,  | benzene,              | 220.0000,                    | 27.7483,                        | 112.8033,                         | 79.4484,                       | 163.8772,                      | 136.0181,                        | 0.0528,                       | 0.0936,                         |
| 39,  | benz(a)anthracene,    | 2.0000,                      | 0.0000,                         | 0.0513,                           | 1.9487,                        | 0.0000,                        | 0.0618,                          | 0.0013,                       | 0.0008,                         |
| 40,  | benzo(a)pyrene,       | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 50,  | benzo(b)fluoranthene, | 1.1000,                      | 0.0000,                         | 0.0250,                           | 1.0750,                        | 0.0000,                        | 0.0302,                          | 0.0007,                       | 0.0004,                         |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 33,  | 9h-fluorene,          | 25.0000,                     | 0.0037,                         | 1.3406,                           | 23.6557,                       | 0.0221,                        | 1.6165,                          | 0.0157,                       | 0.0131,                         |
| 38,  | chrysene,             | 2.1000,                      | 0.0000,                         | 0.0025,                           | 2.0975,                        | 0.0000,                        | 0.0030,                          | 0.0014,                       | 0.0008,                         |
| 44,  | ethylbenzene,         | 7.7000,                      | 0.0772,                         | 4.1617,                           | 3.4611,                        | 0.4562,                        | 5.0182,                          | 0.0023,                       | 0.0030,                         |
| 36,  | fluoranthene,         | 6.3000,                      | 0.0001,                         | 0.1808,                           | 6.1191,                        | 0.0006,                        | 0.2180,                          | 0.0041,                       | 0.0028,                         |
| 28,  | naphthalene,          | 1300.0000,                   | 4.0117,                         | 255.4320,                         | 1040.5563,                     | 23.6923,                       | 307.9995,                        | 0.6911,                       | 0.7474,                         |
| 34,  | phenanthrene,         | 27.0000,                     | 0.0022,                         | 1.5275,                           | 25.4703,                       | 0.0132,                        | 1.8419,                          | 0.0169,                       | 0.0132,                         |
| 37,  | pyrene,               | 4.9000,                      | 0.0000,                         | 0.0911,                           | 4.8089,                        | 0.0003,                        | 0.1098,                          | 0.0032,                       | 0.0022,                         |
| 43,  | toluene,              | 140.0000,                    | 5.2353,                         | 76.9259,                          | 57.8387,                       | 30.9191,                       | 92.7572,                         | 0.0384,                       | 0.0578,                         |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.350000,  
NAPL Volume Frac.(l/l), 0.003112,  
Soil Volume Frac.(l/l), 0.646888,  
Porosity (Volume Frac.), 0.353112,

Bulk Density (kg/l), 2.0670,  
NAPL Density (kg/l), 0.8966,

NAPL Saturation (%), 0.8814,

Numerical Accuracy Information

~~~~~  
 The solution converged in 105 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene      142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene      142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene      152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene      228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene      252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene      252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane      163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
=====
  
```

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/14/2009 1:07:31 PM

Sample Name Identification, BP-SO-B05-20
Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.3800,

NAPLANAL ANALYSIS RESULTS:

~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
=====  
29, 1-methylnaphthalene, 4.4000, 0.0119, 4.3881, 0.0000, 0.0636, 5.4029, 0.0030,  
30, 2-methylnaphthalene, 11.0000, 0.0298, 10.9702, 0.0000, 0.1589, 13.5074, 0.0075,  
32, acenaphthene, 0.8400, 0.0042, 0.8358, 0.0000, 0.0224, 1.0291, 0.0078,  
31, acenaphthylene, 3.7000, 0.0339, 3.6661, 0.0000, 0.1806, 4.5140, 0.0550,  
35, anthracene, 0.5400, 0.0009, 0.5391, 0.0000, 0.0047, 0.6638, 0.1262,  
42, benzene, 21.0000, 4.5765, 16.4235, 0.0000, 24.3638, 20.2220, 0.0167,  
39, benz(a)anthracene, 0.5100, 0.0000, 0.5100, 0.0000, 0.0000, 0.6279, 0.0096,  
40, benzo(a)pyrene, 0.4500, 0.0000, 0.4500, 0.0000, 0.0000, 0.5541, 0.0101,  
50, benzo(b)fluoranthene, 0.3700, 0.0000, 0.3700, 0.0000, 0.0001, 0.4556, 0.0071,  
3, bromodichloromethane, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
33, 9h-fluorene, 2.8000, 0.0088, 2.7912, 0.0000, 0.0471, 3.4367, 0.0334,  
38, chrysene, 0.5100, 0.0001, 0.5099, 0.0000, 0.0003, 0.6279, 0.2089,  
44, ethylbenzene, 0.5900, 0.0121, 0.5779, 0.0000, 0.0647, 0.7115, 0.0005,  
36, fluoranthene, 1.2000, 0.0007, 1.1993, 0.0000, 0.0039, 1.4766, 0.0226,  
28, naphthalene, 120.0000, 2.0976, 117.9024, 0.0000, 11.1670, 145.1714, 0.4218,  
34, phenanthrene, 3.2000, 0.0053, 3.1947, 0.0000, 0.0281, 3.9336, 0.0336,  
37, pyrene, 0.8900, 0.0005, 0.8895, 0.0000, 0.0029, 1.0952, 0.0261,  
43, toluene, 12.0000, 0.8589, 11.1411, 0.0000, 4.5726, 13.7178, 0.0102,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.620000,  
Water Volume Frac.(l/l), 0.380000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.380000,

Bulk Density (kg/l), 2.0230,

Dilution Factor (Vol. fac.), 0.1975,

No NAPL (i.e., NAPL Saturation Equals Zero)

---

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 138000.0000   |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 550000.0000   |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 55000.0000    |
| 3   | 75-27-4  | bromodichloromethane | 163.8000                    | 1.9700            | 2.0600                          | 6716.9000                  | 61.0000       |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 20000.0000    |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

\*\*\*\*\*  
 END OF RECORDS



NAPLANAL Version 1.0.0

Date and Time: 7/15/2009 2:18:18 PM

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Sample Name Identification, BP-SO-B06-8

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1900,

NAPLANAL ANALYSIS RESULTS:  
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL,<br>, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|---------------------------------|
| 29,  | 1-methylnaphthalene,  | 18.0000,                     | 0.0132,                         | 12.6912,                          | 5.2955,                        | 0.1627,                        | 13.8265,                         | 0.0069,                       | 0.0063,                         |
| 30,  | 2-methylnaphthalene,  | 44.0000,                     | 0.0322,                         | 30.8799,                          | 13.0879,                       | 0.3958,                        | 33.6423,                         | 0.0171,                       | 0.0156,                         |
| 32,  | acenapthene,          | 8.2000,                      | 0.0022,                         | 1.1218,                           | 7.0760,                        | 0.0266,                        | 1.2222,                          | 0.0092,                       | 0.0078,                         |
| 31,  | acenapthylene,        | 8.2000,                      | 0.0027,                         | 0.7474,                           | 7.4500,                        | 0.0326,                        | 0.8142,                          | 0.0097,                       | 0.0083,                         |
| 35,  | anthracene,           | 1.9000,                      | 0.0000,                         | 0.0104,                           | 1.8896,                        | 0.0001,                        | 0.0113,                          | 0.0025,                       | 0.0018,                         |
| 42,  | benzene,              | 440.0000,                    | 32.3677,                        | 302.8363,                         | 104.7960,                      | 397.5025,                      | 329.9271,                        | 0.1369,                       | 0.2271,                         |
| 39,  | benz(a)anthracene,    | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 40,  | benzo(a)pyrene,       | 0.0600,                      | 0.0000,                         | 0.0023,                           | 0.0577,                        | 0.0000,                        | 0.0026,                          | 0.0001,                       | 0.0000,                         |
| 50,  | benzo(b)fluoranthene, | 0.1000,                      | 0.0000,                         | 0.0045,                           | 0.0955,                        | 0.0000,                        | 0.0049,                          | 0.0001,                       | 0.0001,                         |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 33,  | 9h-fluorene,          | 10.0000,                     | 0.0013,                         | 1.0340,                           | 8.9647,                        | 0.0154,                        | 1.1265,                          | 0.0117,                       | 0.0091,                         |
| 38,  | chrysene,             | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 44,  | ethylbenzene,         | 5.5000,                      | 0.0313,                         | 3.8823,                           | 1.5864,                        | 0.3845,                        | 4.2296,                          | 0.0021,                       | 0.0025,                         |
| 36,  | fluoranthene,         | 3.2000,                      | 0.0000,                         | 0.1815,                           | 3.0184,                        | 0.0005,                        | 0.1978,                          | 0.0039,                       | 0.0025,                         |
| 28,  | naphthalene,          | 710.0000,                    | 1.6105,                         | 236.0087,                         | 472.3808,                      | 19.7786,                       | 257.1214,                        | 0.6169,                       | 0.6239,                         |
| 34,  | phenanthrene,         | 12.0000,                     | 0.0008,                         | 1.3053,                           | 10.6939,                       | 0.0102,                        | 1.4221,                          | 0.0140,                       | 0.0102,                         |
| 37,  | pyrene,               | 2.8000,                      | 0.0000,                         | 0.1039,                           | 2.6961,                        | 0.0003,                        | 0.1132,                          | 0.0035,                       | 0.0023,                         |
| 43,  | toluene,              | 170.0000,                    | 3.5933,                         | 121.5162,                         | 44.8905,                       | 44.1289,                       | 132.3867,                        | 0.0586,                       | 0.0825,                         |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.190000,  
NAPL Volume Frac.(l/l), 0.001787,  
Soil Volume Frac.(l/l), 0.808213,  
Porosity (Volume Frac.), 0.191787,

Bulk Density (kg/l), 2.3334,  
NAPL Density (kg/l), 0.8933,

NAPL Saturation (%), 0.9316,

Numerical Accuracy Information

~~~~~  
 The solution converged in 399 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)  (kg/l)  (atm-L/mol)  (mg/L)  (ml/g)
=====
  29  90-12-0  1-methylnaphthalene  142.2000  0.9000  0.2600  25.8000  8500.0000
  30  91-57-6  2-methylnaphthalene  142.2000  0.9000  0.5180  25.4000  8500.0000
  32  83-32-9  acenaphthene  154.2100  0.9000  0.0920  3.4200  4600.0000
  31  208-96-8  acenaphthylene  152.2000  0.9000  1.4800  3.9300  2500.0000
  35  120-12-7  anthracene  178.2400  0.9000  1.0200  0.0450  14000.0000
  42  71-43-2  benzene  78.1100  0.8740  5.5900  1750.0000  83.0000
  39  56-55-3  benz(a)anthracene  228.3000  0.9000  0.0012  0.0057  1380000.0000
  40  50-32-8  benzo(a)pyrene  252.3200  0.9000  0.0016  0.0012  5500000.0000
  50  205-99-2  benzo(b)fluoranthene  252.3200  0.9000  0.0119  0.0140  550000.0000
  3  75-27-4  bromodichloromethane  163.8000  1.9700  2.0600  6716.9000  61.0000
  33  86-73-7  9h-fluorene  166.2200  0.9000  0.0642  1.6900  7300.0000
  38  218-01-9  chrysene  228.3000  0.9000  0.0011  0.0018  200000.0000
  44  100-41-4  ethylbenzene  106.1700  0.8670  6.4300  152.0000  1100.0000
  36  206-44-0  fluoranthene  202.2600  0.9000  0.0065  0.2060  38000.0000
  28  91-20-3  naphthalene  128.1800  0.9000  1.1500  31.7000  1300.0000
  34  85-01-8  phenanthrene  178.2400  0.9000  0.1590  1.0000  14000.0000
  37  129-00-0  pyrene  202.2600  0.9000  0.0050  0.1320  38000.0000
  43  108-88-3  toluene  92.1400  0.8620  6.3700  535.0000  300.0000
=====
  
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/15/2009 2:34:19 PM

Sample Name Identification, BP-SO-B06-12

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.2600,

NAPLANAL ANALYSIS RESULTS:
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL,<br>, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|---------------------------------|
| 29,  | 1-methylnaphthalene,  | 16.0000,                     | 0.0102,                         | 6.4732,                           | 9.5166,                        | 0.0865,                        | 7.3511,                          | 0.0044,                       | 0.0034,                         |
| 30,  | 2-methylnaphthalene,  | 40.0000,                     | 0.0252,                         | 16.0329,                          | 23.9420,                       | 0.2142,                        | 18.2071,                         | 0.0111,                       | 0.0084,                         |
| 32,  | acenapthene,          | 2.0000,                      | 0.0002,                         | 0.0861,                           | 1.9136,                        | 0.0021,                        | 0.0978,                          | 0.0009,                       | 0.0006,                         |
| 31,  | acenapthylene,        | 8.4000,                      | 0.0012,                         | 0.2325,                           | 8.1663,                        | 0.0106,                        | 0.2640,                          | 0.0038,                       | 0.0027,                         |
| 35,  | anthracene,           | 1.2000,                      | 0.0000,                         | 0.0019,                           | 1.1981,                        | 0.0000,                        | 0.0021,                          | 0.0006,                       | 0.0003,                         |
| 42,  | benzene,              | 1500.0000,                   | 101.3321,                       | 630.2439,                         | 768.4240,                      | 862.3059,                      | 715.7139,                        | 0.3549,                       | 0.4927,                         |
| 39,  | benz(a)anthracene,    | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 40,  | benzo(a)pyrene,       | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 50,  | benzo(b)fluoranthene, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 33,  | 9h-fluorene,          | 9.1000,                      | 0.0005,                         | 0.2884,                           | 8.8110,                        | 0.0045,                        | 0.3276,                          | 0.0041,                       | 0.0027,                         |
| 38,  | chrysene,             | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 44,  | ethylbenzene,         | 29.0000,                     | 0.1435,                         | 11.8280,                          | 17.0285,                       | 1.2211,                        | 13.4320,                         | 0.0079,                       | 0.0080,                         |
| 36,  | fluoranthene,         | 2.1000,                      | 0.0000,                         | 0.0352,                           | 2.0647,                        | 0.0001,                        | 0.0400,                          | 0.0010,                       | 0.0005,                         |
| 28,  | naphthalene,          | 720.0000,                    | 0.9167,                         | 89.3036,                          | 629.7796,                      | 7.8011,                        | 101.4145,                        | 0.2909,                       | 0.2461,                         |
| 34,  | phenanthrene,         | 8.5000,                      | 0.0003,                         | 0.2846,                           | 8.2151,                        | 0.0023,                        | 0.3232,                          | 0.0038,                       | 0.0023,                         |
| 37,  | pyrene,               | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 43,  | toluene,              | 770.0000,                    | 14.5997,                        | 328.2077,                         | 427.1926,                      | 124.2391,                      | 372.7173,                        | 0.1973,                       | 0.2322,                         |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.260000,  
NAPL Volume Frac.(l/l), 0.004790,  
Soil Volume Frac.(l/l), 0.735210,  
Porosity (Volume Frac.), 0.264790,

Bulk Density (kg/l), 2.2125,  
NAPL Density (kg/l), 0.8804,

NAPL Saturation (%), 1.8091,

Numerical Accuracy Information

~~~~~  
 The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene  142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene  142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene        154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene      152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene          178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene             78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene   228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene      252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene 252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane 163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene        166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene            228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene        106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene        202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene         128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene        178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene              202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene             92.1400      0.8620      6.3700      535.0000      300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/15/2009 2:49:12 PM

Sample Name Identification, BP-SO-B06-16
Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.3700,

NAPLANAL ANALYSIS RESULTS:

~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 1.0000, 0.0026, 0.9974, 0.0000, 0.0143, 1.2184, 0.0006,  
30, 2-methylnaphthalene, 2.6000, 0.0068, 2.5932, 0.0000, 0.0373, 3.1680, 0.0017,  
32, acenaphthene, 0.1600, 0.0008, 0.1592, 0.0000, 0.0042, 0.1945, 0.0014,  
31, acenaphthylene, 0.6100, 0.0054, 0.6046, 0.0000, 0.0295, 0.7386, 0.0087,  
35, anthracene, 0.1200, 0.0002, 0.1198, 0.0000, 0.0010, 0.1464, 0.0268,  
42, benzene, 760.0000, 160.1654, 599.8346, 0.0000, 882.8577, 732.7719, 0.5818,  
39, benz(a)anthracene, 0.0840, 0.0000, 0.0840, 0.0000, 0.0000, 0.1026, 0.0015,  
40, benzo(a)pyrene, 0.0650, 0.0000, 0.0650, 0.0000, 0.0000, 0.0794, 0.0014,  
50, benzo(b)fluoranthene, 0.0870, 0.0000, 0.0870, 0.0000, 0.0000, 0.1063, 0.0016,  
3, bromodichloromethane, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
33, 9h-fluorene, 0.5800, 0.0018, 0.5782, 0.0000, 0.0097, 0.7064, 0.0066,  
38, chrysene, 0.0800, 0.0000, 0.0800, 0.0000, 0.0000, 0.0977, 0.0313,  
44, ethylbenzene, 12.0000, 0.2370, 11.7630, 0.0000, 1.3064, 14.3700, 0.0099,  
36, fluoranthene, 0.2600, 0.0002, 0.2598, 0.0000, 0.0008, 0.3174, 0.0047,  
28, naphthalene, 16.0000, 0.2682, 15.7318, 0.0000, 1.4783, 19.2183, 0.0538,  
34, phenanthrene, 0.8600, 0.0014, 0.8586, 0.0000, 0.0075, 1.0489, 0.0086,  
37, pyrene, 0.2200, 0.0001, 0.2199, 0.0000, 0.0007, 0.2686, 0.0062,  
43, toluene, 310.0000, 21.3256, 288.6744, 0.0000, 117.5504, 352.6513, 0.2534,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.630000,  
Water Volume Frac.(l/l), 0.370000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.370000,

Bulk Density (kg/l), 2.0395,

Dilution Factor (Vol. fac.), 0.1533,

No NAPL (i.e., NAPL Saturation Equals Zero)

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The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 1380000.0000  |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 5500000.0000  |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 550000.0000   |
| 3   | 75-27-4  | bromodichloromethane | 163.8000                    | 1.9700            | 2.0600                          | 6716.9000                  | 61.0000       |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 200000.0000   |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

\*\*\*\*\*

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/31/2009 8:10:39 AM

Sample Name Identification, BP-SO-B07-12-DL

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1600,

NAPLANAL ANALYSIS RESULTS:

| ID# | Name                 | Total mass, (mg/kg)* | Mass, in water, (mg/kg)* | Mass in, in soil, (mg/kg)* | Mass, in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|--------------------------|----------------------------|-------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 11.0000              | 0.0071                   | 8.4171                     | 2.5758                  | 0.1062                  | 9.0272                    | 0.0049                 | 0.0041                  |
| 30  | 2-methylnaphthalene  | 29.0000              | 0.0187                   | 22.1089                    | 6.8724                  | 0.2790                  | 23.7115                   | 0.0130                 | 0.0110                  |
| 32  | acenaphthene         | 1.9000               | 0.0005                   | 0.3376                     | 1.5619                  | 0.0079                  | 0.3621                    | 0.0030                 | 0.0023                  |
| 31  | acenaphthylene       | 8.7000               | 0.0030                   | 1.0464                     | 7.6505                  | 0.0449                  | 1.1223                    | 0.0145                 | 0.0114                  |
| 35  | anthracene           | 1.1000               | 0.0000                   | 0.0082                     | 1.0918                  | 0.0001                  | 0.0088                    | 0.0021                 | 0.0014                  |
| 42  | benzene              | 680.0000             | 43.9947                  | 507.2641                   | 128.7412                | 655.4631                | 544.0344                  | 0.2443                 | 0.3746                  |
| 39  | benz(a)anthracene    | 0.4600               | 0.0000                   | 0.0313                     | 0.4287                  | 0.0000                  | 0.0336                    | 0.0008                 | 0.0004                  |
| 40  | benzo(a)pyrene       | 0.3000               | 0.0000                   | 0.0158                     | 0.2842                  | 0.0000                  | 0.0169                    | 0.0005                 | 0.0003                  |
| 50  | benzo(b)fluoranthene | 0.2700               | 0.0000                   | 0.0164                     | 0.2536                  | 0.0000                  | 0.0176                    | 0.0005                 | 0.0002                  |
| 3   | bromodichloromethane | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 33  | 9h-fluorene          | 6.8000               | 0.0009                   | 0.9240                     | 5.8751                  | 0.0136                  | 0.9909                    | 0.0111                 | 0.0080                  |
| 38  | chrysene             | 0.4400               | 0.0000                   | 0.0015                     | 0.4385                  | 0.0000                  | 0.0016                    | 0.0008                 | 0.0004                  |
| 44  | ethylbenzene         | 8.4000               | 0.0421                   | 6.4307                     | 1.9272                  | 0.6270                  | 6.8969                    | 0.0037                 | 0.0041                  |
| 36  | fluoranthene         | 1.4000               | 0.0000                   | 0.1061                     | 1.2939                  | 0.0003                  | 0.1138                    | 0.0025                 | 0.0015                  |
| 28  | naphthalene          | 420.0000             | 0.9402                   | 169.8009                   | 249.2588                | 14.0084                 | 182.1093                  | 0.4730                 | 0.4419                  |
| 34  | phenanthrene         | 7.2000               | 0.0005                   | 1.0272                     | 6.1722                  | 0.0079                  | 1.1017                    | 0.0117                 | 0.0079                  |
| 37  | pyrene               | 1.4000               | 0.0000                   | 0.0699                     | 1.3301                  | 0.0002                  | 0.0750                    | 0.0025                 | 0.0015                  |
| 43  | toluene              | 250.0000             | 4.6325                   | 193.0605                   | 52.3070                 | 69.0183                 | 207.0550                  | 0.0993                 | 0.1290                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.160000,  
NAPL Volume Frac.(l/l), 0.001256,  
Soil Volume Frac.(l/l), 0.838744,  
Porosity (Volume Frac.), 0.161256,

Bulk Density (kg/l), 2.3838,  
NAPL Density (kg/l), 0.8882,

NAPL Saturation (%),

0.7790,

Numerical Accuracy Information

The solution converged in 924 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with  
CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 138000.0000   |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 550000.0000   |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 55000.0000    |
| 3   | 75-27-4  | bromodichloromethane | 163.8000                    | 1.9700            | 2.0600                          | 6716.9000                  | 61.0000       |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 20000.0000    |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

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END OF RECORDS



NAPLANAL Version 1.0.0

Date and Time: 7/16/2009 1:39:57 PM

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Sample Name Identification, BP-SO-B08-6

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.2400,

NAPLANAL ANALYSIS RESULTS:  
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ID#,	Name,	Total, mass, (mg/kg)*,	Mass, in water, (mg/kg)*,	Mass in, in soil, (mg/kg)*,	Mass, in NAPL, (mg/kg)*,	Conc., in water, (mg/L),	Sorbed, in soil, (mg/kg)^,	Conc., in NAPL, (kg/L),	Mole fraction, in NAPL,
29,	1-methylnaphthalene,	2.7000,	0.0032,	2.2866,	0.4102,	0.0301,	2.5601,	0.0012,	0.0012,
30,	2-methylnaphthalene,	1.8000,	0.0021,	1.5207,	0.2771,	0.0200,	1.7027,	0.0008,	0.0008,
32,	acenaphthene,	14.0000,	0.0098,	3.7688,	10.2214,	0.0917,	4.2197,	0.0296,	0.0268,
31,	acenaphthylene,	2.7000,	0.0024,	0.5103,	2.1872,	0.0229,	0.5714,	0.0063,	0.0058,
35,	anthracene,	1.8000,	0.0000,	0.0227,	1.7773,	0.0002,	0.0254,	0.0051,	0.0040,
42,	benzene,	130.0000,	14.4553,	100.5796,	14.9651,	135.6776,	112.6124,	0.0433,	0.0775,
39,	benz(a)anthracene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
40,	benzo(a)pyrene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
50,	benzo(b)fluoranthene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
3,	bromodichloromethane,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
33,	9h-fluorene,	6.7000,	0.0023,	1.4167,	5.2810,	0.0217,	1.5861,	0.0153,	0.0129,
38,	chrysene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
44,	ethylbenzene,	36.0000,	0.3290,	30.3405,	5.3305,	3.0882,	33.9703,	0.0154,	0.0203,
36,	fluoranthene,	1.9000,	0.0001,	0.2332,	1.6668,	0.0007,	0.2610,	0.0048,	0.0033,
28,	naphthalene,	550.0000,	2.6992,	294.1547,	253.1461,	25.3343,	329.3459,	0.7328,	0.7992,
34,	phenanthrene,	8.0000,	0.0015,	1.7686,	6.2299,	0.0141,	1.9802,	0.0180,	0.0141,
37,	pyrene,	1.4000,	0.0000,	0.1152,	1.2848,	0.0003,	0.1289,	0.0037,	0.0026,
43,	toluene,	54.0000,	1.7914,	45.0526,	7.1560,	16.8141,	50.4424,	0.0207,	0.0314,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.240000,
NAPL Volume Frac.(l/l), 0.000778,
Soil Volume Frac.(l/l), 0.759222,
Porosity (Volume Frac.), 0.240778,

Bulk Density (kg/l), 2.2526,
NAPL Density (kg/l), 0.8972,

NAPL Saturation (%), 0.3232,

Numerical Accuracy Information

~~~~~  
 The solution converged in 398 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene      142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene      142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene      152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene      228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene      252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene      252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane      163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
=====
  
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/16/2009 1:29:52 PM

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Sample Name Identification, BP-SO-B08-10  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1300,

NAPLANAL ANALYSIS RESULTS:

~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,
, , (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/L), (mg/kg)^, ,
=====
29, 1-methylnaphthalene, 2.2000, 0.0015, 2.1985, 0.0000, 0.0273, 2.3225, 0.0025,
30, 2-methylnaphthalene, 5.4000, 0.0036, 5.3964, 0.0000, 0.0671, 5.7007, 0.0063,
32, acenaphthene, 2.7000, 0.0033, 2.6967, 0.0000, 0.0619, 2.8488, 0.0432,
31, acenaphthylene, 1.2000, 0.0027, 1.1973, 0.0000, 0.0506, 1.2648, 0.0307,
35, anthracene, 0.5200, 0.0002, 0.5198, 0.0000, 0.0039, 0.5491, 0.2079,
42, benzene, 15.0000, 0.9542, 14.0458, 0.0000, 17.8768, 14.8378, 0.0244,
39, benz(a)anthracene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
40, benzo(a)pyrene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
50, benzo(b)fluoranthene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
3, bromodichloromethane, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
33, 9h-fluorene, 1.7000, 0.0013, 1.6987, 0.0000, 0.0246, 1.7945, 0.0347,
38, chrysene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
44, ethylbenzene, 0.5300, 0.0027, 0.5273, 0.0000, 0.0506, 0.5570, 0.0008,
36, fluoranthene, 0.7500, 0.0001, 0.7499, 0.0000, 0.0021, 0.7922, 0.0241,
28, naphthalene, 90.0000, 0.3887, 89.6113, 0.0000, 7.2819, 94.6642, 0.5480,
34, phenanthrene, 2.3000, 0.0009, 2.2991, 0.0000, 0.0173, 2.4287, 0.0414,
37, pyrene, 0.5100, 0.0001, 0.5099, 0.0000, 0.0014, 0.5387, 0.0256,
43, toluene, 6.7000, 0.1236, 6.5764, 0.0000, 2.3157, 6.9472, 0.0103,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.870000,
Water Volume Frac.(l/l), 0.130000,
NAPL Volume Frac.(l/l), 0.000000,
Porosity (Volume Frac.), 0.130000,

Bulk Density (kg/l), 2.4355,

Dilution Factor (Vol. fac.), 1.3856,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

ID#	CAS#	Name	Molecular Weight (g/mol)	Density (kg/l)	Henry's Constant (atm-L/mol)	Water Solubility (mg/L)	Koc (ml/g)
29	90-12-0	1-methylnaphthalene	142.2000	0.9000	0.2600	25.8000	8500.0000
30	91-57-6	2-methylnaphthalene	142.2000	0.9000	0.5180	25.4000	8500.0000
32	83-32-9	acenaphthene	154.2100	0.9000	0.0920	3.4200	4600.0000
31	208-96-8	acenaphthylene	152.2000	0.9000	1.4800	3.9300	2500.0000
35	120-12-7	anthracene	178.2400	0.9000	1.0200	0.0450	14000.0000
42	71-43-2	benzene	78.1100	0.8740	5.5900	1750.0000	83.0000
39	56-55-3	benz(a)anthracene	228.3000	0.9000	0.0012	0.0057	138000.0000
40	50-32-8	benzo(a)pyrene	252.3200	0.9000	0.0016	0.0012	550000.0000
50	205-99-2	benzo(b)fluoranthene	252.3200	0.9000	0.0119	0.0140	55000.0000
3	75-27-4	bromodichloromethane	163.8000	1.9700	2.0600	6716.9000	61.0000
33	86-73-7	9h-fluorene	166.2200	0.9000	0.0642	1.6900	7300.0000
38	218-01-9	chrysene	228.3000	0.9000	0.0011	0.0018	200000.0000
44	100-41-4	ethylbenzene	106.1700	0.8670	6.4300	152.0000	1100.0000
36	206-44-0	fluoranthene	202.2600	0.9000	0.0065	0.2060	38000.0000
28	91-20-3	naphthalene	128.1800	0.9000	1.1500	31.7000	1300.0000
34	85-01-8	phenanthrene	178.2400	0.9000	0.1590	1.0000	14000.0000
37	129-00-0	pyrene	202.2600	0.9000	0.0050	0.1320	38000.0000
43	108-88-3	toluene	92.1400	0.8620	6.3700	535.0000	300.0000

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/31/2009 8:14:25 AM

Sample Name Identification, BP-SO-B08-16
Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.2900,

NAPLANAL ANALYSIS RESULTS:

~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 2.0000, 0.0036, 1.9964, 0.0000, 0.0271, 2.3041, 0.0022,  
30, 2-methylnaphthalene, 5.0000, 0.0091, 4.9909, 0.0000, 0.0678, 5.7602, 0.0057,  
32, acenaphthene, 1.1000, 0.0037, 1.0963, 0.0000, 0.0275, 1.2653, 0.0170,  
31, acenaphthylene, 1.5000, 0.0092, 1.4908, 0.0000, 0.0688, 1.7206, 0.0371,  
35, anthracene, 0.3500, 0.0004, 0.3496, 0.0000, 0.0029, 0.4035, 0.1357,  
42, benzene, 50.0000, 7.8309, 42.1691, 0.0000, 58.6370, 48.6688, 0.0710,  
39, benz(a)anthracene, 0.1800, 0.0000, 0.1800, 0.0000, 0.0000, 0.2077, 0.0056,  
40, benzo(a)pyrene, 0.1000, 0.0000, 0.1000, 0.0000, 0.0000, 0.1154, 0.0037,  
50, benzo(b)fluoranthene, 0.1100, 0.0000, 0.1100, 0.0000, 0.0000, 0.1270, 0.0035,  
3, bromodichloromethane, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
33, 9h-fluorene, 1.4000, 0.0029, 1.3971, 0.0000, 0.0221, 1.6124, 0.0277,  
38, chrysene, 0.1500, 0.0000, 0.1500, 0.0000, 0.0001, 0.1731, 0.1019,  
44, ethylbenzene, 5.6000, 0.0774, 5.5226, 0.0000, 0.5794, 6.3738, 0.0081,  
36, fluoranthene, 0.5500, 0.0002, 0.5498, 0.0000, 0.0017, 0.6345, 0.0172,  
28, naphthalene, 74.0000, 0.8671, 73.1329, 0.0000, 6.4927, 84.4051, 0.4339,  
34, phenanthrene, 1.9000, 0.0021, 1.8979, 0.0000, 0.0156, 2.1904, 0.0331,  
37, pyrene, 0.4100, 0.0002, 0.4098, 0.0000, 0.0012, 0.4730, 0.0200,  
43, toluene, 53.0000, 2.5899, 50.4101, 0.0000, 19.3933, 58.1799, 0.0768,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.710000,  
Water Volume Frac.(l/l), 0.290000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.290000,

Bulk Density (kg/l), 2.1715,

Dilution Factor (Vol. fac.), 1.1183,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
29  90-12-0      1-methylnaphthalene      142.2000      0.9000      0.2600      25.8000      8500.0000
30  91-57-6      2-methylnaphthalene      142.2000      0.9000      0.5180      25.4000      8500.0000
32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
31  208-96-8      acenaphthylene      152.2000      0.9000      1.4800      3.9300      2500.0000
35  120-12-7      anthracene      178.2400      0.9000      1.0200      0.0450      14000.0000
42  71-43-2      benzene      78.1100      0.8740      5.5900      1750.0000      83.0000
39  56-55-3      benz(a)anthracene      228.3000      0.9000      0.0012      0.0057      138000.0000
40  50-32-8      benzo(a)pyrene      252.3200      0.9000      0.0016      0.0012      550000.0000
50  205-99-2      benzo(b)fluoranthene      252.3200      0.9000      0.0119      0.0140      55000.0000
3  75-27-4      bromodichloromethane      163.8000      1.9700      2.0600      6716.9000      61.0000
33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
38  218-01-9      chrysene      228.3000      0.9000      0.0011      0.0018      20000.0000
44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
28  91-20-3      naphthalene      128.1800      0.9000      1.1500      31.7000      1300.0000
34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
37  129-00-0      pyrene      202.2600      0.9000      0.0050      0.1320      38000.0000
43  108-88-3      toluene      92.1400      0.8620      6.3700      535.0000      300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/16/2009 2:14:36 PM

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Sample Name Identification, BP-SO-B09-8  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1900,

NAPLANAL ANALYSIS RESULTS:

~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,
, , (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/L), (mg/kg)^, ,
=====
29, 1-methylnaphthalene, 3.7000, 0.0038, 3.6962, 0.0000, 0.0473, 4.0233, 0.0037,
30, 2-methylnaphthalene, 7.6000, 0.0079, 7.5921, 0.0000, 0.0972, 8.2641, 0.0078,
32, acenaphthene, 2.0000, 0.0038, 1.9962, 0.0000, 0.0472, 2.1729, 0.0280,
31, acenaphthylene, 1.4000, 0.0049, 1.3951, 0.0000, 0.0607, 1.5185, 0.0313,
35, anthracene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
42, benzene, 56.0000, 5.3966, 50.6034, 0.0000, 66.3645, 55.0826, 0.0768,
39, benz(a)anthracene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
40, benzo(a)pyrene, 0.8700, 0.0000, 0.8700, 0.0000, 0.0000, 0.9470, 0.0291,
50, benzo(b)fluoranthene, 1.1000, 0.0000, 1.1000, 0.0000, 0.0002, 1.1973, 0.0315,
3, bromodichloromethane, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
33, 9h-fluorene, 6.1000, 0.0074, 6.0926, 0.0000, 0.0908, 6.6319, 0.1089,
38, chrysene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
44, ethylbenzene, 4.1000, 0.0327, 4.0673, 0.0000, 0.4025, 4.4273, 0.0054,
36, fluoranthene, 1.8000, 0.0004, 1.7996, 0.0000, 0.0052, 1.9589, 0.0507,
28, naphthalene, 82.0000, 0.5546, 81.4454, 0.0000, 6.8196, 88.6547, 0.4359,
34, phenanthrene, 2.9000, 0.0018, 2.8982, 0.0000, 0.0225, 3.1547, 0.0457,
37, pyrene, 2.3000, 0.0005, 2.2995, 0.0000, 0.0066, 2.5030, 0.1011,
43, toluene, 33.0000, 0.9458, 32.0542, 0.0000, 11.6305, 34.8915, 0.0441,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.810000,
Water Volume Frac.(l/l), 0.190000,
NAPL Volume Frac.(l/l), 0.000000,
Porosity (Volume Frac.), 0.190000,

Bulk Density (kg/l), 2.3365,

Dilution Factor (Vol. fac.), 1.0264,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
29  90-12-0  1-methylnaphthalene  142.2000  0.9000  0.2600  25.8000  8500.0000
30  91-57-6  2-methylnaphthalene  142.2000  0.9000  0.5180  25.4000  8500.0000
32  83-32-9  acenaphthene  154.2100  0.9000  0.0920  3.4200  4600.0000
31  208-96-8  acenaphthylene  152.2000  0.9000  1.4800  3.9300  2500.0000
35  120-12-7  anthracene  178.2400  0.9000  1.0200  0.0450  14000.0000
42  71-43-2  benzene  78.1100  0.8740  5.5900  1750.0000  83.0000
39  56-55-3  benz(a)anthracene  228.3000  0.9000  0.0012  0.0057  1380000.0000
40  50-32-8  benzo(a)pyrene  252.3200  0.9000  0.0016  0.0012  5500000.0000
50  205-99-2  benzo(b)fluoranthene  252.3200  0.9000  0.0119  0.0140  550000.0000
3  75-27-4  bromodichloromethane  163.8000  1.9700  2.0600  6716.9000  61.0000
33  86-73-7  9h-fluorene  166.2200  0.9000  0.0642  1.6900  7300.0000
38  218-01-9  chrysene  228.3000  0.9000  0.0011  0.0018  200000.0000
44  100-41-4  ethylbenzene  106.1700  0.8670  6.4300  152.0000  1100.0000
36  206-44-0  fluoranthene  202.2600  0.9000  0.0065  0.2060  38000.0000
28  91-20-3  naphthalene  128.1800  0.9000  1.1500  31.7000  1300.0000
34  85-01-8  phenanthrene  178.2400  0.9000  0.1590  1.0000  14000.0000
37  129-00-0  pyrene  202.2600  0.9000  0.0050  0.1320  38000.0000
43  108-88-3  toluene  92.1400  0.8620  6.3700  535.0000  300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/16/2009 2:05:10 PM

Sample Name Identification, BP-SO-B09-14

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.1800,

NAPLANAL ANALYSIS RESULTS:

ID#,	Name,	Total, mass, (mg/kg)*,	Mass, in water, (mg/kg)*,	Mass in, in soil, (mg/kg)*,	Mass, in NAPL, (mg/kg)*,	Conc., in water, (mg/L),	Sorbed, in soil, (mg/kg)^,	Conc., in NAPL, (kg/L),	Mole fraction, in NAPL, ,
==,	====,	====,	====,	====,	====,	====,	====,	====,	====,
29,	1-methylnaphthalene,	5.9000,	0.0008,	0.7585,	5.1407,	0.0098,	0.8298,	0.0005,	0.0004,
30,	2-methylnaphthalene,	14.0000,	0.0018,	1.7755,	12.2227,	0.0229,	1.9424,	0.0013,	0.0009,
32,	acenapthene,	3.1000,	0.0001,	0.0300,	3.0700,	0.0007,	0.0328,	0.0003,	0.0002,
31,	acenapthylene,	5.4000,	0.0001,	0.0331,	5.3667,	0.0015,	0.0363,	0.0006,	0.0004,
35,	anthracene,	27.0000,	0.0000,	0.0091,	26.9909,	0.0001,	0.0100,	0.0029,	0.0016,
42,	benzene,	6100.0000,	93.0033,	907.3018,	5099.6948,	1195.8847,	992.5843,	0.5456,	0.6834,
39,	benz(a)anthracene,	29.0000,	0.0000,	0.0953,	28.9047,	0.0000,	0.1042,	0.0031,	0.0013,
40,	benzo(a)pyrene,	26.0000,	0.0000,	0.0649,	25.9351,	0.0000,	0.0710,	0.0028,	0.0011,
50,	benzo(b)fluoranthene,	37.0000,	0.0000,	0.1077,	36.8923,	0.0000,	0.1178,	0.0039,	0.0015,
3,	bromodichloromethane,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
33,	9h-fluorene,	18.0000,	0.0001,	0.1269,	17.8729,	0.0019,	0.1388,	0.0019,	0.0011,
38,	chrysene,	31.0000,	0.0000,	0.0047,	30.9953,	0.0000,	0.0051,	0.0033,	0.0014,
44,	ethylbenzene,	47.0000,	0.0476,	6.1481,	40.8044,	0.6115,	6.7260,	0.0044,	0.0040,
36,	fluoranthene,	89.0000,	0.0001,	0.3283,	88.6716,	0.0009,	0.3592,	0.0095,	0.0046,
28,	naphthalene,	110.0000,	0.0215,	3.2819,	106.6966,	0.2762,	3.5904,	0.0114,	0.0087,
34,	phenanthrene,	94.0000,	0.0004,	0.7011,	93.2984,	0.0055,	0.7670,	0.0100,	0.0055,
37,	pyrene,	52.0000,	0.0000,	0.1231,	51.8769,	0.0004,	0.1347,	0.0055,	0.0027,
43,	toluene,	2900.0000,	11.7011,	412.5947,	2475.7042,	150.4589,	451.3768,	0.2649,	0.2812,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.180000,
NAPL Volume Frac.(l/l), 0.021635,
Soil Volume Frac.(l/l), 0.798365,
Porosity (Volume Frac.), 0.201635,

Bulk Density (kg/l), 2.3145,
NAPL Density (kg/l), 0.8719,

NAPL Saturation (%), 10.7296,

Numerical Accuracy Information

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 The solution converged in 16 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene  142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene  142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene  154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene  152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene  178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene  78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene  228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene  252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene  252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane  163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene  166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene  228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene  106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene  202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene  128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene  178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene  202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene  92.1400      0.8620      6.3700      535.0000      300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/16/2009 1:56:52 PM

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Sample Name Identification, BP-SO-B09-18

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.3300,

NAPLANAL ANALYSIS RESULTS:  
~~~~~

ID#,	Name,	Total, mass, (mg/kg)*,	Mass, in water, (mg/kg)*,	Mass in, in soil, (mg/kg)*,	Mass, in NAPL, (mg/kg)*,	Conc., in water, (mg/L),	Sorbed, in soil, (mg/kg)^,	Conc., in NAPL, (kg/L),	Mole fraction, in NAPL, ,
29,	1-methylnaphthalene,	3.0000,	0.0008,	0.3392,	2.6600,	0.0048,	0.4078,	0.0003,	0.0002,
30,	2-methylnaphthalene,	7.7000,	0.0019,	0.8587,	6.8394,	0.0121,	1.0322,	0.0007,	0.0005,
32,	acenapthene,	0.7100,	0.0000,	0.0059,	0.7040,	0.0002,	0.0071,	0.0001,	0.0000,
31,	acenapthylene,	1.1000,	0.0000,	0.0058,	1.0941,	0.0003,	0.0070,	0.0001,	0.0001,
35,	anthracene,	7.8000,	0.0000,	0.0023,	7.7977,	0.0000,	0.0027,	0.0008,	0.0004,
42,	benzene,	5600.0000,	167.2341,	724.0832,	4708.6826,	1048.6112,	870.3473,	0.4779,	0.5992,
39,	benz(a)anthracene,	11.0000,	0.0000,	0.0313,	10.9687,	0.0000,	0.0376,	0.0011,	0.0005,
40,	benzo(a)pyrene,	8.4000,	0.0000,	0.0181,	8.3819,	0.0000,	0.0218,	0.0009,	0.0003,
50,	benzo(b)fluoranthene,	13.0000,	0.0000,	0.0327,	12.9673,	0.0000,	0.0393,	0.0013,	0.0005,
3,	bromodichloromethane,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
33,	9h-fluorene,	2.5000,	0.0000,	0.0153,	2.4847,	0.0003,	0.0183,	0.0003,	0.0001,
38,	chrysene,	11.0000,	0.0000,	0.0014,	10.9986,	0.0000,	0.0017,	0.0011,	0.0005,
44,	ethylbenzene,	350.0000,	0.7014,	40.2479,	309.0507,	4.3980,	48.3780,	0.0314,	0.0289,
36,	fluoranthene,	26.0000,	0.0000,	0.0829,	25.9170,	0.0003,	0.0997,	0.0026,	0.0013,
28,	naphthalene,	130.0000,	0.0496,	3.3655,	126.5849,	0.3112,	4.0453,	0.0128,	0.0098,
34,	phenanthrene,	21.0000,	0.0002,	0.1355,	20.8643,	0.0012,	0.1629,	0.0021,	0.0012,
37,	pyrene,	20.0000,	0.0000,	0.0409,	19.9590,	0.0001,	0.0492,	0.0020,	0.0010,
43,	toluene,	3800.0000,	30.3291,	474.6415,	3295.0294,	190.1729,	570.5186,	0.3344,	0.3555,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.330000,
NAPL Volume Frac.(l/l), 0.020389,
Soil Volume Frac.(l/l), 0.649611,
Porosity (Volume Frac.), 0.350389,

Bulk Density (kg/l), 2.0692,
NAPL Density (kg/l), 0.8698,

NAPL Saturation (%), 5.8189,

Numerical Accuracy Information

~~~~~  
 The solution converged in 16 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25°C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
      (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
  29  90-12-0      1-methylnaphthalene  142.2000      0.9000      0.2600      25.8000      8500.0000
  30  91-57-6      2-methylnaphthalene  142.2000      0.9000      0.5180      25.4000      8500.0000
  32  83-32-9      acenaphthene      154.2100      0.9000      0.0920      3.4200      4600.0000
  31  208-96-8      acenaphthylene    152.2000      0.9000      1.4800      3.9300      2500.0000
  35  120-12-7      anthracene        178.2400      0.9000      1.0200      0.0450      14000.0000
  42  71-43-2      benzene           78.1100      0.8740      5.5900      1750.0000      83.0000
  39  56-55-3      benz(a)anthracene  228.3000      0.9000      0.0012      0.0057      1380000.0000
  40  50-32-8      benzo(a)pyrene    252.3200      0.9000      0.0016      0.0012      5500000.0000
  50  205-99-2      benzo(b)fluoranthene  252.3200      0.9000      0.0119      0.0140      550000.0000
   3  75-27-4      bromodichloromethane  163.8000      1.9700      2.0600      6716.9000      61.0000
  33  86-73-7      9h-fluorene      166.2200      0.9000      0.0642      1.6900      7300.0000
  38  218-01-9      chrysene          228.3000      0.9000      0.0011      0.0018      200000.0000
  44  100-41-4      ethylbenzene      106.1700      0.8670      6.4300      152.0000      1100.0000
  36  206-44-0      fluoranthene      202.2600      0.9000      0.0065      0.2060      38000.0000
  28  91-20-3      naphthalene       128.1800      0.9000      1.1500      31.7000      1300.0000
  34  85-01-8      phenanthrene      178.2400      0.9000      0.1590      1.0000      14000.0000
  37  129-00-0      pyrene            202.2600      0.9000      0.0050      0.1320      38000.0000
  43  108-88-3      toluene           92.1400      0.8620      6.3700      535.0000      300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/16/2009 2:30:38 PM

-----  
Sample Name Identification, BP-SO-B10-4  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.2000,

NAPLANAL ANALYSIS RESULTS:

~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,
, , (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/L), (mg/kg)^, ,
=====
29, 1-methylnaphthalene, 1.4000, 0.0016, 1.3984, 0.0000, 0.0180, 1.5304, 0.0009,
30, 2-methylnaphthalene, 2.1000, 0.0023, 2.0977, 0.0000, 0.0270, 2.2956, 0.0014,
32, acenaphthene, 4.3000, 0.0088, 4.2912, 0.0000, 0.1021, 4.6960, 0.0396,
31, acenaphthylene, 3.6000, 0.0135, 3.5865, 0.0000, 0.1570, 3.9248, 0.0531,
35, anthracene, 1.1000, 0.0007, 1.0993, 0.0000, 0.0086, 1.2030, 0.2536,
42, benzene, 6.1000, 0.6226, 5.4774, 0.0000, 7.2219, 5.9942, 0.0055,
39, benz(a)anthracene, 0.8900, 0.0000, 0.8900, 0.0000, 0.0001, 0.9740, 0.0164,
40, benzo(a)pyrene, 0.6500, 0.0000, 0.6500, 0.0000, 0.0000, 0.7113, 0.0143,
50, benzo(b)fluoranthene, 0.6400, 0.0000, 0.6400, 0.0000, 0.0001, 0.7004, 0.0121,
3, bromodichloromethane, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
33, 9h-fluorene, 3.0000, 0.0039, 2.9961, 0.0000, 0.0449, 3.2788, 0.0353,
38, chrysene, 0.8100, 0.0000, 0.8100, 0.0000, 0.0004, 0.8864, 0.3270,
44, ethylbenzene, 0.5700, 0.0048, 0.5652, 0.0000, 0.0562, 0.6185, 0.0005,
36, fluoranthene, 2.9000, 0.0007, 2.8993, 0.0000, 0.0083, 3.1728, 0.0538,
28, naphthalene, 25.0000, 0.1801, 24.8199, 0.0000, 2.0893, 27.1614, 0.0875,
34, phenanthrene, 3.5000, 0.0024, 3.4976, 0.0000, 0.0273, 3.8276, 0.0363,
37, pyrene, 2.1000, 0.0005, 2.0995, 0.0000, 0.0060, 2.2975, 0.0608,
43, toluene, 1.9000, 0.0579, 1.8421, 0.0000, 0.6720, 2.0159, 0.0017,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.800000,
Water Volume Frac.(l/l), 0.200000,
NAPL Volume Frac.(l/l), 0.000000,
Porosity (Volume Frac.), 0.200000,

Bulk Density (kg/l), 2.3200,

Dilution Factor (Vol. fac.), 0.3282,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

ID#	CAS#	Name	Molecular Weight (g/mol)	Density (kg/l)	Henry's Constant (atm-L/mol)	Water Solubility (mg/L)	Koc (ml/g)
29	90-12-0	1-methylnaphthalene	142.2000	0.9000	0.2600	25.8000	8500.0000
30	91-57-6	2-methylnaphthalene	142.2000	0.9000	0.5180	25.4000	8500.0000
32	83-32-9	acenaphthene	154.2100	0.9000	0.0920	3.4200	4600.0000
31	208-96-8	acenaphthylene	152.2000	0.9000	1.4800	3.9300	2500.0000
35	120-12-7	anthracene	178.2400	0.9000	1.0200	0.0450	14000.0000
42	71-43-2	benzene	78.1100	0.8740	5.5900	1750.0000	83.0000
39	56-55-3	benz(a)anthracene	228.3000	0.9000	0.0012	0.0057	138000.0000
40	50-32-8	benzo(a)pyrene	252.3200	0.9000	0.0016	0.0012	550000.0000
50	205-99-2	benzo(b)fluoranthene	252.3200	0.9000	0.0119	0.0140	55000.0000
3	75-27-4	bromodichloromethane	163.8000	1.9700	2.0600	6716.9000	61.0000
33	86-73-7	9h-fluorene	166.2200	0.9000	0.0642	1.6900	7300.0000
38	218-01-9	chrysene	228.3000	0.9000	0.0011	0.0018	20000.0000
44	100-41-4	ethylbenzene	106.1700	0.8670	6.4300	152.0000	1100.0000
36	206-44-0	fluoranthene	202.2600	0.9000	0.0065	0.2060	38000.0000
28	91-20-3	naphthalene	128.1800	0.9000	1.1500	31.7000	1300.0000
34	85-01-8	phenanthrene	178.2400	0.9000	0.1590	1.0000	14000.0000
37	129-00-0	pyrene	202.2600	0.9000	0.0050	0.1320	38000.0000
43	108-88-3	toluene	92.1400	0.8620	6.3700	535.0000	300.0000

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/16/2009 3:10:17 PM

Sample Name Identification, BP-SO-B11-4-DL
Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.1700,

NAPLANAL ANALYSIS RESULTS:

~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
=====  
29, 1-methylnaphthalene, 0.2000, 0.0002, 0.1998, 0.0000, 0.0025, 0.2153, 0.0020,  
30, 2-methylnaphthalene, 0.5000, 0.0005, 0.4995, 0.0000, 0.0063, 0.5382, 0.0051,  
32, acenaphthene, 0.0250, 0.0000, 0.0250, 0.0000, 0.0006, 0.0269, 0.0035,  
31, acenaphthylene, 0.0240, 0.0001, 0.0239, 0.0000, 0.0010, 0.0258, 0.0053,  
35, anthracene, 0.0200, 0.0000, 0.0200, 0.0000, 0.0002, 0.0215, 0.0694,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
40, benzo(a)pyrene, 0.0110, 0.0000, 0.0110, 0.0000, 0.0000, 0.0119, 0.0036,  
50, benzo(b)fluoranthene, 0.0290, 0.0000, 0.0290, 0.0000, 0.0000, 0.0312, 0.0082,  
3, bromodichloromethane, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
33, 9h-fluorene, 0.7600, 0.0008, 0.7592, 0.0000, 0.0112, 0.8179, 0.1347,  
38, chrysene, 0.0920, 0.0000, 0.0920, 0.0000, 0.0000, 0.0991, 0.5593,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 0.0840, 0.0000, 0.0840, 0.0000, 0.0002, 0.0905, 0.0235,  
28, naphthalene, 1.9000, 0.0112, 1.8888, 0.0000, 0.1565, 2.0348, 0.1003,  
34, phenanthrene, 0.3000, 0.0002, 0.2998, 0.0000, 0.0023, 0.3230, 0.0469,  
37, pyrene, 0.0850, 0.0000, 0.0850, 0.0000, 0.0002, 0.0916, 0.0371,  
43, toluene, 0.0750, 0.0019, 0.0731, 0.0000, 0.0263, 0.0788, 0.0010,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.830000,  
Water Volume Frac.(l/l), 0.170000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.170000,

Bulk Density (kg/l), 2.3695,

Dilution Factor (Vol. fac.), 19.3178,

No NAPL (i.e., NAPL Saturation Equals Zero)

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The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 138000.0000   |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 550000.0000   |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 55000.0000    |
| 3   | 75-27-4  | bromodichloromethane | 163.8000                    | 1.9700            | 2.0600                          | 6716.9000                  | 61.0000       |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 20000.0000    |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

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END OF RECORDS



NAPLANAL Version 1.0.0

Date and Time:

8/31/2009 8:37:28 AM

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Sample Name Identification, CT-SO-B01-10  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1000,  
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NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,
, , (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/L), (mg/kg)^, ,
====, =====, =====, =====, =====, =====, =====, =====, =====,
29, 1-methylnaphthalene, 0.2300, 0.0001, 0.2299, 0.0000, 0.0028, 0.2395, 0.0002,
30, 2-methylnaphthalene, 0.3600, 0.0002, 0.3598, 0.0000, 0.0044, 0.3749, 0.0004,
32, acenaphthene, 0.0470, 0.0000, 0.0470, 0.0000, 0.0011, 0.0489, 0.0007,
31, acenaphthylene, 0.0860, 0.0001, 0.0859, 0.0000, 0.0036, 0.0895, 0.0019,
35, anthracene, 0.2100, 0.0001, 0.2099, 0.0000, 0.0016, 0.2187, 0.0739,
42, benzene, 0.4900, 0.0236, 0.4664, 0.0000, 0.5855, 0.4860, 0.0007,
39, benz(a)anthracene, 0.7700, 0.0000, 0.7700, 0.0000, 0.0001, 0.8023, 0.0217,
40, benzo(a)pyrene, 0.4700, 0.0000, 0.4700, 0.0000, 0.0000, 0.4897, 0.0158,
50, benzo(b)fluoranthene, 1.0000, 0.0000, 1.0000, 0.0000, 0.0002, 1.0419, 0.0288,
51, benzo(k)fluoranthene, 0.0370, 0.0000, 0.0370, 0.0000, 0.0000, 0.0386, 0.0035,
38, chrysene, 1.1000, 0.0000, 1.1000, 0.0000, 0.0006, 1.1461, 0.6776,
44, ethylbenzene, 0.0540, 0.0002, 0.0538, 0.0000, 0.0051, 0.0561, 0.0001,
36, fluoranthene, 2.6000, 0.0003, 2.5997, 0.0000, 0.0071, 2.7087, 0.0737,
33, 9h-fluorene, 0.1100, 0.0001, 0.1099, 0.0000, 0.0016, 0.1145, 0.0020,
28, naphthalene, 0.9300, 0.0030, 0.9270, 0.0000, 0.0743, 0.9659, 0.0050,
34, phenanthrene, 2.0000, 0.0006, 1.9994, 0.0000, 0.0149, 2.0832, 0.0317,
37, pyrene, 1.4000, 0.0002, 1.3998, 0.0000, 0.0038, 1.4585, 0.0619,
43, toluene, 0.3800, 0.0052, 0.3748, 0.0000, 0.1302, 0.3905, 0.0005,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.900000,
Water Volume Frac.(l/l), 0.100000,
NAPL Volume Frac.(l/l), 0.000000,
Porosity (Volume Frac.), 0.100000,

Bulk Density (kg/l), 2.4850,

Dilution Factor (Vol. fac.), 1.1285,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID#      CAS#      Name                Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
              (g/mol)          (kg/l)    (atm-L/mol)      (mg/L)           (ml/g)
=====
 29  90-12-0    1-methylnaphthalene  142.2000         0.9000    0.2600           25.8000          8500.0000
 30  91-57-6    2-methylnaphthalene  142.2000         0.9000    0.5180           25.4000          8500.0000
 32  83-32-9    acenaphthene         154.2100         0.9000    0.0920           3.4200           4600.0000
 31  208-96-8    acenaphthylene       152.2000         0.9000    1.4800           3.9300           2500.0000
 35  120-12-7    anthracene           178.2400         0.9000    1.0200           0.0450           14000.0000
 42  71-43-2    benzene               78.1100          0.8740    5.5900           1750.0000         83.0000
 39  56-55-3    benz(a)anthracene    228.3000         0.9000    0.0012           0.0057           138000.0000
 40  50-32-8    benzo(a)pyrene       252.3200         0.9000    0.0016           0.0012           550000.0000
 50  205-99-2    benzo(b)fluoranthene 252.3200         0.9000    0.0119           0.0140           55000.0000
 51  207-08-9    benzo(k)fluoranthene 252.3200         0.9000    0.0394           0.0043           55000.0000
 38  218-01-9    chrysene             228.3000         0.9000    0.0011           0.0018           20000.0000
 44  100-41-4    ethylbenzene         106.1700         0.8670    6.4300           152.0000          110.0000
 36  206-44-0    fluoranthene         202.2600         0.9000    0.0065           0.2060           38000.0000
 33  86-73-7    9h-fluorene         166.2200         0.9000    0.0642           1.6900           7300.0000
 28  91-20-3    naphthalene          128.1800         0.9000    1.1500           31.7000           1300.0000
 34  85-01-8    phenanthrene         178.2400         0.9000    0.1590           1.0000           14000.0000
 37  129-00-0    pyrene               202.2600         0.9000    0.0050           0.1320           38000.0000
 43  108-88-3    toluene              92.1400          0.8620    6.3700           535.0000          300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/5/2009 8:10:57 AM

Sample Name Identification, CT-SO-B01-14
Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.1500,

NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.1200, 0.0001, 0.1199, 0.0000, 0.0015, 0.1279, 0.0002,  
30, 2-methylnaphthalene, 1.2900, 0.0010, 1.2890, 0.0000, 0.0162, 1.3748, 0.0021,  
32, acenaphthene, 0.0520, 0.0001, 0.0519, 0.0000, 0.0012, 0.0554, 0.0012,  
31, acenaphthylene, 0.1400, 0.0004, 0.1396, 0.0000, 0.0060, 0.1489, 0.0051,  
35, anthracene, 0.2600, 0.0001, 0.2599, 0.0000, 0.0020, 0.2772, 0.1472,  
42, benzene, 0.6300, 0.0468, 0.5832, 0.0000, 0.7495, 0.6220, 0.0014,  
39, benz(a)anthracene, 0.5400, 0.0000, 0.5400, 0.0000, 0.0000, 0.5760, 0.0245,  
40, benzo(a)pyrene, 0.3900, 0.0000, 0.3900, 0.0000, 0.0000, 0.4160, 0.0211,  
50, benzo(b)fluoranthene, 0.6800, 0.0000, 0.6800, 0.0000, 0.0001, 0.7253, 0.0315,  
51, benzo(k)fluoranthene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
38, chrysene, 0.6000, 0.0000, 0.6000, 0.0000, 0.0003, 0.6399, 0.5947,  
44, ethylbenzene, 0.0580, 0.0003, 0.0577, 0.0000, 0.0056, 0.0615, 0.0001,  
36, fluoranthene, 1.3000, 0.0002, 1.2998, 0.0000, 0.0036, 1.3863, 0.0593,  
33, 9h-fluorene, 0.1900, 0.0002, 0.1898, 0.0000, 0.0028, 0.2025, 0.0055,  
28, naphthalene, 2.0000, 0.0102, 1.9898, 0.0000, 0.1633, 2.1223, 0.0172,  
34, phenanthrene, 1.1000, 0.0005, 1.0995, 0.0000, 0.0084, 1.1727, 0.0280,  
37, pyrene, 0.8400, 0.0001, 0.8399, 0.0000, 0.0024, 0.8958, 0.0597,  
43, toluene, 0.5100, 0.0111, 0.4989, 0.0000, 0.1774, 0.5321, 0.0011,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.850000,  
Water Volume Frac.(l/l), 0.150000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.150000,

Bulk Density (kg/l), 2.4025,

Dilution Factor (Vol. fac.), 2.3456,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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ID#      CAS#      Name                Molecular Weight   Density   Henry's Constant   Water Solubility   Koc
              (g/mol)          (kg/l)          (atm-L/mol)      (mg/L)          (ml/g)
=====
 29  90-12-0    1-methylnaphthalene  142.2000          0.9000          0.2600            25.8000           8500.0000
 30  91-57-6    2-methylnaphthalene  142.2000          0.9000          0.5180            25.4000           8500.0000
 32  83-32-9    acenaphthene         154.2100          0.9000          0.0920            3.4200            4600.0000
 31  208-96-8    acenaphthylene       152.2000          0.9000          1.4800            3.9300            2500.0000
 35  120-12-7    anthracene           178.2400          0.9000          1.0200            0.0450            14000.0000
 42  71-43-2    benzene              78.1100           0.8740          5.5900            1750.0000          83.0000
 39  56-55-3    benz(a)anthracene    228.3000          0.9000          0.0012            0.0057            138000.0000
 40  50-32-8    benzo(a)pyrene       252.3200          0.9000          0.0016            0.0012            550000.0000
 50  205-99-2    benzo(b)fluoranthene 252.3200          0.9000          0.0119            0.0140            55000.0000
 51  207-08-9    benzo(k)fluoranthene 252.3200          0.9000          0.0394            0.0043            55000.0000
 38  218-01-9    chrysene             228.3000          0.9000          0.0011            0.0018            20000.0000
 44  100-41-4    ethylbenzene         106.1700          0.8670          6.4300            152.0000           1100.0000
 36  206-44-0    fluoranthene         202.2600          0.9000          0.0065            0.2060            38000.0000
 33  86-73-7    9h-fluorene         166.2200          0.9000          0.0642            1.6900             7300.0000
 28  91-20-3    naphthalene          128.1800          0.9000          1.1500            31.7000            1300.0000
 34  85-01-8    phenanthrene         178.2400          0.9000          0.1590            1.0000            14000.0000
 37  129-00-0    pyrene               202.2600          0.9000          0.0050            0.1320            38000.0000
 43  108-88-3    toluene              92.1400           0.8620          6.3700            535.0000           300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/5/2009 8:19:46 AM

Sample Name Identification, CT-SO-B01-18

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.2400,

NAPLANAL ANALYSIS RESULTS:

| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 160.0000             | 0.0874                  | 61.6381                | 98.2745                | 0.8154                  | 69.3121                   | 0.0252                 | 0.0316                  |
| 30  | 2-methylnaphthalene  | 410.0000             | 0.2219                  | 156.4344               | 253.3437               | 2.0695                  | 175.9108                  | 0.0649                 | 0.0815                  |
| 32  | acenaphthene         | 19.0000              | 0.0020                  | 0.7568                 | 18.2412                | 0.0185                  | 0.8510                    | 0.0047                 | 0.0054                  |
| 31  | acenaphthylene       | 350.0000             | 0.0432                  | 8.9526                 | 341.0043               | 0.4027                  | 10.0672                   | 0.0874                 | 0.1025                  |
| 35  | anthracene           | 140.0000             | 0.0002                  | 0.2010                 | 139.7989               | 0.0016                  | 0.2260                    | 0.0358                 | 0.0359                  |
| 42  | benzene              | 15.0000              | 0.8829                  | 6.0790                 | 8.0381                 | 8.2359                  | 6.8358                    | 0.0021                 | 0.0047                  |
| 39  | benz(a)anthracene    | 130.0000             | 0.0000                  | 1.7965                 | 128.2035               | 0.0001                  | 2.0201                    | 0.0328                 | 0.0257                  |
| 40  | benzo(a)pyrene       | 99.0000              | 0.0000                  | 1.0421                 | 97.9579                | 0.0000                  | 1.1718                    | 0.0251                 | 0.0178                  |
| 50  | benzo(b)fluoranthene | 130.0000             | 0.0000                  | 1.5937                 | 128.4063               | 0.0003                  | 1.7921                    | 0.0329                 | 0.0233                  |
| 51  | benzo(k)fluoranthene | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 38  | chrysene             | 110.0000             | 0.0000                  | 0.0705                 | 109.9295               | 0.0000                  | 0.0793                    | 0.0282                 | 0.0220                  |
| 44  | ethylbenzene         | 1.3000               | 0.0055                  | 0.5054                 | 0.7891                 | 0.0517                  | 0.5683                    | 0.0002                 | 0.0003                  |
| 36  | fluoranthene         | 350.0000             | 0.0017                  | 5.4237                 | 344.5746               | 0.0160                  | 6.0989                    | 0.0883                 | 0.0779                  |
| 33  | 9h-fluorene          | 190.0000             | 0.0092                  | 5.5669                 | 184.4239               | 0.0858                  | 6.2600                    | 0.0472                 | 0.0507                  |
| 28  | naphthalene          | 1100.0000            | 1.1783                  | 127.0606               | 971.7611               | 10.9908                 | 142.8799                  | 0.2489                 | 0.3467                  |
| 34  | phenanthrene         | 490.0000             | 0.0131                  | 15.1677                | 474.8192               | 0.1218                  | 17.0561                   | 0.1216                 | 0.1218                  |
| 37  | pyrene               | 200.0000             | 0.0006                  | 1.9970                 | 198.0023               | 0.0059                  | 2.2457                    | 0.0507                 | 0.0448                  |
| 43  | toluene              | 26.0000              | 0.4261                  | 10.6046                | 14.9692                | 3.9750                  | 11.9249                   | 0.0038                 | 0.0074                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.240000,  
NAPL Volume Frac.(l/l), 0.008740,  
Soil Volume Frac.(l/l), 0.751260,  
Porosity (Volume Frac.), 0.248740,

Bulk Density (kg/l), 2.2387,  
NAPL Density (kg/l), 0.8998,

NAPL Saturation (%),

3.5135,

Numerical Accuracy Information

The solution converged in 100 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/5/2009 8:38:48 AM

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Sample Name Identification, CT-SO-B02-12  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1600,  
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NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,
, , (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/L), (mg/kg)^, ,
====, =====, =====, =====, =====, =====, =====, =====, =====,
29, 1-methylnaphthalene, 0.0750, 0.0001, 0.0749, 0.0000, 0.0009, 0.0803, 0.0004,
30, 2-methylnaphthalene, 0.1800, 0.0002, 0.1798, 0.0000, 0.0023, 0.1928, 0.0011,
32, acenaphthene, 0.0270, 0.0000, 0.0270, 0.0000, 0.0006, 0.0289, 0.0022,
31, acenaphthylene, 0.0870, 0.0002, 0.0868, 0.0000, 0.0037, 0.0930, 0.0112,
35, anthracene, 0.0720, 0.0000, 0.0720, 0.0000, 0.0006, 0.0771, 0.1451,
42, benzene, 0.1700, 0.0135, 0.1565, 0.0000, 0.2020, 0.1677, 0.0014,
39, benz(a)anthracene, 0.1500, 0.0000, 0.1500, 0.0000, 0.0000, 0.1608, 0.0242,
40, benzo(a)pyrene, 0.1100, 0.0000, 0.1100, 0.0000, 0.0000, 0.1179, 0.0212,
50, benzo(b)fluoranthene, 0.1900, 0.0000, 0.1900, 0.0000, 0.0000, 0.2037, 0.0313,
51, benzo(k)fluoranthene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
38, chrysene, 0.1600, 0.0000, 0.1600, 0.0000, 0.0001, 0.1715, 0.5645,
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
36, fluoranthene, 0.3600, 0.0001, 0.3599, 0.0000, 0.0010, 0.3858, 0.0584,
33, 9h-fluorene, 0.0590, 0.0001, 0.0589, 0.0000, 0.0009, 0.0632, 0.0061,
28, naphthalene, 1.5000, 0.0082, 1.4918, 0.0000, 0.1230, 1.5990, 0.0460,
34, phenanthrene, 0.2900, 0.0001, 0.2899, 0.0000, 0.0022, 0.3107, 0.0263,
37, pyrene, 0.2400, 0.0000, 0.2400, 0.0000, 0.0007, 0.2572, 0.0608,
43, toluene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.840000,
Water Volume Frac.(l/l), 0.160000,
NAPL Volume Frac.(l/l), 0.000000,
Porosity (Volume Frac.), 0.160000,

Bulk Density (kg/l), 2.3860,

Dilution Factor (Vol. fac.), 10.8490,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name                Molecular Weight   Density   Henry's Constant   Water Solubility   Koc
              (g/mol)          (kg/l)          (atm-L/mol)      (mg/L)          (ml/g)
=====
 29  90-12-0    1-methylnaphthalene  142.2000          0.9000          0.2600            25.8000           8500.0000
 30  91-57-6    2-methylnaphthalene  142.2000          0.9000          0.5180            25.4000           8500.0000
 32  83-32-9    acenaphthene         154.2100          0.9000          0.0920            3.4200            4600.0000
 31  208-96-8    acenaphthylene       152.2000          0.9000          1.4800            3.9300            2500.0000
 35  120-12-7    anthracene           178.2400          0.9000          1.0200            0.0450            14000.0000
 42  71-43-2    benzene              78.1100           0.8740          5.5900            1750.0000          83.0000
 39  56-55-3    benz(a)anthracene    228.3000          0.9000          0.0012            0.0057            138000.0000
 40  50-32-8    benzo(a)pyrene       252.3200          0.9000          0.0016            0.0012            550000.0000
 50  205-99-2    benzo(b)fluoranthene 252.3200          0.9000          0.0119            0.0140            55000.0000
 51  207-08-9    benzo(k)fluoranthene 252.3200          0.9000          0.0394            0.0043            55000.0000
 38  218-01-9    chrysene             228.3000          0.9000          0.0011            0.0018            20000.0000
 44  100-41-4    ethylbenzene         106.1700          0.8670          6.4300            152.0000           1100.0000
 36  206-44-0    fluoranthene         202.2600          0.9000          0.0065            0.2060            38000.0000
 33  86-73-7    9h-fluorene         166.2200          0.9000          0.0642            1.6900            7300.0000
 28  91-20-3    naphthalene         128.1800          0.9000          1.1500            31.7000            1300.0000
 34  85-01-8    phenanthrene        178.2400          0.9000          0.1590            1.0000            14000.0000
 37  129-00-0    pyrene              202.2600          0.9000          0.0050            0.1320            38000.0000
 43  108-88-3    toluene             92.1400           0.8620          6.3700            535.0000           300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/31/2009 8:39:08 AM

Sample Name Identification, CT-SO-B02-16

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.1400,

NAPLANAL ANALYSIS RESULTS:
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| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc. in water, (mg/L) | Sorbed in soil, (mg/kg)^ | Conc. in NAPL, (kg/L) | Mole fraction in NAPL |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|------------------------|--------------------------|-----------------------|-----------------------|
| 29  | 1-methylnaphthalene  | 6.7000               | 0.0048                  | 6.6142                 | 0.0810                 | 0.0826                 | 7.0208                   | 0.0024                | 0.0032                |
| 30  | 2-methylnaphthalene  | 16.0000              | 0.0114                  | 15.7921                | 0.1965                 | 0.1972                 | 16.7628                  | 0.0057                | 0.0078                |
| 32  | acenaphthene         | 0.9100               | 0.0010                  | 0.7670                 | 0.1420                 | 0.0177                 | 0.8141                   | 0.0041                | 0.0052                |
| 31  | acenaphthylene       | 14.0000              | 0.0266                  | 10.8100                | 3.1635                 | 0.4590                 | 11.4744                  | 0.0918                | 0.1168                |
| 35  | anthracene           | 4.5000               | 0.0003                  | 0.7092                 | 3.7905                 | 0.0054                 | 0.7528                   | 0.1100                | 0.1195                |
| 42  | benzene              | 1.6000               | 0.1092                  | 1.4758                 | 0.0150                 | 1.8873                 | 1.5665                   | 0.0004                | 0.0011                |
| 39  | benz(a)anthracene    | 3.8000               | 0.0000                  | 2.4543                 | 1.3457                 | 0.0002                 | 2.6052                   | 0.0390                | 0.0331                |
| 40  | benzo(a)pyrene       | 3.0000               | 0.0000                  | 1.7419                 | 1.2581                 | 0.0000                 | 1.8490                   | 0.0365                | 0.0280                |
| 50  | benzo(b)fluoranthene | 4.0000               | 0.0000                  | 2.4706                 | 1.5294                 | 0.0005                 | 2.6224                   | 0.0444                | 0.0341                |
| 51  | benzo(k)fluoranthene | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                 | 0.0000                   | 0.0000                | 0.0000                |
| 38  | chrysene             | 3.3000               | 0.0000                  | 0.2542                 | 3.0458                 | 0.0001                 | 0.2699                   | 0.0884                | 0.0750                |
| 44  | ethylbenzene         | 0.2900               | 0.0016                  | 0.2850                 | 0.0034                 | 0.0275                 | 0.3025                   | 0.0001                | 0.0002                |
| 36  | fluoranthene         | 9.7000               | 0.0011                  | 6.5176                 | 3.1813                 | 0.0182                 | 6.9183                   | 0.0923                | 0.0884                |
| 33  | 9h-fluorene          | 7.1000               | 0.0048                  | 5.6557                 | 1.4395                 | 0.0822                 | 6.0034                   | 0.0418                | 0.0487                |
| 28  | naphthalene          | 110.0000             | 0.4888                  | 103.4335               | 6.0777                 | 8.4455                 | 109.7915                 | 0.1764                | 0.2664                |
| 34  | phenanthrene         | 16.0000              | 0.0057                  | 12.8933                | 3.1010                 | 0.0978                 | 13.6859                  | 0.0900                | 0.0978                |
| 37  | pyrene               | 6.0000               | 0.0006                  | 3.4054                 | 2.5940                 | 0.0095                 | 3.6147                   | 0.0753                | 0.0721                |
| 43  | toluene              | 4.5000               | 0.0894                  | 4.3633                 | 0.0473                 | 1.5438                 | 4.6315                   | 0.0014                | 0.0029                |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.140000,  
NAPL Volume Frac.(l/l), 0.000083,  
Soil Volume Frac.(l/l), 0.859917,  
Porosity (Volume Frac.), 0.140083,

Bulk Density (kg/l), 2.4189,  
NAPL Density (kg/l), 0.8999,

NAPL Saturation (%),

0.0595,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/6/2009 12:46:57 PM

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Sample Name Identification, CT-SO-B02-20

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.2000,

NAPLANAL ANALYSIS RESULTS:  
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ID#	Name	Total mass, (mg/kg)*	Mass, in water, (mg/kg)*	Mass in, in soil, (mg/kg)*	Mass, in NAPL, (mg/kg)*	Conc., in water, (mg/L)	Sorbed, in soil, (mg/kg)^	Conc., in NAPL, (kg/L)	Mole fraction, in NAPL,
29	1-methylnaphthalene	29.0000	0.0233	20.9235	8.0532	0.2697	22.9211	0.0090	0.0105
30	2-methylnaphthalene	72.0000	0.0576	51.7218	20.2207	0.6666	56.6598	0.0225	0.0262
32	acenaphthene	3.8000	0.0011	0.5571	3.2417	0.0133	0.6103	0.0036	0.0039
31	acenaphthylene	64.0000	0.0237	6.2752	57.7011	0.2750	6.8743	0.0641	0.0700
35	anthracene	26.0000	0.0001	0.1539	25.8460	0.0012	0.1686	0.0287	0.0268
42	benzene	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
39	benz(a)anthracene	23.0000	0.0000	1.2618	21.7382	0.0001	1.3823	0.0242	0.0176
40	benzo(a)pyrene	16.0000	0.0000	0.6753	15.3247	0.0000	0.7398	0.0170	0.0112
50	benzo(b)fluoranthene	24.0000	0.0000	1.1736	22.8264	0.0002	1.2856	0.0254	0.0167
51	benzo(k)fluoranthene	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
38	chrysene	20.0000	0.0000	0.0530	19.9470	0.0000	0.0580	0.0222	0.0161
44	ethylbenzene	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
36	fluoranthene	57.0000	0.0009	3.4890	53.5101	0.0101	3.8221	0.0595	0.0488
33	9h-fluorene	34.0000	0.0049	3.7783	30.2168	0.0567	4.1391	0.0336	0.0335
28	naphthalene	660.0000	1.6828	231.2919	427.0252	19.4903	253.3741	0.4746	0.6148
34	phenanthrene	82.0000	0.0065	9.5820	72.4115	0.0750	10.4968	0.0805	0.0750
37	pyrene	33.0000	0.0003	1.3234	31.6762	0.0038	1.4498	0.0352	0.0289
43	toluene	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.200000,
NAPL Volume Frac.(l/l), 0.002084,
Soil Volume Frac.(l/l), 0.797916,
Porosity (Volume Frac.), 0.202084,

Bulk Density (kg/l), 2.3164,
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

1.0313,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

ID#	CAS#	Name	Molecular Weight (g/mol)	Density (kg/l)	Henry's Constant (atm-L/mol)	Water Solubility (mg/L)	Koc (ml/g)
29	90-12-0	1-methylnaphthalene	142.2000	0.9000	0.2600	25.8000	8500.0000
30	91-57-6	2-methylnaphthalene	142.2000	0.9000	0.5180	25.4000	8500.0000
32	83-32-9	acenaphthene	154.2100	0.9000	0.0920	3.4200	4600.0000
31	208-96-8	acenaphthylene	152.2000	0.9000	1.4800	3.9300	2500.0000
35	120-12-7	anthracene	178.2400	0.9000	1.0200	0.0450	14000.0000
42	71-43-2	benzene	78.1100	0.8740	5.5900	1750.0000	83.0000
39	56-55-3	benz(a)anthracene	228.3000	0.9000	0.0012	0.0057	138000.0000
40	50-32-8	benzo(a)pyrene	252.3200	0.9000	0.0016	0.0012	550000.0000
50	205-99-2	benzo(b)fluoranthene	252.3200	0.9000	0.0119	0.0140	55000.0000
51	207-08-9	benzo(k)fluoranthene	252.3200	0.9000	0.0394	0.0043	55000.0000
38	218-01-9	chrysene	228.3000	0.9000	0.0011	0.0018	20000.0000
44	100-41-4	ethylbenzene	106.1700	0.8670	6.4300	152.0000	1100.0000
36	206-44-0	fluoranthene	202.2600	0.9000	0.0065	0.2060	38000.0000
33	86-73-7	9h-fluorene	166.2200	0.9000	0.0642	1.6900	7300.0000
28	91-20-3	naphthalene	128.1800	0.9000	1.1500	31.7000	1300.0000
34	85-01-8	phenanthrene	178.2400	0.9000	0.1590	1.0000	14000.0000
37	129-00-0	pyrene	202.2600	0.9000	0.0050	0.1320	38000.0000
43	108-88-3	toluene	92.1400	0.8620	6.3700	535.0000	300.0000

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time:

8/6/2009 1:00:41 PM

Sample Name Identification, CT-SO-B03-10
Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.1400,

NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.0097, 0.0000, 0.0097, 0.0000, 0.0001, 0.0103, 0.0001,  
30, 2-methylnaphthalene, 0.0190, 0.0000, 0.0190, 0.0000, 0.0002, 0.0202, 0.0001,  
32, acenaphthene, 0.0038, 0.0000, 0.0038, 0.0000, 0.0001, 0.0040, 0.0004,  
31, acenaphthylene, 0.0150, 0.0000, 0.0150, 0.0000, 0.0006, 0.0159, 0.0023,  
35, anthracene, 0.0260, 0.0000, 0.0260, 0.0000, 0.0002, 0.0276, 0.0611,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 0.1400, 0.0000, 0.1400, 0.0000, 0.0000, 0.1486, 0.0264,  
40, benzo(a)pyrene, 0.1100, 0.0000, 0.1100, 0.0000, 0.0000, 0.1168, 0.0247,  
50, benzo(b)fluoranthene, 0.2200, 0.0000, 0.2200, 0.0000, 0.0000, 0.2335, 0.0423,  
51, benzo(k)fluoranthene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
38, chrysene, 0.1700, 0.0000, 0.1700, 0.0000, 0.0001, 0.1804, 0.6995,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 0.3200, 0.0001, 0.3199, 0.0000, 0.0009, 0.3396, 0.0605,  
33, 9h-fluorene, 0.0082, 0.0000, 0.0082, 0.0000, 0.0001, 0.0087, 0.0010,  
28, naphthalene, 0.1600, 0.0008, 0.1592, 0.0000, 0.0130, 0.1690, 0.0057,  
34, phenanthrene, 0.1600, 0.0001, 0.1599, 0.0000, 0.0012, 0.1698, 0.0169,  
37, pyrene, 0.2000, 0.0000, 0.2000, 0.0000, 0.0006, 0.2123, 0.0591,  
17, trichloroethylene (tce),  
0.0026, 0.0001, 0.0025, 0.0000, 0.0021, 0.0026, 0.0000,  
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(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.860000,  
Water Volume Frac.(l/l), 0.140000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.140000,

Bulk Density (kg/l), 2.4190,

Dilution Factor (Vol. fac.), 12.9554,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                    | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|-------------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene     | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene     | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenapthene             | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenapthylene           | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene              | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene                 | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene       | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 1380000.0000  |
| 40  | 50-32-8  | benzo(a)pyrene          | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 5500000.0000  |
| 50  | 205-99-2 | benzo(b)fluoranthene    | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 550000.0000   |
| 51  | 207-08-9 | benzo(k)fluoranthene    | 252.3200                    | 0.9000            | 0.0394                          | 0.0043                     | 550000.0000   |
| 38  | 218-01-9 | chrysene                | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 200000.0000   |
| 44  | 100-41-4 | ethylbenzene            | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene            | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 33  | 86-73-7  | 9h-fluorene             | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 28  | 91-20-3  | naphthalene             | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene            | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene                  | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 17  | 79-01-6  | trichloroethylene (tce) | 131.5000                    | 1.4600            | 9.3700                          | 1384.9000                  | 126.0000      |

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/6/2009 1:19:36 PM

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Sample Name Identification, CT-SO-B03-20

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1600,

NAPLANAL ANALYSIS RESULTS:  
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ID#	Name	Total mass, (mg/kg)*	Mass, in water, (mg/kg)*	Mass in, in soil, (mg/kg)*	Mass, in NAPL, (mg/kg)*	Conc., in water, (mg/L)	Sorbed, in soil, (mg/kg)^	Conc., in NAPL, (kg/L)	Mole fraction, in NAPL,
29	1-methylnaphthalene	8.8000	0.0065	7.6504	1.1432	0.0965	8.2042	0.0026	0.0037
30	2-methylnaphthalene	13.0000	0.0095	11.2786	1.7119	0.1423	12.0951	0.0039	0.0056
32	acenaphthene	4.0000	0.0019	1.2268	2.7713	0.0286	1.3156	0.0063	0.0084
31	acenaphthylene	33.0000	0.0208	7.2167	25.7625	0.3096	7.7392	0.0586	0.0788
35	anthracene	30.0000	0.0002	0.4532	29.5466	0.0035	0.4860	0.0672	0.0771
42	benzene	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
39	benz(a)anthracene	24.0000	0.0000	3.1216	20.8784	0.0002	3.3476	0.0475	0.0426
40	benzo(a)pyrene	18.0000	0.0000	1.8349	16.1651	0.0000	1.9677	0.0368	0.0298
50	benzo(b)fluoranthene	25.0000	0.0000	2.9235	22.0765	0.0006	3.1351	0.0502	0.0407
51	benzo(k)fluoranthene	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
38	chrysene	21.0000	0.0000	0.1427	20.8573	0.0001	0.1531	0.0474	0.0425
44	ethylbenzene	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
36	fluoranthene	79.0000	0.0022	11.3597	67.6382	0.0321	12.1820	0.1539	0.1556
33	9h-fluorene	32.0000	0.0077	7.7938	24.1985	0.1145	8.3580	0.0550	0.0677
28	naphthalene	81.0000	0.2604	47.0298	33.7098	3.8796	50.4344	0.0767	0.1224
34	phenanthrene	120.0000	0.0157	30.5002	89.4841	0.2336	32.7082	0.2035	0.2336
37	pyrene	44.0000	0.0008	4.2750	39.7242	0.0121	4.5845	0.0904	0.0914
17	trichloroethylene (tce)	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.160000,
NAPL Volume Frac.(l/l), 0.001048,
Soil Volume Frac.(l/l), 0.838952,
Porosity (Volume Frac.), 0.161048,

Bulk Density (kg/l), 2.3842,

NAPL Density (kg/l), 0.9000,
 NAPL Saturation (%), 0.6508,

Numerical Accuracy Information

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 The solution converged in 214 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25°C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                    | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|-------------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene     | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene     | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenapthene             | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenapthylene           | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene              | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene                 | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene       | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 1380000.0000  |
| 40  | 50-32-8  | benzo(a)pyrene          | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 5500000.0000  |
| 50  | 205-99-2 | benzo(b)fluoranthene    | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 550000.0000   |
| 51  | 207-08-9 | benzo(k)fluoranthene    | 252.3200                    | 0.9000            | 0.0394                          | 0.0043                     | 550000.0000   |
| 38  | 218-01-9 | chrysene                | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 200000.0000   |
| 44  | 100-41-4 | ethylbenzene            | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene            | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 33  | 86-73-7  | 9h-fluorene             | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 28  | 91-20-3  | naphthalene             | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene            | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene                  | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 17  | 79-01-6  | trichloroethylene (tce) | 131.5000                    | 1.4600            | 9.3700                          | 1384.9000                  | 126.0000      |

\*\*\*\*\*

END OF RECORDS



NAPLANAL Version 1.0.0

Date and Time: 8/6/2009 1:26:08 PM

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Sample Name Identification, CT-SO-B03-22  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1400,  
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NAPLANAL ANALYSIS RESULTS:

~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,
, , (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/L), (mg/kg)^, ,
====, =====, =====, =====, =====, =====, =====, =====, =====,
29, 1-methylnaphthalene, 3.6000, 0.0026, 3.5974, 0.0000, 0.0449, 3.8184, 0.0020,
30, 2-methylnaphthalene, 8.5000, 0.0061, 8.4939, 0.0000, 0.1061, 9.0156, 0.0047,
32, acenaphthene, 0.7700, 0.0010, 0.7690, 0.0000, 0.0177, 0.8162, 0.0059,
31, acenaphthylene, 7.1000, 0.0174, 7.0826, 0.0000, 0.3007, 7.5177, 0.0868,
35, anthracene, 1.7000, 0.0007, 1.6993, 0.0000, 0.0129, 1.8036, 0.3249,
42, benzene, 0.3200, 0.0221, 0.2979, 0.0000, 0.3810, 0.3163, 0.0002,
39, benz(a)anthracene, 0.5400, 0.0000, 0.5400, 0.0000, 0.0000, 0.5732, 0.0083,
40, benzo(a)pyrene, 0.3800, 0.0000, 0.3800, 0.0000, 0.0000, 0.4033, 0.0069,
50, benzo(b)fluoranthene, 0.5500, 0.0000, 0.5500, 0.0000, 0.0001, 0.5838, 0.0086,
51, benzo(k)fluoranthene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
38, chrysene, 0.5600, 0.0000, 0.5600, 0.0000, 0.0003, 0.5944, 0.1874,
44, ethylbenzene, 0.0950, 0.0005, 0.0945, 0.0000, 0.0091, 0.1003, 0.0001,
36, fluoranthene, 3.2000, 0.0005, 3.1995, 0.0000, 0.0089, 3.3960, 0.0492,
33, 9h-fluorene, 5.3000, 0.0045, 5.2955, 0.0000, 0.0770, 5.6209, 0.0517,
28, naphthalene, 43.0000, 0.2022, 42.7978, 0.0000, 3.4944, 45.4268, 0.1251,
34, phenanthrene, 11.0000, 0.0048, 10.9952, 0.0000, 0.0834, 11.6706, 0.0946,
37, pyrene, 1.8000, 0.0003, 1.7997, 0.0000, 0.0050, 1.9103, 0.0432,
43, toluene, 0.4300, 0.0086, 0.4214, 0.0000, 0.1491, 0.4473, 0.0003,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.860000,
Water Volume Frac.(l/l), 0.140000,
NAPL Volume Frac.(l/l), 0.000000,
Porosity (Volume Frac.), 0.140000,

Bulk Density (kg/l), 2.4190,

Dilution Factor (Vol. fac.), 0.1348,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
          (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
29  90-12-0  1-methylnaphthalene  142.2000  0.9000  0.2600  25.8000  8500.0000
30  91-57-6  2-methylnaphthalene  142.2000  0.9000  0.5180  25.4000  8500.0000
32  83-32-9  acenaphthene  154.2100  0.9000  0.0920  3.4200  4600.0000
31  208-96-8  acenaphthylene  152.2000  0.9000  1.4800  3.9300  2500.0000
35  120-12-7  anthracene  178.2400  0.9000  1.0200  0.0450  14000.0000
42  71-43-2  benzene  78.1100  0.8740  5.5900  1750.0000  83.0000
39  56-55-3  benz(a)anthracene  228.3000  0.9000  0.0012  0.0057  138000.0000
40  50-32-8  benzo(a)pyrene  252.3200  0.9000  0.0016  0.0012  550000.0000
50  205-99-2  benzo(b)fluoranthene  252.3200  0.9000  0.0119  0.0140  55000.0000
51  207-08-9  benzo(k)fluoranthene  252.3200  0.9000  0.0394  0.0043  55000.0000
38  218-01-9  chrysene  228.3000  0.9000  0.0011  0.0018  20000.0000
44  100-41-4  ethylbenzene  106.1700  0.8670  6.4300  152.0000  1100.0000
36  206-44-0  fluoranthene  202.2600  0.9000  0.0065  0.2060  38000.0000
33  86-73-7  9h-fluorene  166.2200  0.9000  0.0642  1.6900  7300.0000
28  91-20-3  naphthalene  128.1800  0.9000  1.1500  31.7000  1300.0000
34  85-01-8  phenanthrene  178.2400  0.9000  0.1590  1.0000  14000.0000
37  129-00-0  pyrene  202.2600  0.9000  0.0050  0.1320  38000.0000
43  108-88-3  toluene  92.1400  0.8620  6.3700  535.0000  300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/6/2009 3:58:09 PM

Sample Name Identification, CT-SO-B04-10
Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.2400,

NAPLANAL ANALYSIS RESULTS:

ID#	Name	Total mass, (mg/kg)*	Mass, in water, (mg/kg)*	Mass, in soil, (mg/kg)*	Mass, in NAPL, (mg/kg)*	Conc., in water, (mg/L)	Sorbed, in soil, (mg/kg)^	Mole fraction, in NAPL,
29	1-methylnaphthalene,	0.9700,	0.0014,	0.9686,	0.0000,	0.0128,	1.0841,	0.0005,
30	2-methylnaphthalene,	1.9000,	0.0027,	1.8973,	0.0000,	0.0250,	2.1234,	0.0011,
32	acenapthene,	0.1600,	0.0004,	0.1596,	0.0000,	0.0039,	0.1786,	0.0012,
31	acenapthylene,	0.7700,	0.0037,	0.7663,	0.0000,	0.0343,	0.8577,	0.0095,
35	anthracene,	0.7000,	0.0006,	0.6994,	0.0000,	0.0056,	0.7827,	0.1347,
42	benzene,	0.5100,	0.0640,	0.4460,	0.0000,	0.6013,	0.4991,	0.0004,
39	benz(a)anthracene,	1.5000,	0.0000,	1.5000,	0.0000,	0.0001,	1.6787,	0.0231,
40	benzo(a)pyrene,	1.2000,	0.0000,	1.2000,	0.0000,	0.0000,	1.3430,	0.0221,
50	benzo(b)fluoranthene,	2.2000,	0.0000,	2.2000,	0.0000,	0.0004,	2.4621,	0.0347,
51	benzo(k)fluoranthene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
38	chrysene,	1.8000,	0.0001,	1.7999,	0.0000,	0.0010,	2.0144,	0.6067,
44	ethylbenzene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
36	fluoranthene,	3.1000,	0.0010,	3.0990,	0.0000,	0.0091,	3.4683,	0.0480,
33	9h-fluorene,	0.6300,	0.0010,	0.6290,	0.0000,	0.0096,	0.7039,	0.0062,
28	naphthalene,	9.6000,	0.0872,	9.5128,	0.0000,	0.8190,	10.6464,	0.0280,
34	phenanthrene,	3.5000,	0.0030,	3.4970,	0.0000,	0.0280,	3.9137,	0.0303,
37	pyrene,	2.2000,	0.0007,	2.1993,	0.0000,	0.0065,	2.4614,	0.0532,
43	toluene,	0.2800,	0.0107,	0.2693,	0.0000,	0.1005,	0.3014,	0.0002,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.760000,
Water Volume Frac.(l/l), 0.240000,
NAPL Volume Frac.(l/l), 0.000000,
Porosity (Volume Frac.), 0.240000,

Bulk Density (kg/l), 2.2540,

Dilution Factor (Vol. fac.), 0.0844,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name                Molecular Weight   Density   Henry's Constant   Water Solubility   Koc
              (g/mol)          (kg/l)          (atm-L/mol)      (mg/L)          (ml/g)
=====
 29  90-12-0    1-methylnaphthalene  142.2000          0.9000          0.2600            25.8000           8500.0000
 30  91-57-6    2-methylnaphthalene  142.2000          0.9000          0.5180            25.4000           8500.0000
 32  83-32-9    acenaphthene         154.2100          0.9000          0.0920            3.4200            4600.0000
 31  208-96-8    acenaphthylene       152.2000          0.9000          1.4800            3.9300            2500.0000
 35  120-12-7    anthracene            178.2400          0.9000          1.0200            0.0450            14000.0000
 42  71-43-2    benzene               78.1100           0.8740          5.5900            1750.0000          83.0000
 39  56-55-3    benz(a)anthracene    228.3000          0.9000          0.0012            0.0057            138000.0000
 40  50-32-8    benzo(a)pyrene       252.3200          0.9000          0.0016            0.0012            550000.0000
 50  205-99-2    benzo(b)fluoranthene 252.3200          0.9000          0.0119            0.0140            55000.0000
 51  207-08-9    benzo(k)fluoranthene 252.3200          0.9000          0.0394            0.0043            55000.0000
 38  218-01-9    chrysene              228.3000          0.9000          0.0011            0.0018            20000.0000
 44  100-41-4    ethylbenzene         106.1700          0.8670          6.4300            152.0000           1100.0000
 36  206-44-0    fluoranthene         202.2600          0.9000          0.0065            0.2060            38000.0000
 33  86-73-7    9h-fluorene         166.2200          0.9000          0.0642            1.6900            7300.0000
 28  91-20-3    naphthalene          128.1800          0.9000          1.1500            31.7000            1300.0000
 34  85-01-8    phenanthrene         178.2400          0.9000          0.1590            1.0000            14000.0000
 37  129-00-0    pyrene               202.2600          0.9000          0.0050            0.1320            38000.0000
 43  108-88-3    toluene              92.1400           0.8620          6.3700            535.0000           300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time:

8/6/2009 4:01:59 PM

Sample Name Identification, CT-SO-B04-14
 Model used: Liquid saturated & water moisture content known,
 Water moisture content (Volume Frac.), 0.1800,

NAPLANAL ANALYSIS RESULTS:

ID#	Name	Total mass, (mg/kg)*	Mass, in water, (mg/kg)*	Mass, in soil, (mg/kg)*	Mass, in NAPL, (mg/kg)*	Conc., in water, (mg/L)	Sorbed, in soil, (mg/kg)^	Mole fraction, in NAPL,
29,	1-methylnaphthalene,	0.4800,	0.0005,	0.4795,	0.0000,	0.0061,	0.5193,	0.0003,
30,	2-methylnaphthalene,	1.2000,	0.0012,	1.1988,	0.0000,	0.0153,	1.2981,	0.0008,
32,	acenapthene,	0.1200,	0.0002,	0.1198,	0.0000,	0.0028,	0.1297,	0.0011,
31,	acenapthylene,	0.6700,	0.0022,	0.6678,	0.0000,	0.0289,	0.7231,	0.0100,
35,	anthracene,	0.8600,	0.0005,	0.8595,	0.0000,	0.0066,	0.9307,	0.2008,
42,	benzene,	0.7500,	0.0681,	0.6819,	0.0000,	0.8897,	0.7384,	0.0007,
39,	benz(a)anthracene,	1.4000,	0.0000,	1.4000,	0.0000,	0.0001,	1.5160,	0.0262,
40,	benzo(a)pyrene,	1.2000,	0.0000,	1.2000,	0.0000,	0.0000,	1.2994,	0.0268,
50,	benzo(b)fluoranthene,	1.6000,	0.0000,	1.6000,	0.0000,	0.0003,	1.7325,	0.0306,
51,	benzo(k)fluoranthene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,
38,	chrysene,	1.3000,	0.0001,	1.2999,	0.0000,	0.0007,	1.4076,	0.5316,
44,	ethylbenzene,	0.0400,	0.0003,	0.0397,	0.0000,	0.0039,	0.0430,	0.0000,
36,	fluoranthene,	2.8000,	0.0006,	2.7994,	0.0000,	0.0080,	3.0313,	0.0526,
33,	9h-fluorene,	0.6000,	0.0007,	0.5993,	0.0000,	0.0089,	0.6490,	0.0072,
28,	naphthalene,	7.1000,	0.0450,	7.0550,	0.0000,	0.5876,	7.6394,	0.0252,
34,	phenanthrene,	2.6000,	0.0015,	2.5985,	0.0000,	0.0201,	2.8137,	0.0273,
37,	pyrene,	2.0000,	0.0004,	1.9996,	0.0000,	0.0057,	2.1652,	0.0587,
43,	toluene,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,	0.0000,

(mg/kg)* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.820000,
 Water Volume Frac.(l/l), 0.180000,
 NAPL Volume Frac.(l/l), 0.000000,
 Porosity (Volume Frac.), 0.180000,

Bulk Density (kg/l), 2.3530,

Dilution Factor (Vol. fac.), 0.3595,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
 Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID#      CAS#      Name      Molecular Weight  Density  Henry's Constant  Water Solubility  Koc
          (g/mol)      (kg/l)      (atm-L/mol)      (mg/L)      (ml/g)
=====
 29  90-12-0    1-methylnaphthalene    142.2000    0.9000    0.2600    25.8000    8500.0000
 30  91-57-6    2-methylnaphthalene    142.2000    0.9000    0.5180    25.4000    8500.0000
 32  83-32-9    acenaphthene           154.2100    0.9000    0.0920    3.4200    4600.0000
 31  208-96-8    acenaphthylene         152.2000    0.9000    1.4800    3.9300    2500.0000
 35  120-12-7    anthracene              178.2400    0.9000    1.0200    0.0450    14000.0000
 42  71-43-2     benzene                  78.1100    0.8740    5.5900    1750.0000    83.0000
 39  56-55-3     benz(a)anthracene       228.3000    0.9000    0.0012    0.0057    138000.0000
 40  50-32-8     benzo(a)pyrene          252.3200    0.9000    0.0016    0.0012    550000.0000
 50  205-99-2     benzo(b)fluoranthene    252.3200    0.9000    0.0119    0.0140    55000.0000
 51  207-08-9     benzo(k)fluoranthene    252.3200    0.9000    0.0394    0.0043    55000.0000
 38  218-01-9     chrysene                 228.3000    0.9000    0.0011    0.0018    20000.0000
 44  100-41-4     ethylbenzene             106.1700    0.8670    6.4300    152.0000    1100.0000
 36  206-44-0     fluoranthene             202.2600    0.9000    0.0065    0.2060    38000.0000
 33  86-73-7     9h-fluorene              166.2200    0.9000    0.0642    1.6900    7300.0000
 28  91-20-3     naphthalene              128.1800    0.9000    1.1500    31.7000    1300.0000
 34  85-01-8     phenanthrene             178.2400    0.9000    0.1590    1.0000    14000.0000
 37  129-00-0     pyrene                   202.2600    0.9000    0.0050    0.1320    38000.0000
 43  108-88-3     toluene                   92.1400    0.8620    6.3700    535.0000    300.0000
=====
  
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/6/2009 4:41:41 PM

Sample Name Identification, CT-SO-B04-18

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.2400,

NAPLANAL ANALYSIS RESULTS:

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| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 30.0000              | 0.0315                  | 22.3833                | 7.5853                 | 0.2950                  | 25.0758                   | 0.0092                 | 0.0114                  |
| 30  | 2-methylnaphthalene  | 74.0000              | 0.0773                  | 54.9931                | 18.9296                | 0.7248                  | 61.6083                   | 0.0229                 | 0.0285                  |
| 32  | acenaphthene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 31  | acenaphthylene       | 49.0000              | 0.0257                  | 5.3842                 | 43.5901                | 0.2413                  | 6.0318                    | 0.0528                 | 0.0614                  |
| 35  | anthracene           | 28.0000              | 0.0002                  | 0.1881                 | 27.8117                | 0.0015                  | 0.2107                    | 0.0337                 | 0.0334                  |
| 42  | benzene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 31.0000              | 0.0000                  | 1.9173                 | 29.0827                | 0.0002                  | 2.1480                    | 0.0352                 | 0.0273                  |
| 40  | benzo(a)pyrene       | 24.0000              | 0.0000                  | 1.1439                 | 22.8560                | 0.0000                  | 1.2816                    | 0.0277                 | 0.0194                  |
| 50  | benzo(b)fluoranthene | 33.0000              | 0.0000                  | 1.8206                 | 31.1793                | 0.0004                  | 2.0396                    | 0.0378                 | 0.0265                  |
| 51  | benzo(k)fluoranthene | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 38  | chrysene             | 27.0000              | 0.0000                  | 0.0812                 | 26.9188                | 0.0000                  | 0.0910                    | 0.0326                 | 0.0253                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 73.0000              | 0.0016                  | 5.0332                 | 67.9653                | 0.0148                  | 5.6386                    | 0.0823                 | 0.0720                  |
| 33  | 9h-fluorene          | 40.0000              | 0.0081                  | 4.9732                 | 35.0186                | 0.0763                  | 5.5715                    | 0.0424                 | 0.0452                  |
| 28  | naphthalene          | 470.0000             | 1.6392                  | 178.3847               | 289.9761               | 15.3725                 | 199.8428                  | 0.3511                 | 0.4849                  |
| 34  | phenanthrene         | 110.0000             | 0.0123                  | 14.3705                | 95.6172                | 0.1150                  | 16.0991                   | 0.1158                 | 0.1150                  |
| 37  | pyrene               | 49.0000              | 0.0007                  | 2.2198                 | 46.7795                | 0.0065                  | 2.4868                    | 0.0566                 | 0.0496                  |
| 43  | toluene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.240000,  
NAPL Volume Frac.(l/l), 0.001859,  
Soil Volume Frac.(l/l), 0.758141,  
Porosity (Volume Frac.), 0.241859,

Bulk Density (kg/l), 2.2507,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

0.7686,

Numerical Accuracy Information

The solution converged in 75 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS



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Sample Name Identification,          CT-SO-B05-8
Model used:  Unsaturated sample,
Porosity (Volume Frac.),           0.3000,
Water moisture content (Volume Frac.), 0.0990,
Fraction organic carbon (foc),      0.0100,
    
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NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Mass, Conc., Sorbed, Conc., Mole
frac., , mass, in water, in air, in soil, in NAPL, in water, in soil, in air, in
NAPL, , (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/kg)*, (mg/L), (mg/kg)^, (mg/L), ,
, , =====, =====, =====, =====, =====, =====, =====, =====, =====, =====
29, 1-methylnaphthalene, 0.0650, 0.0000, 0.0000, 0.0650, 0.0000, 0.0008, 0.0684, 0.0000, 0.0001,
30, 2-methylnaphthalene, 0.1400, 0.0001, 0.0000, 0.1399, 0.0000, 0.0017, 0.1474, 0.0000, 0.0003,
32, acenapthene, 0.0110, 0.0000, 0.0000, 0.0110, 0.0000, 0.0003, 0.0116, 0.0000, 0.0003,
31, acenapthylene, 0.0240, 0.0001, 0.0000, 0.0239, 0.0000, 0.0010, 0.0252, 0.0001, 0.0010,
35, anthracene, 0.0810, 0.0000, 0.0000, 0.0810, 0.0000, 0.0006, 0.0853, 0.0000, 0.0547,
42, benzene, 0.1200, 0.0070, 0.0033, 0.1096, 0.0000, 0.1391, 0.1155, 0.0323, 0.0003,
39, benz(a)anthracene, 0.5100, 0.0000, 0.0000, 0.5100, 0.0000, 0.0000, 0.5373, 0.0000, 0.0276,
40, benzo(a)pyrene, 0.3600, 0.0000, 0.0000, 0.3600, 0.0000, 0.0000, 0.3793, 0.0000, 0.0232,
50, benzo(b)fluoranthene, 0.7900, 0.0000, 0.0000, 0.7900, 0.0000, 0.0002, 0.8323, 0.0000, 0.0436,
51, benzo(k)fluoranthene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
38, chrysene, 0.5800, 0.0000, 0.0000, 0.5800, 0.0000, 0.0003, 0.6110, 0.0000, 0.6854,
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
36, fluoranthene, 1.1000, 0.0002, 0.0000, 1.0998, 0.0000, 0.0030, 1.1587, 0.0000, 0.0598,
33, 9h-fluorene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,
28, naphthalene, 0.7400, 0.0030, 0.0003, 0.7367, 0.0000, 0.0597, 0.7761, 0.0029, 0.0076,
34, phenanthrene, 0.6800, 0.0003, 0.0000, 0.6797, 0.0000, 0.0051, 0.7161, 0.0000, 0.0207,
37, pyrene, 0.8800, 0.0001, 0.0000, 0.8799, 0.0000, 0.0024, 0.9270, 0.0000, 0.0746,
43, toluene, 0.3200, 0.0055, 0.0030, 0.3115, 0.0000, 0.1094, 0.3281, 0.0290, 0.0008,

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(mg/kg)\* --- mg per kg of soil sample (wet soil)  
(mg/kg)^ --- mg per kg of solid (dry soil)

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Soil Volume Frac.(l/l), 0.700000,
Water Volume Frac.(l/l), 0.099000,
NAPL Volume Frac.(l/l), 0.000000,
Air Volume Frac.(l/l), 0.201000,
Porosity (Volume Frac.), 0.300000,

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Bulk Density (kg/l), 1.9543,

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Dilution Factor (Vol. fac.), 3.0380,

No NAPL (i.e., NAPL Saturation Equals Zero)

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The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)

Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenapthene          | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenapthylene        | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 1380000.0000  |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 5500000.0000  |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 550000.0000   |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                    | 0.9000            | 0.0394                          | 0.0043                     | 550000.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 200000.0000   |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time:

8/31/2009 9:07:51 AM

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Sample Name Identification, CT-SO-B05-16  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1600,  
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NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.4300, 0.0004, 0.4296, 0.0000, 0.0054, 0.4605, 0.0028,  
30, 2-methylnaphthalene, 1.1000, 0.0009, 1.0991, 0.0000, 0.0139, 1.1781, 0.0073,  
32, acenaphthene, 0.0700, 0.0001, 0.0699, 0.0000, 0.0016, 0.0749, 0.0064,  
31, acenaphthylene, 0.6600, 0.0019, 0.6581, 0.0000, 0.0282, 0.7054, 0.0962,  
35, anthracene, 0.0880, 0.0000, 0.0880, 0.0000, 0.0007, 0.0943, 0.2005,  
42, benzene, 1.7000, 0.1355, 1.5645, 0.0000, 2.0204, 1.6770, 0.0155,  
39, benz(a)anthracene, 0.0460, 0.0000, 0.0460, 0.0000, 0.0000, 0.0493, 0.0084,  
40, benzo(a)pyrene, 0.0320, 0.0000, 0.0320, 0.0000, 0.0000, 0.0343, 0.0070,  
50, benzo(b)fluoranthene, 0.0590, 0.0000, 0.0590, 0.0000, 0.0000, 0.0632, 0.0110,  
51, benzo(k)fluoranthene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
38, chrysene, 0.0520, 0.0000, 0.0520, 0.0000, 0.0000, 0.0557, 0.2075,  
44, ethylbenzene, 0.1100, 0.0007, 0.1093, 0.0000, 0.0106, 0.1171, 0.0009,  
36, fluoranthene, 0.1400, 0.0000, 0.1400, 0.0000, 0.0004, 0.1500, 0.0257,  
33, 9h-fluorene, 0.4200, 0.0004, 0.4196, 0.0000, 0.0062, 0.4497, 0.0488,  
28, naphthalene, 7.7000, 0.0423, 7.6577, 0.0000, 0.6314, 8.2081, 0.2669,  
34, phenanthrene, 0.4500, 0.0002, 0.4498, 0.0000, 0.0034, 0.4821, 0.0461,  
37, pyrene, 0.1100, 0.0000, 0.1100, 0.0000, 0.0003, 0.1179, 0.0315,  
43, toluene, 2.0000, 0.0468, 1.9532, 0.0000, 0.6979, 2.0936, 0.0175,  
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(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.840000,  
Water Volume Frac.(l/l), 0.160000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.160000,

Bulk Density (kg/l), 2.3860,

Dilution Factor (Vol. fac.), 12.3997,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
 29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
 30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
 32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
 31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
 35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
 42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
 39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
 40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
 50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
 51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
 38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
 44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
 36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
 33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
 28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
 34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
 37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
 43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
=====

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END OF RECORDS

Sample Name Identification, CT-SO-B05-20

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.2000,

NAPLANAL ANALYSIS RESULTS:

Table with 10 columns: ID#, Name, Total mass, Mass in water, Mass in soil, Mass in NAPL, Conc. in water, Sorbed, Conc. in NAPL, Mole fraction in NAPL. Lists various chemical compounds and their concentrations.

(mg/kg)\* --- mg per kg of soil sample (wet soil)
(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.200000,
NAPL Volume Frac.(l/l), 0.002039,
Soil Volume Frac.(l/l), 0.797961,
Porosity (Volume Frac.), 0.202039,

Bulk Density (kg/l), 2.3164,
NAPL Density (kg/l), 0.8998,

NAPL Saturation (%),

1.0093,

Numerical Accuracy Information

The solution converged in 102 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

\*\*\*\*\*  
END OF RECORDS

**NAPLANAL MODEL  
OFFSHORE INVESTIGATION OUTPUT FILES**

NAPLANAL Version 1.0.0

Date and Time:

8/3/2009 11:37:22 AM

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Sample Name Identification, BH-SED-01-00  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5300,

NAPLANAL ANALYSIS RESULTS:

-----  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.1700, 0.0008, 0.1692, 0.0000, 0.0028, 0.2411, 0.0002,  
30, 2-methylnaphthalene, 0.3400, 0.0017, 0.3383, 0.0000, 0.0057, 0.4823, 0.0004,  
32, acenapthene, 0.0730, 0.0007, 0.0723, 0.0000, 0.0022, 0.1031, 0.0013,  
31, acenapthylene, 0.2300, 0.0038, 0.2262, 0.0000, 0.0129, 0.3224, 0.0065,  
35, anthracene, 0.3100, 0.0009, 0.3091, 0.0000, 0.0031, 0.4406, 0.1375,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 0.6800, 0.0000, 0.6800, 0.0000, 0.0001, 0.9693, 0.0242,  
40, benzo(a)pyrene, 1.1000, 0.0000, 1.1000, 0.0000, 0.0000, 1.5681, 0.0467,  
50, benzo(b)fluoranthene, 1.3000, 0.0001, 1.2999, 0.0000, 0.0003, 1.8530, 0.0473,  
51, benzo(k)fluoranthene, 0.4400, 0.0000, 0.4400, 0.0000, 0.0001, 0.6272, 0.0521,  
33, 9h-fluorene, 0.2000, 0.0012, 0.1988, 0.0000, 0.0039, 0.2835, 0.0045,  
38, chrysene, 0.6800, 0.0001, 0.6799, 0.0000, 0.0005, 0.9692, 0.5292,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 1.3000, 0.0015, 1.2985, 0.0000, 0.0049, 1.8511, 0.0465,  
28, naphthalene, 3.7000, 0.1173, 3.5827, 0.0000, 0.3929, 5.1073, 0.0244,  
34, phenanthrene, 0.6100, 0.0018, 0.6082, 0.0000, 0.0062, 0.8669, 0.0122,  
37, pyrene, 1.2000, 0.0013, 1.1987, 0.0000, 0.0045, 1.7087, 0.0670,  
43, toluene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.470000,  
Water Volume Frac.(l/l), 0.530000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.530000,

Bulk Density (kg/l), 1.7755,

Dilution Factor (Vol. fac.), 0.9659,

No NAPL (i.e., NAPL Saturation Equals Zero)



The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
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END OF RECORDS

Sample Name Identification, BH-SED-01

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.5300,

NAPLANAL ANALYSIS RESULTS:

Table with 10 columns: ID#, Name, Total mass, Mass in water, Mass in soil, Mass in NAPL, Conc. in water, Sorbed, Conc. in NAPL, Mole fraction in NAPL. Rows include various polycyclic aromatic hydrocarbons like 1-methylnaphthalene, 2-methylnaphthalene, acenaphthene, etc.

(mg/kg)\* --- mg per kg of soil sample (wet soil)
(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.530000,
NAPL Volume Frac.(l/l), 0.000005,
Soil Volume Frac.(l/l), 0.469995,
Porosity (Volume Frac.), 0.530005,

Bulk Density (kg/l), 1.7755,
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0009,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/24/2009 9:00:19 AM

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Sample Name Identification, BH-SED-02-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.3400,

NAPLANAL ANALYSIS RESULTS:  
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL,<br>, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|---------------------------------|
| 29,  | 1-methylnaphthalene,  | 2.6000,                      | 0.0056,                         | 2.4693,                           | 0.1251,                        | 0.0347,                        | 2.9499,                          | 0.0009,                       | 0.0013,                         |
| 30,  | 2-methylnaphthalene,  | 5.0000,                      | 0.0109,                         | 4.7450,                           | 0.2441,                        | 0.0667,                        | 5.6686,                          | 0.0018,                       | 0.0026,                         |
| 32,  | acenaphthene,         | 1.5000,                      | 0.0036,                         | 0.8475,                           | 0.6489,                        | 0.0220,                        | 1.0125,                          | 0.0047,                       | 0.0064,                         |
| 31,  | acenaphthylene,       | 5.5000,                      | 0.0193,                         | 2.4800,                           | 3.0007,                        | 0.1185,                        | 2.9627,                          | 0.0215,                       | 0.0302,                         |
| 35,  | anthracene,           | 7.3000,                      | 0.0004,                         | 0.3160,                           | 6.9835,                        | 0.0027,                        | 0.3775,                          | 0.0501,                       | 0.0599,                         |
| 42,  | benzene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 39,  | benz(a)anthracene,    | 12.0000,                     | 0.0001,                         | 3.6731,                           | 8.3268,                        | 0.0003,                        | 4.3880,                          | 0.0597,                       | 0.0558,                         |
| 40,  | benzo(a)pyrene,       | 13.0000,                     | 0.0000,                         | 3.2613,                           | 9.7386,                        | 0.0001,                        | 3.8961,                          | 0.0699,                       | 0.0590,                         |
| 50,  | benzo(b)fluoranthene, | 19.0000,                     | 0.0002,                         | 5.3378,                           | 13.6620,                       | 0.0012,                        | 6.3767,                          | 0.0980,                       | 0.0828,                         |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 33,  | 9h-fluorene,          | 4.4000,                      | 0.0057,                         | 2.1411,                           | 2.2532,                        | 0.0350,                        | 2.5578,                          | 0.0162,                       | 0.0207,                         |
| 38,  | chrysene,             | 11.0000,                     | 0.0000,                         | 0.2177,                           | 10.7823,                       | 0.0001,                        | 0.2600,                          | 0.0773,                       | 0.0722,                         |
| 44,  | ethylbenzene,         | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 36,  | fluoranthene,         | 33.0000,                     | 0.0056,                         | 10.9320,                          | 22.0624,                       | 0.0344,                        | 13.0597,                         | 0.1583,                       | 0.1668,                         |
| 28,  | naphthalene,          | 85.0000,                     | 1.0109,                         | 67.5725,                          | 16.4166,                       | 6.2096,                        | 80.7243,                         | 0.1178,                       | 0.1959,                         |
| 34,  | phenanthrene,         | 20.0000,                     | 0.0139,                         | 10.0210,                          | 9.9651,                        | 0.0855,                        | 11.9714,                         | 0.0715,                       | 0.0855,                         |
| 37,  | pyrene,               | 28.0000,                     | 0.0035,                         | 6.7469,                           | 21.2496,                       | 0.0212,                        | 8.0601,                          | 0.1524,                       | 0.1607,                         |
| 43,  | toluene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.340000,  
NAPL Volume Frac.(l/l), 0.000291,  
Soil Volume Frac.(l/l), 0.659709,  
Porosity (Volume Frac.), 0.340291,

Bulk Density (kg/l), 2.0885,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0856,

Numerical Accuracy Information

~~~~~  
 The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 1380000.0000
40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 5500000.0000
50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 550000.0000
3 75-27-4 bromodichloromethane 163.8000 1.9700 2.0600 6716.9000 61.0000
33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 200000.0000
44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
=====

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END OF RECORDS

Sample Name Identification, BH-SED-02-4

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.1500,

NAPLANAL ANALYSIS RESULTS:

Table with 10 columns: ID#, Name, Total mass, Mass in water, Mass in soil, Mass in NAPL, Conc. in water, Sorbed, Conc. in NAPL, Mole fraction in NAPL. Lists various chemical compounds and their concentrations.

(mg/kg)\* --- mg per kg of soil sample (wet soil)
(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.150000,
NAPL Volume Frac.(l/l), 0.000256,
Soil Volume Frac.(l/l), 0.849744,
Porosity (Volume Frac.), 0.150256,

Bulk Density (kg/l), 2.4021,
NAPL Density (kg/l), 0.8997,

NAPL Saturation (%),

0.1706,

Numerical Accuracy Information

The solution converged in 6 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with  
CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenapthene          | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenapthylene        | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 138000.0000   |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 550000.0000   |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 55000.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                    | 0.9000            | 0.0394                          | 0.0043                     | 55000.0000    |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 20000.0000    |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

\*\*\*\*\*  
END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 11:46:30 AM

Sample Name Identification, BH-SED-03A

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.4300,

NAPLANAL ANALYSIS RESULTS:

| ID# | Name                 | Total mass, (mg/kg)* | Mass, in water, (mg/kg)* | Mass in, in soil, (mg/kg)* | Mass, in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|--------------------------|----------------------------|-------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 2.1000               | 0.0069                   | 2.0591                     | 0.0340                  | 0.0311                  | 2.6455                    | 0.0008                 | 0.0012                  |
| 30  | 2-methylnaphthalene  | 3.0000               | 0.0099                   | 2.9409                     | 0.0493                  | 0.0445                  | 3.7783                    | 0.0012                 | 0.0018                  |
| 32  | acenaphthene         | 1.3000               | 0.0064                   | 1.0354                     | 0.2582                  | 0.0289                  | 1.3302                    | 0.0061                 | 0.0085                  |
| 31  | acenaphthylene       | 2.3000               | 0.0186                   | 1.6364                     | 0.6450                  | 0.0841                  | 2.1024                    | 0.0151                 | 0.0214                  |
| 35  | anthracene           | 2.3000               | 0.0006                   | 0.2805                     | 2.0190                  | 0.0026                  | 0.3603                    | 0.0474                 | 0.0572                  |
| 42  | benzene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 6.7000               | 0.0001                   | 3.8539                     | 2.8460                  | 0.0004                  | 4.9514                    | 0.0668                 | 0.0629                  |
| 40  | benzo(a)pyrene       | 7.6000               | 0.0000                   | 3.8525                     | 3.7475                  | 0.0001                  | 4.9496                    | 0.0880                 | 0.0750                  |
| 50  | benzo(b)fluoranthene | 11.0000              | 0.0003                   | 5.9984                     | 5.0013                  | 0.0014                  | 7.7065                    | 0.1174                 | 0.1001                  |
| 51  | benzo(k)fluoranthene | 1.3000               | 0.0000                   | 0.3500                     | 0.9500                  | 0.0001                  | 0.4496                    | 0.0223                 | 0.0190                  |
| 33  | 9h-fluorene          | 1.0000               | 0.0029                   | 0.7425                     | 0.2546                  | 0.0131                  | 0.9540                    | 0.0060                 | 0.0077                  |
| 38  | chrysene             | 5.5000               | 0.0000                   | 0.3210                     | 5.1790                  | 0.0002                  | 0.4124                    | 0.1216                 | 0.1145                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 12.0000              | 0.0054                   | 7.2369                     | 4.7577                  | 0.0245                  | 9.2977                    | 0.1117                 | 0.1188                  |
| 28  | naphthalene          | 90.0000              | 1.7902                   | 81.7407                    | 6.4691                  | 8.0783                  | 105.0176                  | 0.1518                 | 0.2548                  |
| 34  | phenanthrene         | 4.8000               | 0.0074                   | 3.6200                     | 1.1727                  | 0.0332                  | 4.6508                    | 0.0275                 | 0.0332                  |
| 37  | pyrene               | 9.8000               | 0.0036                   | 4.8354                     | 4.9610                  | 0.0163                  | 6.2123                    | 0.1164                 | 0.1239                  |
| 43  | toluene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.430000,  
NAPL Volume Frac.(l/l), 0.000083,  
Soil Volume Frac.(l/l), 0.569917,  
Porosity (Volume Frac.), 0.430083,

Bulk Density (kg/l), 1.9404,  
NAPL Density (kg/l), 0.9000,



NAPL Saturation (%),

0.0192,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphtene          | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphtylene        | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time:

8/3/2009 11:49:48 AM

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Sample Name Identification, BH-SED-3A-12  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.4000,

NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.6500, 0.0019, 0.6481, 0.0000, 0.0095, 0.8111, 0.0011,  
30, 2-methylnaphthalene, 1.6000, 0.0047, 1.5953, 0.0000, 0.0235, 1.9966, 0.0027,  
32, acenaphthene, 0.0930, 0.0005, 0.0925, 0.0000, 0.0025, 0.1158, 0.0022,  
31, acenaphthylene, 0.3200, 0.0032, 0.3168, 0.0000, 0.0159, 0.3965, 0.0119,  
35, anthracene, 0.1800, 0.0003, 0.1797, 0.0000, 0.0016, 0.2249, 0.1051,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 0.4600, 0.0000, 0.4600, 0.0000, 0.0000, 0.5757, 0.0215,  
40, benzo(a)pyrene, 0.4600, 0.0000, 0.4600, 0.0000, 0.0000, 0.5757, 0.0257,  
50, benzo(b)fluoranthene, 0.6700, 0.0000, 0.6700, 0.0000, 0.0002, 0.8385, 0.0321,  
51, benzo(k)fluoranthene, 0.2800, 0.0000, 0.2800, 0.0000, 0.0001, 0.3504, 0.0436,  
33, 9h-fluorene, 0.3500, 0.0012, 0.3488, 0.0000, 0.0060, 0.4365, 0.0104,  
38, chrysene, 0.3800, 0.0000, 0.3800, 0.0000, 0.0002, 0.4755, 0.3888,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 0.7000, 0.0005, 0.6995, 0.0000, 0.0023, 0.8755, 0.0329,  
28, naphthalene, 29.0000, 0.5505, 28.4495, 0.0000, 2.7390, 35.6065, 0.2543,  
34, phenanthrene, 0.5400, 0.0010, 0.5390, 0.0000, 0.0048, 0.6746, 0.0142,  
37, pyrene, 0.7300, 0.0005, 0.7295, 0.0000, 0.0024, 0.9130, 0.0536,  
43, toluene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.600000,  
Water Volume Frac.(l/l), 0.400000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.400000,

Bulk Density (kg/l), 1.9900,

Dilution Factor (Vol. fac.), 1.9433,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:02:32 PM

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Sample Name Identification, BH-SED-3B-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.6400,

NAPLANAL ANALYSIS RESULTS:  
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| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 30  | 2-methylnaphthalene  | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 32  | acenaphthene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 31  | acenaphthylene       | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 35  | anthracene           | 36.0000              | 0.0001                  | 0.0130                 | 35.9869                | 0.0002                  | 0.0223                    | 0.0043                 | 0.0035                  |
| 42  | benzene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 57.0000              | 0.0000                  | 0.2007                 | 56.7992                | 0.0000                  | 0.3431                    | 0.0068                 | 0.0044                  |
| 40  | benzo(a)pyrene       | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 50  | benzo(b)fluoranthene | 57.0000              | 0.0000                  | 0.1779                 | 56.8221                | 0.0001                  | 0.3040                    | 0.0068                 | 0.0039                  |
| 51  | benzo(k)fluoranthene | 30.0000              | 0.0000                  | 0.0288                 | 29.9712                | 0.0000                  | 0.0493                    | 0.0036                 | 0.0021                  |
| 33  | 9h-fluorene          | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 38  | chrysene             | 54.0000              | 0.0000                  | 0.0087                 | 53.9913                | 0.0000                  | 0.0149                    | 0.0065                 | 0.0041                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 140.0000             | 0.0010                  | 0.5536                 | 139.4454               | 0.0025                  | 0.9462                    | 0.0167                 | 0.0121                  |
| 28  | naphthalene          | 7200.0000            | 12.2908                 | 229.4781               | 6958.2312              | 30.1701                 | 392.2107                  | 0.8319                 | 0.9517                  |
| 34  | phenanthrene         | 88.0000              | 0.0035                  | 0.7033                 | 87.2932                | 0.0086                  | 1.2021                    | 0.0104                 | 0.0086                  |
| 37  | pyrene               | 110.0000             | 0.0005                  | 0.2791                 | 109.7204               | 0.0013                  | 0.4771                    | 0.0131                 | 0.0095                  |
| 43  | toluene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.640000,  
NAPL Volume Frac.(l/l), 0.013141,  
Soil Volume Frac.(l/l), 0.346859,  
Porosity (Volume Frac.), 0.653141,

Bulk Density (kg/l), 1.5710,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

2.0120,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS

Sample Name Identification, BH-SED-03B-2  
 Model used: Liquid saturated & water moisture content known,  
 Water moisture content (Volume Frac.), 0.5400,

NAPLANAL ANALYSIS RESULTS:

| ID# | Name                 | Total mass, (mg/kg)* | Mass, in water, (mg/kg)* | Mass, in soil, (mg/kg)* | Mass, in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|--------------------------|-------------------------|-------------------------|-------------------------|---------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 0.4300               | 0.0022                   | 0.4278                  | 0.0000                  | 0.0073                  | 0.6173                    | 0.0003                  |
| 30  | 2-methylnaphthalene  | 0.7500               | 0.0039                   | 0.7461                  | 0.0000                  | 0.0127                  | 1.0766                    | 0.0005                  |
| 32  | acenapthene          | 0.1500               | 0.0014                   | 0.1486                  | 0.0000                  | 0.0047                  | 0.2144                    | 0.0015                  |
| 31  | acenapthylene        | 0.3000               | 0.0052                   | 0.2948                  | 0.0000                  | 0.0170                  | 0.4254                    | 0.0047                  |
| 35  | anthracene           | 0.3600               | 0.0011                   | 0.3589                  | 0.0000                  | 0.0037                  | 0.5178                    | 0.0891                  |
| 42  | benzene              | 0.0000               | 0.0000                   | 0.0000                  | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                  |
| 39  | benz(a)anthracene    | 1.6000               | 0.0001                   | 1.5999                  | 0.0000                  | 0.0002                  | 2.3087                    | 0.0318                  |
| 40  | benzo(a)pyrene       | 1.7000               | 0.0000                   | 1.7000                  | 0.0000                  | 0.0000                  | 2.4531                    | 0.0403                  |
| 50  | benzo(b)fluoranthene | 2.5000               | 0.0002                   | 2.4998                  | 0.0000                  | 0.0007                  | 3.6072                    | 0.0508                  |
| 3   | bromodichloromethane | 0.0000               | 0.0000                   | 0.0000                  | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                  |
| 33  | 9h-fluorene          | 0.2000               | 0.0012                   | 0.1988                  | 0.0000                  | 0.0039                  | 0.2869                    | 0.0025                  |
| 38  | chrysene             | 1.1000               | 0.0002                   | 1.0998                  | 0.0000                  | 0.0008                  | 1.5869                    | 0.4779                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                   | 0.0000                  | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                  |
| 36  | fluoranthene         | 2.2000               | 0.0026                   | 2.1974                  | 0.0000                  | 0.0083                  | 3.1709                    | 0.0439                  |
| 28  | naphthalene          | 51.0000              | 1.6806                   | 49.3194                 | 0.0000                  | 5.4744                  | 71.1672                   | 0.1872                  |
| 34  | phenanthrene         | 0.6400               | 0.0020                   | 0.6380                  | 0.0000                  | 0.0066                  | 0.9206                    | 0.0071                  |
| 37  | pyrene               | 2.0000               | 0.0023                   | 1.9977                  | 0.0000                  | 0.0076                  | 2.8826                    | 0.0623                  |
| 43  | toluene              | 0.0000               | 0.0000                   | 0.0000                  | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)  
 (mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.460000,  
 Water Volume Frac.(l/l), 0.540000,  
 NAPL Volume Frac.(l/l), 0.000000,  
 Porosity (Volume Frac.), 0.540000,

Bulk Density (kg/l), 1.7590,

Dilution Factor (Vol. fac.), 0.0841,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)

Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 1380000.0000  |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 5500000.0000  |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 550000.0000   |
| 3   | 75-27-4  | bromodichloromethane | 163.8000                    | 1.9700            | 2.0600                          | 6716.9000                  | 61.0000       |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 200000.0000   |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/24/2009 10:55:20 AM

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Sample Name Identification, BH-SED-3C-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.6900,

NAPLANAL ANALYSIS RESULTS:  
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL,<br>, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|---------------------------------|
| 29,  | 1-methylnaphthalene,  | 4.1000,                      | 0.0356,                         | 3.5941,                           | 0.4703,                        | 0.0778,                        | 6.6172,                          | 0.0023,                       | 0.0030,                         |
| 30,  | 2-methylnaphthalene,  | 5.9000,                      | 0.0511,                         | 5.1627,                           | 0.6862,                        | 0.1118,                        | 9.5052,                          | 0.0033,                       | 0.0044,                         |
| 32,  | acenapthene,          | 4.2000,                      | 0.0256,                         | 1.4017,                           | 2.7727,                        | 0.0561,                        | 2.5806,                          | 0.0134,                       | 0.0164,                         |
| 31,  | acenapthylene,        | 2.0000,                      | 0.0162,                         | 0.4808,                           | 1.5030,                        | 0.0354,                        | 0.8852,                          | 0.0073,                       | 0.0090,                         |
| 35,  | anthracene,           | 5.0000,                      | 0.0005,                         | 0.0861,                           | 4.9134,                        | 0.0011,                        | 0.1584,                          | 0.0238,                       | 0.0251,                         |
| 42,  | benzene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 39,  | benz(a)anthracene,    | 13.0000,                     | 0.0001,                         | 1.8958,                           | 11.1041,                       | 0.0003,                        | 3.4904,                          | 0.0537,                       | 0.0444,                         |
| 40,  | benzo(a)pyrene,       | 15.0000,                     | 0.0000,                         | 1.7211,                           | 13.2788,                       | 0.0001,                        | 3.1688,                          | 0.0643,                       | 0.0480,                         |
| 50,  | benzo(b)fluoranthene, | 21.0000,                     | 0.0004,                         | 2.7584,                           | 18.2412,                       | 0.0009,                        | 5.0785,                          | 0.0883,                       | 0.0660,                         |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 33,  | 9h-fluorene,          | 2.8000,                      | 0.0086,                         | 0.7506,                           | 2.0408,                        | 0.0189,                        | 1.3819,                          | 0.0099,                       | 0.0112,                         |
| 38,  | chrysene,             | 10.0000,                     | 0.0000,                         | 0.0775,                           | 9.9224,                        | 0.0001,                        | 0.1427,                          | 0.0480,                       | 0.0397,                         |
| 44,  | ethylbenzene,         | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 36,  | fluoranthene,         | 30.0000,                     | 0.0107,                         | 4.8259,                           | 25.1635,                       | 0.0234,                        | 8.8850,                          | 0.1218,                       | 0.1135,                         |
| 28,  | naphthalene,          | 190.0000,                    | 7.2609,                         | 112.2681,                         | 70.4709,                       | 15.8999,                       | 206.6987,                        | 0.3410,                       | 0.5016,                         |
| 34,  | phenanthrene,         | 6.9000,                      | 0.0116,                         | 1.9299,                           | 4.9585,                        | 0.0254,                        | 3.5532,                          | 0.0240,                       | 0.0254,                         |
| 37,  | pyrene,               | 23.0000,                     | 0.0056,                         | 2.5165,                           | 20.4779,                       | 0.0122,                        | 4.6332,                          | 0.0991,                       | 0.0924,                         |
| 43,  | toluene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.690000,  
NAPL Volume Frac.(l/l), 0.000312,  
Soil Volume Frac.(l/l), 0.309688,  
Porosity (Volume Frac.), 0.690312,

Bulk Density (kg/l), 1.5110,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0452,



Numerical Accuracy Information

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 The solution converged in 5 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
 29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
 30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
 32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
 31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
 35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
 42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
 39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 1380000.0000
 40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 5500000.0000
 50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 550000.0000
 3 75-27-4 bromodichloromethane 163.8000 1.9700 2.0600 6716.9000 61.0000
 33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
 38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 200000.0000
 44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
 36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
 28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
 34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
 37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
 43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/27/2009 4:13:36 PM

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Sample Name Identification, BH-SED-03C-2\_RE

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.6000,

NAPLANAL ANALYSIS RESULTS:  
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL,<br>, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|---------------------------------|
| ==,  | ====,                 | ====,                        | ====,                           | ====,                             | ====,                          | ====,                          | ====,                            | ====,                         | ====,                           |
| 29,  | 1-methylnaphthalene,  | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 30,  | 2-methylnaphthalene,  | 66.0000,                     | 0.1409,                         | 20.8923,                          | 44.9668,                       | 0.3877,                        | 32.9555,                         | 0.0150,                       | 0.0153,                         |
| 32,  | acenapthene,          | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 31,  | acenapthylene,        | 15.0000,                     | 0.0067,                         | 0.2904,                           | 14.7029,                       | 0.0183,                        | 0.4581,                          | 0.0049,                       | 0.0047,                         |
| 35,  | anthracene,           | 7.4000,                      | 0.0000,                         | 0.0080,                           | 7.3920,                        | 0.0001,                        | 0.0126,                          | 0.0025,                       | 0.0020,                         |
| 42,  | benzene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 39,  | benz(a)anthracene,    | 11.0000,                     | 0.0000,                         | 0.1148,                           | 10.8852,                       | 0.0000,                        | 0.1810,                          | 0.0036,                       | 0.0023,                         |
| 40,  | benzo(a)pyrene,       | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 50,  | benzo(b)fluoranthene, | 14.0000,                     | 0.0000,                         | 0.1295,                           | 13.8705,                       | 0.0000,                        | 0.2043,                          | 0.0046,                       | 0.0027,                         |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 33,  | 9h-fluorene,          | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 38,  | chrysene,             | 8.0000,                      | 0.0000,                         | 0.0039,                           | 7.9961,                        | 0.0000,                        | 0.0061,                          | 0.0027,                       | 0.0017,                         |
| 44,  | ethylbenzene,         | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 36,  | fluoranthene,         | 24.0000,                     | 0.0004,                         | 0.2809,                           | 23.7187,                       | 0.0012,                        | 0.4431,                          | 0.0079,                       | 0.0057,                         |
| 28,  | naphthalene,          | 2800.0000,                   | 11.0136,                        | 249.8138,                         | 2539.1726,                     | 30.3119,                       | 394.0551,                        | 0.8464,                       | 0.9562,                         |
| 34,  | phenanthrene,         | 20.0000,                     | 0.0019,                         | 0.4694,                           | 19.5287,                       | 0.0053,                        | 0.7404,                          | 0.0065,                       | 0.0053,                         |
| 37,  | pyrene,               | 18.0000,                     | 0.0002,                         | 0.1356,                           | 17.8642,                       | 0.0006,                        | 0.2139,                          | 0.0060,                       | 0.0043,                         |
| 43,  | toluene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.600000,  
NAPL Volume Frac.(l/l), 0.004954,  
Soil Volume Frac.(l/l), 0.395046,  
Porosity (Volume Frac.), 0.604954,

Bulk Density (kg/l), 1.6513,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.8189,

Numerical Accuracy Information

~~~~~  
 The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
 29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
 30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
 32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
 31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
 35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
 42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
 39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 1380000.0000
 40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 5500000.0000
 50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 550000.0000
 3 75-27-4 bromodichloromethane 163.8000 1.9700 2.0600 6716.9000 61.0000
 33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
 38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 200000.0000
 44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
 36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
 28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
 34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
 37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
 43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:08:39 PM

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Sample Name Identification, BH-SED-03D-2

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.6300,

NAPLANAL ANALYSIS RESULTS:  
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|----------------------------|
| 29,  | 1-methylnaphthalene,  | 0.3300,                      | 0.0025,                         | 0.3268,                           | 0.0007,                        | 0.0063,                        | 0.5368,                          | 0.0001,                       | 0.0002,                    |
| 30,  | 2-methylnaphthalene,  | 0.8100,                      | 0.0061,                         | 0.8021,                           | 0.0019,                        | 0.0155,                        | 1.3174,                          | 0.0004,                       | 0.0006,                    |
| 32,  | acenapthene,          | 0.1500,                      | 0.0020,                         | 0.1431,                           | 0.0049,                        | 0.0051,                        | 0.2350,                          | 0.0010,                       | 0.0015,                    |
| 31,  | acenapthylene,        | 0.9300,                      | 0.0221,                         | 0.8608,                           | 0.0471,                        | 0.0566,                        | 1.4139,                          | 0.0091,                       | 0.0144,                    |
| 35,  | anthracene,           | 0.8700,                      | 0.0020,                         | 0.4343,                           | 0.4337,                        | 0.0051,                        | 0.7133,                          | 0.0835,                       | 0.1132,                    |
| 42,  | benzene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 39,  | benz(a)anthracene,    | 2.8000,                      | 0.0001,                         | 2.5397,                           | 0.2602,                        | 0.0003,                        | 4.1716,                          | 0.0501,                       | 0.0530,                    |
| 40,  | benzo(a)pyrene,       | 3.4000,                      | 0.0000,                         | 2.9957,                           | 0.4043,                        | 0.0001,                        | 4.9206,                          | 0.0778,                       | 0.0746,                    |
| 50,  | benzo(b)fluoranthene, | 3.6000,                      | 0.0004,                         | 3.2264,                           | 0.3732,                        | 0.0010,                        | 5.2995,                          | 0.0719,                       | 0.0688,                    |
| 51,  | benzo(k)fluoranthene, | 1.7000,                      | 0.0001,                         | 1.2348,                           | 0.4650,                        | 0.0004,                        | 2.0282,                          | 0.0895,                       | 0.0858,                    |
| 33,  | 9h-fluorene,          | 0.6300,                      | 0.0052,                         | 0.5964,                           | 0.0284,                        | 0.0134,                        | 0.9796,                          | 0.0055,                       | 0.0079,                    |
| 38,  | chrysene,             | 2.2000,                      | 0.0002,                         | 0.6792,                           | 1.5205,                        | 0.0006,                        | 1.1157,                          | 0.2928,                       | 0.3099,                    |
| 44,  | ethylbenzene,         | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 36,  | fluoranthene,         | 5.8000,                      | 0.0090,                         | 5.3070,                           | 0.4840,                        | 0.0229,                        | 8.7170,                          | 0.0932,                       | 0.1114,                    |
| 28,  | naphthalene,          | 5.3000,                      | 0.2470,                         | 4.9981,                           | 0.0549,                        | 0.6315,                        | 8.2096,                          | 0.0106,                       | 0.0199,                    |
| 34,  | phenanthrene,         | 1.4000,                      | 0.0061,                         | 1.3339,                           | 0.0599,                        | 0.0157,                        | 2.1910,                          | 0.0115,                       | 0.0157,                    |
| 37,  | pyrene,               | 4.3000,                      | 0.0064,                         | 3.7586,                           | 0.5350,                        | 0.0162,                        | 6.1738,                          | 0.1030,                       | 0.1231,                    |
| 43,  | toluene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.630000,  
NAPL Volume Frac.(l/l), 0.000008,  
Soil Volume Frac.(l/l), 0.369992,  
Porosity (Volume Frac.), 0.630008,

Bulk Density (kg/l), 1.6105,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0013,

Numerical Accuracy Information

~~~~~  
 The solution converged in 16 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
 ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
 29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
 30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
 32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
 31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
 35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
 42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
 39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
 40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
 50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
 51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
 33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
 38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
 44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
 36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
 28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
 34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
 37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
 43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:12:50 PM

-----  
Sample Name Identification, BH-SED-03E-2  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.6300,  
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NAPLANAL ANALYSIS RESULTS:

~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.0660, 0.0005, 0.0655, 0.0000, 0.0013, 0.1076, 0.0001,  
30, 2-methylnaphthalene, 0.1500, 0.0011, 0.1489, 0.0000, 0.0029, 0.2445, 0.0002,  
32, acenaphthene, 0.0310, 0.0004, 0.0306, 0.0000, 0.0011, 0.0502, 0.0005,  
31, acenaphthylene, 0.2000, 0.0050, 0.1950, 0.0000, 0.0128, 0.3203, 0.0049,  
35, anthracene, 0.2300, 0.0011, 0.2289, 0.0000, 0.0027, 0.3761, 0.0905,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 0.8300, 0.0000, 0.8300, 0.0000, 0.0001, 1.3632, 0.0263,  
40, benzo(a)pyrene, 1.1000, 0.0000, 1.1000, 0.0000, 0.0000, 1.8068, 0.0415,  
50, benzo(b)fluoranthene, 1.1000, 0.0001, 1.0999, 0.0000, 0.0003, 1.8066, 0.0356,  
51, benzo(k)fluoranthene, 0.5800, 0.0001, 0.5799, 0.0000, 0.0002, 0.9526, 0.0611,  
33, 9h-fluorene, 0.1100, 0.0010, 0.1090, 0.0000, 0.0025, 0.1791, 0.0022,  
38, chrysene, 0.9200, 0.0003, 0.9197, 0.0000, 0.0008, 1.5106, 0.6364,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 0.9800, 0.0017, 0.9783, 0.0000, 0.0042, 1.6070, 0.0311,  
28, naphthalene, 0.8700, 0.0410, 0.8290, 0.0000, 0.1047, 1.3617, 0.0050,  
34, phenanthrene, 0.2900, 0.0013, 0.2887, 0.0000, 0.0034, 0.4742, 0.0051,  
37, pyrene, 1.2000, 0.0020, 1.1980, 0.0000, 0.0052, 1.9677, 0.0595,  
43, toluene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.370000,  
Water Volume Frac.(l/l), 0.630000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.630000,

Bulk Density (kg/l), 1.6105,

Dilution Factor (Vol. fac.), 0.5165,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
 29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
 30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
 32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
 31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
 35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
 42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
 39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
 40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
 50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
 51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
 33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
 38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
 44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
 36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
 28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
 34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
 37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
 43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/24/2009 12:14:16 PM

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Sample Name Identification, BH-SED-04-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.4500,

NAPLANAL ANALYSIS RESULTS:

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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|----------------------------|
| ==,  | ====,                 | ====,                        | ====,                           | ====,                             | ====,                          | ====,                          | ====,                            | ====,                         | ====,                      |
| 29,  | 1-methylnaphthalene,  | 2.5000,                      | 0.0089,                         | 2.4448,                           | 0.0464,                        | 0.0376,                        | 3.1998,                          | 0.0010,                       | 0.0015,                    |
| 30,  | 2-methylnaphthalene,  | 5.9000,                      | 0.0210,                         | 5.7680,                           | 0.1111,                        | 0.0888,                        | 7.5494,                          | 0.0024,                       | 0.0035,                    |
| 32,  | acenapthene,          | 0.8300,                      | 0.0043,                         | 0.6417,                           | 0.1839,                        | 0.0183,                        | 0.8400,                          | 0.0039,                       | 0.0053,                    |
| 31,  | acenapthylene,        | 2.2000,                      | 0.0185,                         | 1.5014,                           | 0.6801,                        | 0.0786,                        | 1.9651,                          | 0.0144,                       | 0.0200,                    |
| 35,  | anthracene,           | 3.1000,                      | 0.0007,                         | 0.3342,                           | 2.7650,                        | 0.0031,                        | 0.4375,                          | 0.0585,                       | 0.0694,                    |
| 42,  | benzene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 39,  | benz(a)anthracene,    | 6.9000,                      | 0.0001,                         | 3.7323,                           | 3.1676,                        | 0.0004,                        | 4.8851,                          | 0.0671,                       | 0.0621,                    |
| 40,  | benzo(a)pyrene,       | 8.5000,                      | 0.0000,                         | 4.0134,                           | 4.4866,                        | 0.0001,                        | 5.2529,                          | 0.0950,                       | 0.0796,                    |
| 50,  | benzo(b)fluoranthene, | 12.0000,                     | 0.0003,                         | 6.1279,                           | 5.8718,                        | 0.0015,                        | 8.0205,                          | 0.1243,                       | 0.1042,                    |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 33,  | 9h-fluorene,          | 1.8000,                      | 0.0054,                         | 1.2874,                           | 0.5072,                        | 0.0231,                        | 1.6850,                          | 0.0107,                       | 0.0137,                    |
| 38,  | chrysene,             | 5.4000,                      | 0.0000,                         | 0.2763,                           | 5.1237,                        | 0.0002,                        | 0.3616,                          | 0.1085,                       | 0.1005,                    |
| 44,  | ethylbenzene,         | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 36,  | fluoranthene,         | 12.0000,                     | 0.0056,                         | 6.8324,                           | 5.1621,                        | 0.0235,                        | 8.9426,                          | 0.1093,                       | 0.1142,                    |
| 28,  | naphthalene,          | 97.0000,                     | 2.0670,                         | 87.0185,                          | 7.9145,                        | 8.7611,                        | 113.8945,                        | 0.1675,                       | 0.2764,                    |
| 34,  | phenanthrene,         | 7.4000,                      | 0.0119,                         | 5.3838,                           | 2.0043,                        | 0.0503,                        | 7.0466,                          | 0.0424,                       | 0.0503,                    |
| 37,  | pyrene,               | 8.3000,                      | 0.0031,                         | 3.8075,                           | 4.4894,                        | 0.0131,                        | 4.9835,                          | 0.0950,                       | 0.0994,                    |
| 43,  | toluene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.450000,  
NAPL Volume Frac.(l/l), 0.000090,  
Soil Volume Frac.(l/l), 0.549910,  
Porosity (Volume Frac.), 0.450090,

Bulk Density (kg/l), 1.9073,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0200,



Numerical Accuracy Information

~~~~~  
 The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
3 75-27-4 bromodichloromethane 163.8000 1.9700 2.0600 6716.9000 61.0000
33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 3800.0000
28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 1400.0000
37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 3800.0000
43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
=====

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 END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:16:51 PM

Sample Name Identification, BH-SED-04-8

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5900,

NAPLANAL ANALYSIS RESULTS:

~~~~~

| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|-------------------------|--------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 37.0000              | 0.0533                  | 8.1705                 | 28.7762                | 0.1501                  | 12.7598                  | 0.0056                 | 0.0058                  |
| 30  | 2-methylnaphthalene  | 93.0000              | 0.1324                  | 20.2882                | 72.5795                | 0.3728                  | 31.6838                  | 0.0142                 | 0.0147                  |
| 32  | acenaphthene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                   | 0.0000                 | 0.0000                  |
| 31  | acenaphthylene       | 25.0000              | 0.0065                  | 0.2936                 | 24.6999                | 0.0183                  | 0.4584                   | 0.0048                 | 0.0047                  |
| 35  | anthracene           | 17.0000              | 0.0000                  | 0.0111                 | 16.9889                | 0.0001                  | 0.0173                   | 0.0033                 | 0.0027                  |
| 42  | benzene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                   | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 35.0000              | 0.0000                  | 0.2206                 | 34.7794                | 0.0000                  | 0.3445                   | 0.0068                 | 0.0044                  |
| 40  | benzo(a)pyrene       | 29.0000              | 0.0000                  | 0.1390                 | 28.8610                | 0.0000                  | 0.2171                   | 0.0056                 | 0.0033                  |
| 50  | benzo(b)fluoranthene | 33.0000              | 0.0000                  | 0.1844                 | 32.8156                | 0.0001                  | 0.2879                   | 0.0064                 | 0.0037                  |
| 51  | benzo(k)fluoranthene | 12.0000              | 0.0000                  | 0.0207                 | 11.9793                | 0.0000                  | 0.0323                   | 0.0023                 | 0.0014                  |
| 33  | 9h-fluorene          | 19.0000              | 0.0019                  | 0.2561                 | 18.7420                | 0.0055                  | 0.4000                   | 0.0037                 | 0.0032                  |
| 38  | chrysene             | 28.0000              | 0.0000                  | 0.0081                 | 27.9919                | 0.0000                  | 0.0127                   | 0.0055                 | 0.0035                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                   | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 67.0000              | 0.0007                  | 0.4740                 | 66.5253                | 0.0019                  | 0.7403                   | 0.0130                 | 0.0095                  |
| 28  | naphthalene          | 4400.0000            | 10.4633                 | 245.3027               | 4144.2340              | 29.4682                 | 383.0869                 | 0.8112                 | 0.9296                  |
| 34  | phenanthrene         | 44.0000              | 0.0025                  | 0.6272                 | 43.3703                | 0.0070                  | 0.9795                   | 0.0085                 | 0.0070                  |
| 37  | pyrene               | 46.0000              | 0.0003                  | 0.2091                 | 45.7906                | 0.0009                  | 0.3265                   | 0.0090                 | 0.0065                  |
| 43  | toluene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                   | 0.0000                 | 0.0000                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.590000,  
NAPL Volume Frac.(l/l), 0.008489,  
Soil Volume Frac.(l/l), 0.401511,  
Porosity (Volume Frac.), 0.598489,

Bulk Density (kg/l), 1.6616,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

1.4185,

Numerical Accuracy Information

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The solution converged in 2 iterations with residual less than 1.0E-6.

-----  
The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with  
CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

=====

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenapthene          | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenapthylene        | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 1380000.0000  |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 5500000.0000  |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 550000.0000   |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                    | 0.9000            | 0.0394                          | 0.0043                     | 550000.0000   |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 200000.0000   |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:19:27 PM

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Sample Name Identification, BH-SED-05-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.3800,

NAPLANAL ANALYSIS RESULTS:  
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| ID# | Name                 | Total mass, (mg/kg)* | Mass, in water, (mg/kg)* | Mass in, in soil, (mg/kg)* | Mass, in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|--------------------------|----------------------------|-------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 30  | 2-methylnaphthalene  | 3.4000               | 0.0080                   | 2.9412                     | 0.4508                  | 0.0426                  | 3.6239                    | 0.0010                 | 0.0017                  |
| 32  | acenaphthene         | 4.5000               | 0.0069                   | 1.3695                     | 3.1236                  | 0.0367                  | 1.6873                    | 0.0072                 | 0.0107                  |
| 31  | acenaphthylene       | 9.0000               | 0.0181                   | 1.9507                     | 7.0313                  | 0.0961                  | 2.4034                    | 0.0162                 | 0.0245                  |
| 35  | anthracene           | 22.0000              | 0.0005                   | 0.3292                     | 21.6703                 | 0.0029                  | 0.4056                    | 0.0498                 | 0.0644                  |
| 42  | benzene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 48.0000              | 0.0001                   | 6.1908                     | 41.8091                 | 0.0006                  | 7.6277                    | 0.0961                 | 0.0970                  |
| 40  | benzo(a)pyrene       | 48.0000              | 0.0000                   | 4.8506                     | 43.1494                 | 0.0001                  | 5.9764                    | 0.0992                 | 0.0906                  |
| 50  | benzo(b)fluoranthene | 53.0000              | 0.0003                   | 6.1450                     | 46.8547                 | 0.0014                  | 7.5712                    | 0.1077                 | 0.0983                  |
| 51  | benzo(k)fluoranthene | 19.0000              | 0.0000                   | 0.7357                     | 18.2643                 | 0.0002                  | 0.9065                    | 0.0420                 | 0.0383                  |
| 33  | 9h-fluorene          | 3.5000               | 0.0027                   | 0.8458                     | 2.6515                  | 0.0143                  | 1.0421                    | 0.0061                 | 0.0084                  |
| 38  | chrysene             | 40.0000              | 0.0000                   | 0.2692                     | 39.7307                 | 0.0002                  | 0.3317                    | 0.0913                 | 0.0921                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 88.0000              | 0.0076                   | 12.5485                    | 75.4438                 | 0.0407                  | 15.4610                   | 0.1735                 | 0.1975                  |
| 28  | naphthalene          | 50.0000              | 0.5115                   | 28.7098                    | 20.7787                 | 2.7210                  | 35.3732                   | 0.0478                 | 0.0858                  |
| 34  | phenanthrene         | 23.0000              | 0.0096                   | 5.8021                     | 17.1883                 | 0.0511                  | 7.1487                    | 0.0395                 | 0.0511                  |
| 37  | pyrene               | 59.0000              | 0.0035                   | 5.6822                     | 53.3143                 | 0.0184                  | 7.0011                    | 0.1226                 | 0.1396                  |
| 43  | toluene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.380000,  
NAPL Volume Frac.(l/l), 0.000879,  
Soil Volume Frac.(l/l), 0.619121,  
Porosity (Volume Frac.), 0.380879,

Bulk Density (kg/l), 2.0215,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

0.2308,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with  
CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 138000.0000   |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 550000.0000   |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 55000.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                    | 0.9000            | 0.0394                          | 0.0043                     | 55000.0000    |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 20000.0000    |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

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END OF RECORDS

Sample Name Identification, BH-SED-05-8

Model used: Liquid saturated & water moisture content known,  
 Water moisture content (Volume Frac.), 0.4600,

NAPLANAL ANALYSIS RESULTS:

| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 3.4000               | 0.0102                  | 2.6946                 | 0.6952                 | 0.0419                  | 3.5637                    | 0.0015                 | 0.0016                  |
| 30  | 2-methylnaphthalene  | 8.6000               | 0.0257                  | 6.7940                 | 1.7803                 | 0.1057                  | 8.9852                    | 0.0038                 | 0.0042                  |
| 32  | acenaphthene         | 9.7000               | 0.0138                  | 1.9768                 | 7.7094                 | 0.0568                  | 2.6144                    | 0.0166                 | 0.0166                  |
| 31  | acenaphthylene       | 1.8000               | 0.0032                  | 0.2508                 | 1.5459                 | 0.0133                  | 0.3317                    | 0.0033                 | 0.0034                  |
| 35  | anthracene           | 6.7000               | 0.0001                  | 0.0590                 | 6.6409                 | 0.0006                  | 0.0780                    | 0.0143                 | 0.0124                  |
| 42  | benzene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 7.1000               | 0.0000                  | 0.5659                 | 6.5341                 | 0.0001                  | 0.7484                    | 0.0141                 | 0.0095                  |
| 40  | benzo(a)pyrene       | 6.2000               | 0.0000                  | 0.3825                 | 5.8175                 | 0.0000                  | 0.5058                    | 0.0125                 | 0.0077                  |
| 50  | benzo(b)fluoranthene | 6.4000               | 0.0000                  | 0.4559                 | 5.9440                 | 0.0001                  | 0.6030                    | 0.0128                 | 0.0078                  |
| 51  | benzo(k)fluoranthene | 3.5000               | 0.0000                  | 0.0806                 | 3.4194                 | 0.0000                  | 0.1065                    | 0.0074                 | 0.0045                  |
| 33  | 9h-fluorene          | 7.2000               | 0.0050                  | 1.1312                 | 6.0638                 | 0.0205                  | 1.4961                    | 0.0131                 | 0.0121                  |
| 38  | chrysene             | 5.5000               | 0.0000                  | 0.0217                 | 5.4783                 | 0.0000                  | 0.0287                    | 0.0118                 | 0.0080                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 16.0000              | 0.0012                  | 1.4184                 | 14.5804                | 0.0049                  | 1.8758                    | 0.0314                 | 0.0240                  |
| 28  | naphthalene          | 590.0000             | 6.4594                  | 260.8013               | 322.7393               | 26.5321                 | 344.9171                  | 0.6956                 | 0.8370                  |
| 34  | phenanthrene         | 23.0000              | 0.0087                  | 3.7907                 | 19.2006                | 0.0358                  | 5.0132                    | 0.0414                 | 0.0358                  |
| 37  | pyrene               | 10.0000              | 0.0005                  | 0.5867                 | 9.4128                 | 0.0020                  | 0.7760                    | 0.0203                 | 0.0155                  |
| 43  | toluene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.460000,  
 NAPL Volume Frac.(l/l), 0.000877,  
 Soil Volume Frac.(l/l), 0.539123,  
 Porosity (Volume Frac.), 0.460877,

Bulk Density (kg/l), 1.8895,  
 NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

0.1902,

Numerical Accuracy Information

The solution converged in 26 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:25:35 PM

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Sample Name Identification, BH-SED-06-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5900,

NAPLANAL ANALYSIS RESULTS:  
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| ID# | Name                 | Total mass, (mg/kg)* | Mass, in water, (mg/kg)* | Mass in, in soil, (mg/kg)* | Mass, in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|--------------------------|----------------------------|-------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 0.5600               | 0.0033                   | 0.5144                     | 0.0423                  | 0.0093                  | 0.7940                    | 0.0002                 | 0.0004                  |
| 30  | 2-methylnaphthalene  | 1.0000               | 0.0059                   | 0.9174                     | 0.0767                  | 0.0167                  | 1.4162                    | 0.0004                 | 0.0007                  |
| 32  | acenaphthene         | 3.0000               | 0.0157                   | 1.3295                     | 1.6547                  | 0.0446                  | 2.0524                    | 0.0087                 | 0.0130                  |
| 31  | acenaphthylene       | 6.4000               | 0.0466                   | 2.1414                     | 4.2120                  | 0.1322                  | 3.3058                    | 0.0220                 | 0.0336                  |
| 35  | anthracene           | 13.0000              | 0.0014                   | 0.3521                     | 12.6466                 | 0.0039                  | 0.5435                    | 0.0661                 | 0.0863                  |
| 42  | benzene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 24.0000              | 0.0002                   | 5.1226                     | 18.8772                 | 0.0006                  | 7.9078                    | 0.0987                 | 0.1005                  |
| 40  | benzo(a)pyrene       | 26.0000              | 0.0000                   | 4.4413                     | 21.5586                 | 0.0001                  | 6.8562                    | 0.1127                 | 0.1039                  |
| 50  | benzo(b)fluoranthene | 12.0000              | 0.0002                   | 2.3252                     | 9.6745                  | 0.0007                  | 3.5895                    | 0.0506                 | 0.0466                  |
| 51  | benzo(k)fluoranthene | 12.0000              | 0.0001                   | 0.8249                     | 11.1750                 | 0.0002                  | 1.2735                    | 0.0584                 | 0.0538                  |
| 33  | 9h-fluorene          | 1.9000               | 0.0052                   | 0.6990                     | 1.1958                  | 0.0148                  | 1.0791                    | 0.0063                 | 0.0087                  |
| 38  | chrysene             | 23.0000              | 0.0001                   | 0.2821                     | 22.7178                 | 0.0002                  | 0.4355                    | 0.1188                 | 0.1210                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 44.0000              | 0.0147                   | 10.2754                    | 33.7099                 | 0.0417                  | 15.8623                   | 0.1762                 | 0.2026                  |
| 28  | naphthalene          | 16.0000              | 0.4655                   | 11.1364                    | 4.3981                  | 1.3224                  | 17.1914                   | 0.0230                 | 0.0417                  |
| 34  | phenanthrene         | 11.0000              | 0.0163                   | 4.1979                     | 6.7859                  | 0.0463                  | 6.4803                    | 0.0355                 | 0.0463                  |
| 37  | pyrene               | 28.0000              | 0.0065                   | 4.5742                     | 23.4192                 | 0.0186                  | 7.0613                    | 0.1224                 | 0.1408                  |
| 43  | toluene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.590000,  
NAPL Volume Frac.(l/l), 0.000321,  
Soil Volume Frac.(l/l), 0.409679,  
Porosity (Volume Frac.), 0.590321,

Bulk Density (kg/l), 1.6759,  
NAPL Density (kg/l), 0.9000,



NAPL Saturation (%),

0.0543,

Numerical Accuracy Information

The solution converged in 3 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/28/2009 8:34:09 AM

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Sample Name Identification, BH-SED-06-6

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5800,

NAPLANAL ANALYSIS RESULTS:  
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL,<br>, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|---------------------------------|
| ==,  | ====,                 | ====,                        | ====,                           | ====,                             | ====,                          | ====,                          | ====,                            | ====,                         | ====,                           |
| 29,  | 1-methylnaphthalene,  | 17.0000,                     | 0.0622,                         | 10.0961,                          | 6.8417,                        | 0.1812,                        | 15.3999,                         | 0.0056,                       | 0.0070,                         |
| 30,  | 2-methylnaphthalene,  | 30.0000,                     | 0.1091,                         | 17.7044,                          | 12.1865,                       | 0.3177,                        | 27.0051,                         | 0.0101,                       | 0.0125,                         |
| 32,  | acenaphthene,         | 44.0000,                     | 0.0445,                         | 3.9092,                           | 40.0463,                       | 0.1296,                        | 5.9628,                          | 0.0330,                       | 0.0379,                         |
| 31,  | acenaphthylene,       | 7.3000,                      | 0.0089,                         | 0.4242,                           | 6.8670,                        | 0.0259,                        | 0.6470,                          | 0.0057,                       | 0.0066,                         |
| 35,  | anthracene,           | 47.0000,                     | 0.0006,                         | 0.1584,                           | 46.8410,                       | 0.0017,                        | 0.2416,                          | 0.0386,                       | 0.0384,                         |
| 42,  | benzene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 39,  | benz(a)anthracene,    | 48.0000,                     | 0.0001,                         | 1.5320,                           | 46.4680,                       | 0.0002,                        | 2.3368,                          | 0.0383,                       | 0.0297,                         |
| 40,  | benzo(a)pyrene,       | 37.0000,                     | 0.0000,                         | 0.9035,                           | 36.0965,                       | 0.0000,                        | 1.3781,                          | 0.0298,                       | 0.0209,                         |
| 50,  | benzo(b)fluoranthene, | 55.0000,                     | 0.0001,                         | 1.5605,                           | 53.4394,                       | 0.0004,                        | 2.3802,                          | 0.0441,                       | 0.0309,                         |
| 3,   | bromodichloromethane, | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 33,  | 9h-fluorene,          | 36.0000,                     | 0.0171,                         | 2.3860,                           | 33.5969,                       | 0.0499,                        | 3.6395,                          | 0.0277,                       | 0.0295,                         |
| 38,  | chrysene,             | 42.0000,                     | 0.0000,                         | 0.0633,                           | 41.9367,                       | 0.0000,                        | 0.0965,                          | 0.0346,                       | 0.0268,                         |
| 44,  | ethylbenzene,         | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |
| 36,  | fluoranthene,         | 110.0000,                    | 0.0054,                         | 3.9280,                           | 106.0666,                      | 0.0158,                        | 5.9915,                          | 0.0875,                       | 0.0765,                         |
| 28,  | naphthalene,          | 620.0000,                    | 5.8204,                         | 144.4912,                         | 469.6884,                      | 16.9536,                       | 220.3969,                        | 0.3875,                       | 0.5348,                         |
| 34,  | phenanthrene,         | 120.0000,                    | 0.0314,                         | 8.3862,                           | 111.5824,                      | 0.0914,                        | 12.7918,                         | 0.0921,                       | 0.0914,                         |
| 37,  | pyrene,               | 81.0000,                     | 0.0026,                         | 1.8775,                           | 79.1199,                       | 0.0075,                        | 2.8638,                          | 0.0653,                       | 0.0571,                         |
| 43,  | toluene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                         |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.580000,  
NAPL Volume Frac.(l/l), 0.002048,  
Soil Volume Frac.(l/l), 0.417952,  
Porosity (Volume Frac.), 0.582048,

Bulk Density (kg/l), 1.6894,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.3518,

Numerical Accuracy Information

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 The solution converged in 4 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
 29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
 30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
 32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
 31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
 35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
 42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
 39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 1380000.0000
 40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 5500000.0000
 50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 550000.0000
 3 75-27-4 bromodichloromethane 163.8000 1.9700 2.0600 6716.9000 61.0000
 33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
 38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 200000.0000
 44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
 36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
 28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
 34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
 37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
 43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:32:08 PM

Sample Name Identification, BH-SED-07-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5100,

NAPLANAL ANALYSIS RESULTS:

| ID# | Name                 | Total mass, (mg/kg)* | Mass, in water, (mg/kg)* | Mass in, in soil, (mg/kg)* | Mass, in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|--------------------------|----------------------------|-------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 1.2000               | 0.0045                   | 0.9635                     | 0.2321                  | 0.0158                  | 1.3435                    | 0.0004                 | 0.0006                  |
| 30  | 2-methylnaphthalene  | 1.8000               | 0.0067                   | 1.4408                     | 0.3525                  | 0.0236                  | 2.0091                    | 0.0006                 | 0.0009                  |
| 32  | acenaphthene         | 8.4000               | 0.0155                   | 1.8065                     | 6.5780                  | 0.0548                  | 2.5190                    | 0.0107                 | 0.0160                  |
| 31  | acenaphthylene       | 16.0000              | 0.0372                   | 2.3633                     | 13.5995                 | 0.1318                  | 3.2954                    | 0.0221                 | 0.0335                  |
| 35  | anthracene           | 40.0000              | 0.0011                   | 0.3770                     | 39.6219                 | 0.0038                  | 0.5257                    | 0.0643                 | 0.0834                  |
| 42  | benzene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 61.0000              | 0.0001                   | 5.1776                     | 55.8223                 | 0.0005                  | 7.2198                    | 0.0905                 | 0.0918                  |
| 40  | benzo(a)pyrene       | 56.0000              | 0.0000                   | 3.6838                     | 52.3161                 | 0.0001                  | 5.1368                    | 0.0849                 | 0.0778                  |
| 50  | benzo(b)fluoranthene | 59.0000              | 0.0003                   | 4.4789                     | 54.5208                 | 0.0011                  | 6.2455                    | 0.0884                 | 0.0811                  |
| 51  | benzo(k)fluoranthene | 31.0000              | 0.0001                   | 0.7629                     | 30.2370                 | 0.0002                  | 1.0639                    | 0.0490                 | 0.0450                  |
| 33  | 9h-fluorene          | 4.7000               | 0.0042                   | 0.7820                     | 3.9138                  | 0.0149                  | 1.0904                    | 0.0063                 | 0.0088                  |
| 38  | chrysene             | 63.0000              | 0.0001                   | 0.2663                     | 62.7336                 | 0.0002                  | 0.3713                    | 0.1017                 | 0.1031                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 140.0000             | 0.0137                   | 13.2086                    | 126.7778                | 0.0485                  | 18.4184                   | 0.2056                 | 0.2353                  |
| 28  | naphthalene          | 28.0000              | 0.3879                   | 12.8104                    | 14.8017                 | 1.3741                  | 17.8632                   | 0.0240                 | 0.0433                  |
| 34  | phenanthrene         | 28.0000              | 0.0137                   | 4.8847                     | 23.1016                 | 0.0487                  | 6.8113                    | 0.0375                 | 0.0487                  |
| 37  | pyrene               | 75.0000              | 0.0049                   | 4.6934                     | 70.3018                 | 0.0172                  | 6.5446                    | 0.1140                 | 0.1305                  |
| 43  | toluene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.510000,  
NAPL Volume Frac.(l/l), 0.001114,  
Soil Volume Frac.(l/l), 0.488886,  
Porosity (Volume Frac.), 0.511114,

Bulk Density (kg/l), 1.8066,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

0.2179,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with  
CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 138000.0000   |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 550000.0000   |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 55000.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                    | 0.9000            | 0.0394                          | 0.0043                     | 55000.0000    |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 20000.0000    |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time:

8/3/2009 12:34:48 PM

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Sample Name Identification, BH-SED-07-6  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5800,

NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.1900, 0.0012, 0.1888, 0.0000, 0.0034, 0.2873, 0.0002,  
30, 2-methylnaphthalene, 0.3700, 0.0023, 0.3677, 0.0000, 0.0066, 0.5594, 0.0004,  
32, acenaphthene, 0.2100, 0.0024, 0.2076, 0.0000, 0.0069, 0.3159, 0.0033,  
31, acenaphthylene, 0.1600, 0.0033, 0.1567, 0.0000, 0.0095, 0.2384, 0.0040,  
35, anthracene, 0.3200, 0.0012, 0.3188, 0.0000, 0.0035, 0.4850, 0.1266,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 0.9500, 0.0000, 0.9500, 0.0000, 0.0001, 1.4450, 0.0302,  
40, benzo(a)pyrene, 1.0000, 0.0000, 1.0000, 0.0000, 0.0000, 1.5211, 0.0379,  
50, benzo(b)fluoranthene, 1.1000, 0.0001, 1.0999, 0.0000, 0.0003, 1.6731, 0.0357,  
51, benzo(k)fluoranthene, 0.4500, 0.0000, 0.4500, 0.0000, 0.0001, 0.6844, 0.0476,  
33, 9h-fluorene, 0.1800, 0.0013, 0.1787, 0.0000, 0.0037, 0.2719, 0.0036,  
38, chrysene, 0.7600, 0.0002, 0.7598, 0.0000, 0.0006, 1.1557, 0.5280,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 1.7000, 0.0023, 1.6977, 0.0000, 0.0068, 2.5824, 0.0543,  
28, naphthalene, 9.8000, 0.3777, 9.4223, 0.0000, 1.1025, 14.3324, 0.0572,  
34, phenanthrene, 0.6300, 0.0023, 0.6277, 0.0000, 0.0068, 0.9547, 0.0112,  
37, pyrene, 1.2000, 0.0016, 1.1984, 0.0000, 0.0048, 1.8228, 0.0598,  
43, toluene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.420000,  
Water Volume Frac.(l/l), 0.580000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.580000,

Bulk Density (kg/l), 1.6930,

Dilution Factor (Vol. fac.), 0.6446,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:38:16 PM

Sample Name Identification, BH-SED-08-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.6400,

NAPLANAL ANALYSIS RESULTS:

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| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 0.3000               | 0.0023                  | 0.2916                 | 0.0061                 | 0.0057                  | 0.4872                    | 0.0001                 | 0.0002                  |
| 30  | 2-methylnaphthalene  | 0.6700               | 0.0051                  | 0.6510                 | 0.0139                 | 0.0128                  | 1.0878                    | 0.0003                 | 0.0005                  |
| 32  | acenaphthene         | 0.7600               | 0.0083                  | 0.5707                 | 0.1810                 | 0.0207                  | 0.9536                    | 0.0039                 | 0.0061                  |
| 31  | acenaphthylene       | 2.0000               | 0.0351                  | 1.3088                 | 0.6561                 | 0.0875                  | 2.1870                    | 0.0143                 | 0.0223                  |
| 35  | anthracene           | 3.7000               | 0.0017                  | 0.3642                 | 3.3341                 | 0.0043                  | 0.6085                    | 0.0726                 | 0.0966                  |
| 42  | benzene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 8.1000               | 0.0002                  | 4.1768                 | 3.9230                 | 0.0005                  | 6.9797                    | 0.0854                 | 0.0887                  |
| 40  | benzo(a)pyrene       | 8.8000               | 0.0000                  | 3.9335                 | 4.8664                 | 0.0001                  | 6.5731                    | 0.1059                 | 0.0996                  |
| 50  | benzo(b)fluoranthene | 8.8000               | 0.0005                  | 4.2707                 | 4.5288                 | 0.0013                  | 7.1366                    | 0.0986                 | 0.0927                  |
| 38  | chrysene             | 7.6000               | 0.0001                  | 0.3531                 | 7.2468                 | 0.0003                  | 0.5901                    | 0.1578                 | 0.1639                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 14.0000              | 0.0135                  | 7.6173                 | 6.3692                 | 0.0335                  | 12.7290                   | 0.1386                 | 0.1626                  |
| 33  | 9h-fluorene          | 0.6100               | 0.0039                  | 0.4221                 | 0.1840                 | 0.0097                  | 0.7053                    | 0.0040                 | 0.0057                  |
| 28  | naphthalene          | 12.0000              | 0.5375                  | 10.4142                | 1.0483                 | 1.3387                  | 17.4028                   | 0.0228                 | 0.0422                  |
| 34  | phenanthrene         | 3.6000               | 0.0122                  | 2.5409                 | 1.0469                 | 0.0303                  | 4.2461                    | 0.0228                 | 0.0303                  |
| 37  | pyrene               | 9.1000               | 0.0070                  | 3.9451                 | 5.1479                 | 0.0173                  | 6.5925                    | 0.1121                 | 0.1314                  |
| 43  | toluene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 51  | benzo(k)fluoranthene | 3.6000               | 0.0001                  | 0.8085                 | 2.7914                 | 0.0002                  | 1.3510                    | 0.0608                 | 0.0571                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.640000,  
NAPL Volume Frac.(l/l), 0.000073,  
Soil Volume Frac.(l/l), 0.359927,  
Porosity (Volume Frac.), 0.640073,

Bulk Density (kg/l), 1.5939,  
NAPL Density (kg/l), 0.9000,



NAPL Saturation (%),

0.0114,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenapthene          | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenapthylene        | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 550000.0000 |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 200000.0000 |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 550000.0000 |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:42:28 PM

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Sample Name Identification, BH-SED-08-10  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5500,

NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.0880, 0.0005, 0.0875, 0.0000, 0.0015, 0.1279, 0.0001,  
30, 2-methylnaphthalene, 0.2400, 0.0013, 0.2387, 0.0000, 0.0041, 0.3488, 0.0002,  
32, acenaphthene, 0.0530, 0.0005, 0.0525, 0.0000, 0.0017, 0.0767, 0.0006,  
31, acenaphthylene, 0.2000, 0.0036, 0.1964, 0.0000, 0.0115, 0.2869, 0.0036,  
35, anthracene, 0.2500, 0.0008, 0.2492, 0.0000, 0.0026, 0.3641, 0.0719,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 1.3000, 0.0000, 1.3000, 0.0000, 0.0001, 1.8995, 0.0300,  
40, benzo(a)pyrene, 1.4000, 0.0000, 1.4000, 0.0000, 0.0000, 2.0457, 0.0386,  
50, benzo(b)fluoranthene, 1.6000, 0.0001, 1.5999, 0.0000, 0.0004, 2.3377, 0.0378,  
38, chrysene, 1.2000, 0.0003, 1.1997, 0.0000, 0.0009, 1.7531, 0.6059,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 1.8000, 0.0022, 1.7978, 0.0000, 0.0069, 2.6270, 0.0418,  
33, 9h-fluorene, 0.1500, 0.0009, 0.1491, 0.0000, 0.0030, 0.2178, 0.0022,  
28, naphthalene, 17.0000, 0.5825, 16.4175, 0.0000, 1.8454, 23.9896, 0.0724,  
34, phenanthrene, 0.3100, 0.0010, 0.3090, 0.0000, 0.0032, 0.4515, 0.0040,  
37, pyrene, 1.3000, 0.0016, 1.2984, 0.0000, 0.0050, 1.8973, 0.0471,  
43, toluene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
51, benzo(k)fluoranthene, 0.5700, 0.0000, 0.5700, 0.0000, 0.0002, 0.8328, 0.0438,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.450000,  
Water Volume Frac.(l/l), 0.550000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.550000,

Bulk Density (kg/l), 1.7425,

Dilution Factor (Vol. fac.), 0.2443,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
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END OF RECORDS

Sample Name Identification, BH-SED-09-00

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.4900,

NAPLANAL ANALYSIS RESULTS:

Table with 10 columns: ID#, Name, Total mass, Mass in water, Mass in soil, Mass in NAPL, Conc. in water, Sorbed in soil, Conc. in NAPL, Mole fraction in NAPL. Lists various chemical compounds and their concentrations.

(mg/kg)\* --- mg per kg of soil sample (wet soil)
(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.490000,
NAPL Volume Frac.(l/l), 0.000201,
Soil Volume Frac.(l/l), 0.509799,
Porosity (Volume Frac.), 0.490201,

Bulk Density (kg/l), 1.8411,
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0410,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/28/2009 10:14:22 AM

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Sample Name Identification, BH-SED-10-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5900,

NAPLANAL ANALYSIS RESULTS:  
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| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc. in water, (mg/L) | Sorbed in soil, (mg/kg)^ | Conc. in NAPL, (kg/L) | Mole fraction in NAPL |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|------------------------|--------------------------|-----------------------|-----------------------|
| 29  | 1-methylnaphthalene  | 0.5900               | 0.0036                  | 0.5623                 | 0.0241                 | 0.0102                 | 0.8679                   | 0.0002                | 0.0004                |
| 30  | 2-methylnaphthalene  | 0.6800               | 0.0041                  | 0.6477                 | 0.0282                 | 0.0118                 | 0.9997                   | 0.0003                | 0.0005                |
| 32  | acenaphthene         | 1.8000               | 0.0128                  | 1.0852                 | 0.7020                 | 0.0364                 | 1.6749                   | 0.0070                | 0.0106                |
| 31  | acenaphthylene       | 3.0000               | 0.0319                  | 1.4677                 | 1.5004                 | 0.0906                 | 2.2652                   | 0.0150                | 0.0231                |
| 35  | anthracene           | 4.9000               | 0.0010                  | 0.2490                 | 4.6500                 | 0.0027                 | 0.3844                   | 0.0464                | 0.0610                |
| 42  | benzene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                 | 0.0000                   | 0.0000                | 0.0000                |
| 39  | benz(a)anthracene    | 15.0000              | 0.0002                  | 5.1450                 | 9.8547                 | 0.0006                 | 7.9408                   | 0.0983                | 0.1010                |
| 40  | benzo(a)pyrene       | 15.0000              | 0.0000                  | 4.2578                 | 10.7422                | 0.0001                 | 6.5714                   | 0.1072                | 0.0996                |
| 50  | benzo(b)fluoranthene | 13.0000              | 0.0004                  | 4.1105                 | 8.8891                 | 0.0012                 | 6.3441                   | 0.0887                | 0.0824                |
| 38  | chrysene             | 14.0000              | 0.0001                  | 0.3267                 | 13.6732                | 0.0003                 | 0.5042                   | 0.1364                | 0.1401                |
| 44  | ethylbenzene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                 | 0.0000                   | 0.0000                | 0.0000                |
| 36  | fluoranthene         | 31.0000              | 0.0164                  | 11.4536                | 19.5301                | 0.0465                 | 17.6773                  | 0.1949                | 0.2258                |
| 33  | 9h-fluorene          | 1.7000               | 0.0067                  | 0.8963                 | 0.7970                 | 0.0190                 | 1.3834                   | 0.0080                | 0.0112                |
| 28  | naphthalene          | 9.9000               | 0.3317                  | 7.9387                 | 1.6296                 | 0.9425                 | 12.2525                  | 0.0163                | 0.0297                |
| 34  | phenanthrene         | 6.4000               | 0.0135                  | 3.4706                 | 2.9160                 | 0.0383                 | 5.3564                   | 0.0291                | 0.0383                |
| 37  | pyrene               | 21.0000              | 0.0082                  | 5.7338                 | 15.2580                | 0.0233                 | 8.8495                   | 0.1523                | 0.1764                |
| 43  | toluene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                 | 0.0000                   | 0.0000                | 0.0000                |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.590000,  
NAPL Volume Frac.(l/l), 0.000168,  
Soil Volume Frac.(l/l), 0.409832,  
Porosity (Volume Frac.), 0.590168,

Bulk Density (kg/l), 1.6762,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0285,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 1380000.0000  |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 5500000.0000  |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 550000.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 200000.0000   |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 12:51:20 PM

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Sample Name Identification, BH-SED-10-2

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5900,

NAPLANAL ANALYSIS RESULTS:  
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| ID# | Name                 | Total mass, (mg/kg)* | Mass, in water, (mg/kg)* | Mass in, in soil, (mg/kg)* | Mass, in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|--------------------------|----------------------------|-------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 0.8000               | 0.0050                   | 0.7761                     | 0.0189                  | 0.0141                  | 1.1977                    | 0.0003                 | 0.0005                  |
| 30  | 2-methylnaphthalene  | 1.3000               | 0.0081                   | 1.2607                     | 0.0313                  | 0.0229                  | 1.9455                    | 0.0006                 | 0.0009                  |
| 32  | acenaphthene         | 1.4000               | 0.0120                   | 1.0140                     | 0.3741                  | 0.0340                  | 1.5648                    | 0.0067                 | 0.0099                  |
| 31  | acenaphthylene       | 1.8000               | 0.0244                   | 1.1217                     | 0.6540                  | 0.0692                  | 1.7310                    | 0.0118                 | 0.0176                  |
| 35  | anthracene           | 4.1000               | 0.0014                   | 0.3519                     | 3.7468                  | 0.0039                  | 0.5430                    | 0.0676                 | 0.0862                  |
| 42  | benzene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 8.0000               | 0.0002                   | 3.8234                     | 4.1764                  | 0.0004                  | 5.9004                    | 0.0753                 | 0.0750                  |
| 40  | benzo(a)pyrene       | 7.7000               | 0.0000                   | 3.1572                     | 4.5427                  | 0.0001                  | 4.8724                    | 0.0819                 | 0.0738                  |
| 50  | benzo(b)fluoranthene | 7.9000               | 0.0003                   | 3.5372                     | 4.3624                  | 0.0010                  | 5.4588                    | 0.0787                 | 0.0709                  |
| 38  | chrysene             | 7.2000               | 0.0001                   | 0.2895                     | 6.9104                  | 0.0002                  | 0.4468                    | 0.1247                 | 0.1241                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 17.0000              | 0.0123                   | 8.6126                     | 8.3751                  | 0.0350                  | 13.2911                   | 0.1511                 | 0.1698                  |
| 33  | 9h-fluorene          | 1.6000               | 0.0079                   | 1.0564                     | 0.5357                  | 0.0223                  | 1.6303                    | 0.0097                 | 0.0132                  |
| 28  | naphthalene          | 32.0000              | 1.1537                   | 27.6137                    | 3.2325                  | 3.2780                  | 42.6143                   | 0.0583                 | 0.1034                  |
| 34  | phenanthrene         | 7.0000               | 0.0183                   | 4.7200                     | 2.2616                  | 0.0520                  | 7.2841                    | 0.0408                 | 0.0520                  |
| 37  | pyrene               | 12.0000              | 0.0068                   | 4.7638                     | 7.2294                  | 0.0193                  | 7.3516                    | 0.1304                 | 0.1466                  |
| 43  | toluene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 51  | benzo(k)fluoranthene | 4.3000               | 0.0001                   | 0.8574                     | 3.4426                  | 0.0002                  | 1.3231                    | 0.0621                 | 0.0559                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.590000,  
NAPL Volume Frac.(l/l), 0.000093,  
Soil Volume Frac.(l/l), 0.409907,  
Porosity (Volume Frac.), 0.590093,

Bulk Density (kg/l), 1.6763,  
NAPL Density (kg/l), 0.9000,



NAPL Saturation (%),

0.0157,

Numerical Accuracy Information

The solution converged in 6 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 550000.0000 |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 200000.0000 |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 550000.0000 |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:03:50 PM

Sample Name Identification, BH-SED-11-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5800,

NAPLANAL ANALYSIS RESULTS:

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| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 0.8500               | 0.0050                  | 0.8103                 | 0.0348                 | 0.0145                  | 1.2328                    | 0.0003                 | 0.0006                  |
| 30  | 2-methylnaphthalene  | 1.7000               | 0.0099                  | 1.6195                 | 0.0706                 | 0.0290                  | 2.4640                    | 0.0007                 | 0.0011                  |
| 32  | acenaphthene         | 1.3000               | 0.0089                  | 0.7832                 | 0.5080                 | 0.0259                  | 1.1916                    | 0.0051                 | 0.0076                  |
| 31  | acenaphthylene       | 1.8000               | 0.0183                  | 0.8798                 | 0.9018                 | 0.0535                  | 1.3386                    | 0.0091                 | 0.0136                  |
| 35  | anthracene           | 4.4000               | 0.0008                  | 0.2231                 | 4.1761                 | 0.0024                  | 0.3394                    | 0.0420                 | 0.0539                  |
| 42  | benzene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 14.0000              | 0.0002                  | 4.7938                 | 9.2060                 | 0.0005                  | 7.2935                    | 0.0926                 | 0.0927                  |
| 40  | benzo(a)pyrene       | 12.0000              | 0.0000                  | 3.3998                 | 8.6002                 | 0.0001                  | 5.1727                    | 0.0865                 | 0.0784                  |
| 50  | benzo(b)fluoranthene | 12.0000              | 0.0004                  | 3.7875                 | 8.2121                 | 0.0010                  | 5.7625                    | 0.0826                 | 0.0748                  |
| 38  | chrysene             | 13.0000              | 0.0001                  | 0.3026                 | 12.6973                | 0.0002                  | 0.4604                    | 0.1277                 | 0.1279                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 28.0000              | 0.0142                  | 10.3283                | 17.6576                | 0.0414                  | 15.7140                   | 0.1776                 | 0.2007                  |
| 33  | 9h-fluorene          | 1.8000               | 0.0068                  | 0.9481                 | 0.8452                 | 0.0198                  | 1.4424                    | 0.0085                 | 0.0117                  |
| 28  | naphthalene          | 37.0000              | 1.1909                  | 29.6972                | 6.1119                 | 3.4756                  | 45.1831                   | 0.0615                 | 0.1096                  |
| 34  | phenanthrene         | 5.7000               | 0.0115                  | 3.0875                 | 2.6010                 | 0.0336                  | 4.6976                    | 0.0262                 | 0.0336                  |
| 37  | pyrene               | 19.0000              | 0.0071                  | 5.1779                 | 13.8150                | 0.0207                  | 7.8780                    | 0.1390                 | 0.1571                  |
| 43  | toluene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 51  | benzo(k)fluoranthene | 4.6000               | 0.0001                  | 0.5708                 | 4.0292                 | 0.0002                  | 0.8684                    | 0.0405                 | 0.0367                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.580000,  
NAPL Volume Frac.(l/l), 0.000168,  
Soil Volume Frac.(l/l), 0.419832,  
Porosity (Volume Frac.), 0.580168,

Bulk Density (kg/l), 1.6927,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

0.0290,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphtene          | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphtylene        | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:06:15 PM

Sample Name Identification, BH-SED-11-2

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.5400,

NAPLANAL ANALYSIS RESULTS:

Table with 10 columns: ID#, Name, Total mass, Mass in water, Mass in soil, Mass in NAPL, Conc. in water, Sorbed, Conc. in NAPL, Mole fraction in NAPL. Rows list various compounds like 1-methylnaphthalene, 2-methylnaphthalene, acenaphthene, etc.

(mg/kg)\* --- mg per kg of soil sample (wet soil)
(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.540000,
NAPL Volume Frac.(l/l), 0.005250,
Soil Volume Frac.(l/l), 0.454750,
Porosity (Volume Frac.), 0.545250,

Bulk Density (kg/l), 1.7498,
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

0.9628,

Numerical Accuracy Information

The solution converged in 263 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)
Data for chlorinated compounds are from database of Pankow and Johnson with
CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by
Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons
are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid
densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more
accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are
estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties
for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

Table with 8 columns: ID#, CAS#, Name, Molecular Weight (g/mol), Density (kg/l), Henry's Constant (atm-L/mol), Water Solubility (mg/L), Koc (ml/g). Rows list various aromatic hydrocarbons like 1-methylnaphthalene, 2-methylnaphthalene, acenaphthene, etc.

END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:15:49 PM

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Sample Name Identification, BH-SED-12-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5000,

NAPLANAL ANALYSIS RESULTS:

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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|----------------------------|
| ===, | =====,                | =====,                       | =====,                          | =====,                            | =====,                         | =====,                         | =====,                           | =====,                        | =====,                     |
| 29,  | 1-methylnaphthalene,  | 0.1800,                      | 0.0008,                         | 0.1784,                           | 0.0009,                        | 0.0029,                        | 0.2457,                          | 0.0001,                       | 0.0001,                    |
| 30,  | 2-methylnaphthalene,  | 0.3700,                      | 0.0016,                         | 0.3666,                           | 0.0018,                        | 0.0059,                        | 0.5049,                          | 0.0001,                       | 0.0002,                    |
| 32,  | acenaphthene,         | 0.2400,                      | 0.0018,                         | 0.2221,                           | 0.0161,                        | 0.0066,                        | 0.3059,                          | 0.0012,                       | 0.0019,                    |
| 31,  | acenaphthylene,       | 0.9500,                      | 0.0127,                         | 0.8409,                           | 0.0964,                        | 0.0463,                        | 1.1582,                          | 0.0073,                       | 0.0118,                    |
| 35,  | anthracene,           | 1.2000,                      | 0.0010,                         | 0.3874,                           | 0.8115,                        | 0.0038,                        | 0.5336,                          | 0.0615,                       | 0.0847,                    |
| 42,  | benzene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 39,  | benz(a)anthracene,    | 4.8000,                      | 0.0001,                         | 3.9509,                           | 0.8490,                        | 0.0004,                        | 5.4420,                          | 0.0644,                       | 0.0692,                    |
| 40,  | benzo(a)pyrene,       | 5.5000,                      | 0.0000,                         | 4.2867,                           | 1.2133,                        | 0.0001,                        | 5.9044,                          | 0.0920,                       | 0.0895,                    |
| 50,  | benzo(b)fluoranthene, | 4.9000,                      | 0.0003,                         | 3.9431,                           | 0.9566,                        | 0.0010,                        | 5.4312,                          | 0.0725,                       | 0.0705,                    |
| 38,  | chrysene,             | 4.8000,                      | 0.0002,                         | 0.8428,                           | 3.9570,                        | 0.0006,                        | 1.1609,                          | 0.3000,                       | 0.3225,                    |
| 44,  | ethylbenzene,         | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 36,  | fluoranthene,         | 7.5000,                      | 0.0062,                         | 6.2904,                           | 1.2033,                        | 0.0228,                        | 8.6644,                          | 0.0912,                       | 0.1107,                    |
| 33,  | 9h-fluorene,          | 0.3500,                      | 0.0016,                         | 0.3168,                           | 0.0316,                        | 0.0060,                        | 0.4363,                          | 0.0024,                       | 0.0035,                    |
| 28,  | naphthalene,          | 5.3000,                      | 0.1462,                         | 5.0377,                           | 0.1160,                        | 0.5338,                        | 6.9390,                          | 0.0088,                       | 0.0168,                    |
| 34,  | phenanthrene,         | 1.7000,                      | 0.0042,                         | 1.5497,                           | 0.1461,                        | 0.0152,                        | 2.1346,                          | 0.0111,                       | 0.0152,                    |
| 37,  | pyrene,               | 5.0000,                      | 0.0038,                         | 3.8476,                           | 1.1486,                        | 0.0139,                        | 5.2996,                          | 0.0871,                       | 0.1057,                    |
| 43,  | toluene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 51,  | benzo(k)fluoranthene, | 3.0000,                      | 0.0001,                         | 1.6760,                           | 1.3239,                        | 0.0004,                        | 2.3085,                          | 0.1004,                       | 0.0976,                    |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.500000,  
NAPL Volume Frac.(l/l), 0.000024,  
Soil Volume Frac.(l/l), 0.499976,  
Porosity (Volume Frac.), 0.500024,

Bulk Density (kg/l), 1.8250,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

0.0048,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenapthene          | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenapthylene        | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 550000.0000 |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 200000.0000 |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 550000.0000 |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:19:12 PM

Sample Name Identification, BH-SED-12-4

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.4400,

NAPLANAL ANALYSIS RESULTS:

Table with 10 columns: ID#, Name, Total mass, Mass in water, Mass in soil, Mass in NAPL, Conc. in water, Sorbed in soil, Conc. in NAPL, Mole fraction in NAPL. Rows list various hydrocarbons like 1-methylnaphthalene, 2-methylnaphthalene, acenaphthene, etc.

(mg/kg)\* --- mg per kg of soil sample (wet soil)
(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.440000,
NAPL Volume Frac.(l/l), 0.000004,
Soil Volume Frac.(l/l), 0.559996,
Porosity (Volume Frac.), 0.440004,

Bulk Density (kg/l), 1.9240,
NAPL Density (kg/l), 0.9000,



NAPL Saturation (%), 0.0009,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:23:30 PM

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Sample Name Identification, BH-SED-13A-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.2400,

NAPLANAL ANALYSIS RESULTS:

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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|----------------------------|
| 29,  | 1-methylnaphthalene,  | 0.8300,                      | 0.0012,                         | 0.8264,                           | 0.0025,                        | 0.0109,                        | 0.9248,                          | 0.0003,                       | 0.0004,                    |
| 30,  | 2-methylnaphthalene,  | 1.6000,                      | 0.0022,                         | 1.5929,                           | 0.0049,                        | 0.0210,                        | 1.7827,                          | 0.0005,                       | 0.0008,                    |
| 32,  | acenaphthene,         | 0.9100,                      | 0.0022,                         | 0.8682,                           | 0.0396,                        | 0.0211,                        | 0.9717,                          | 0.0041,                       | 0.0062,                    |
| 31,  | acenaphthylene,       | 1.5000,                      | 0.0066,                         | 1.3931,                           | 0.1003,                        | 0.0624,                        | 1.5591,                          | 0.0105,                       | 0.0159,                    |
| 35,  | anthracene,           | 2.2000,                      | 0.0008,                         | 0.9500,                           | 1.2492,                        | 0.0076,                        | 1.0632,                          | 0.1303,                       | 0.1688,                    |
| 42,  | benzene,              | 0.0790,                      | 0.0099,                         | 0.0689,                           | 0.0002,                        | 0.0929,                        | 0.0771,                          | 0.0000,                       | 0.0001,                    |
| 39,  | benz(a)anthracene,    | 4.0000,                      | 0.0000,                         | 3.5245,                           | 0.4755,                        | 0.0003,                        | 3.9445,                          | 0.0496,                       | 0.0501,                    |
| 40,  | benzo(a)pyrene,       | 3.6000,                      | 0.0000,                         | 3.0568,                           | 0.5432,                        | 0.0001,                        | 3.4211,                          | 0.0567,                       | 0.0518,                    |
| 50,  | benzo(b)fluoranthene, | 3.6000,                      | 0.0001,                         | 3.1241,                           | 0.4758,                        | 0.0006,                        | 3.4964,                          | 0.0496,                       | 0.0454,                    |
| 38,  | chrysene,             | 3.3000,                      | 0.0000,                         | 0.8359,                           | 2.4640,                        | 0.0005,                        | 0.9355,                          | 0.2571,                       | 0.2599,                    |
| 44,  | ethylbenzene,         | 0.0049,                      | 0.0001,                         | 0.0048,                           | 0.0000,                        | 0.0005,                        | 0.0054,                          | 0.0000,                       | 0.0000,                    |
| 36,  | fluoranthene,         | 8.9000,                      | 0.0025,                         | 7.9435,                           | 0.9540,                        | 0.0234,                        | 8.8902,                          | 0.0995,                       | 0.1136,                    |
| 33,  | 9h-fluorene,          | 1.8000,                      | 0.0028,                         | 1.6913,                           | 0.1059,                        | 0.0259,                        | 1.8929,                          | 0.0111,                       | 0.0153,                    |
| 28,  | naphthalene,          | 16.0000,                     | 0.1433,                         | 15.6307,                          | 0.2260,                        | 1.3457,                        | 17.4936,                         | 0.0236,                       | 0.0424,                    |
| 34,  | phenanthrene,         | 7.9000,                      | 0.0063,                         | 7.4526,                           | 0.4410,                        | 0.0596,                        | 8.3408,                          | 0.0460,                       | 0.0596,                    |
| 37,  | pyrene,               | 6.0000,                      | 0.0016,                         | 5.0516,                           | 0.9468,                        | 0.0149,                        | 5.6537,                          | 0.0988,                       | 0.1127,                    |
| 43,  | toluene,              | 0.0570,                      | 0.0022,                         | 0.0547,                           | 0.0001,                        | 0.0204,                        | 0.0612,                          | 0.0000,                       | 0.0000,                    |
| 51,  | benzo(k)fluoranthene, | 1.8000,                      | 0.0000,                         | 1.2033,                           | 0.5967,                        | 0.0002,                        | 1.3467,                          | 0.0623,                       | 0.0569,                    |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.240000,  
NAPL Volume Frac.(l/l), 0.000022,  
Soil Volume Frac.(l/l), 0.759978,  
Porosity (Volume Frac.), 0.240022,

Bulk Density (kg/l), 2.2540,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0090,

Numerical Accuracy Information

The solution converged in 6 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time:

8/3/2009 1:26:23 PM

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Sample Name Identification, BH-SED-13-6  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.1900,  
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NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.1200, 0.0001, 0.1199, 0.0000, 0.0015, 0.1305, 0.0003,  
30, 2-methylnaphthalene, 0.2100, 0.0002, 0.2098, 0.0000, 0.0027, 0.2284, 0.0005,  
32, acenaphthene, 0.1100, 0.0002, 0.1098, 0.0000, 0.0026, 0.1195, 0.0036,  
31, acenaphthylene, 0.2200, 0.0008, 0.2192, 0.0000, 0.0095, 0.2386, 0.0116,  
35, anthracene, 0.2600, 0.0002, 0.2598, 0.0000, 0.0020, 0.2828, 0.2153,  
42, benzene, 0.4900, 0.0472, 0.4428, 0.0000, 0.5807, 0.4820, 0.0016,  
39, benz(a)anthracene, 0.3400, 0.0000, 0.3400, 0.0000, 0.0000, 0.3701, 0.0226,  
40, benzo(a)pyrene, 0.2900, 0.0000, 0.2900, 0.0000, 0.0000, 0.3157, 0.0229,  
50, benzo(b)fluoranthene, 0.3100, 0.0000, 0.3100, 0.0000, 0.0001, 0.3374, 0.0210,  
38, chrysene, 0.3200, 0.0000, 0.3200, 0.0000, 0.0002, 0.3483, 0.4639,  
44, ethylbenzene, 0.0570, 0.0005, 0.0565, 0.0000, 0.0056, 0.0616, 0.0002,  
36, fluoranthene, 0.8700, 0.0002, 0.8698, 0.0000, 0.0025, 0.9468, 0.0580,  
33, 9h-fluorene, 0.3100, 0.0004, 0.3096, 0.0000, 0.0046, 0.3370, 0.0131,  
28, naphthalene, 2.4000, 0.0162, 2.3838, 0.0000, 0.1996, 2.5948, 0.0302,  
34, phenanthrene, 1.1000, 0.0007, 1.0993, 0.0000, 0.0085, 1.1966, 0.0410,  
37, pyrene, 0.6000, 0.0001, 0.5999, 0.0000, 0.0017, 0.6530, 0.0624,  
43, toluene, 0.3000, 0.0086, 0.2914, 0.0000, 0.1057, 0.3172, 0.0009,  
51, benzo(k)fluoranthene, 0.1400, 0.0000, 0.1400, 0.0000, 0.0000, 0.1524, 0.0309,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.810000,  
Water Volume Frac.(l/l), 0.190000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.190000,

Bulk Density (kg/l), 2.3365,

Dilution Factor (Vol. fac.), 3.7946,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
 29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
 30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
 32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
 31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
 35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
 42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
 39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
 40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
 50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
 38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
 44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
 36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
 33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
 28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
 34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
 37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
 43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
 51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time:

8/3/2009 1:28:38 PM

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Sample Name Identification, BH-SED-13B-00  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.7300,  
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NAPLANAL ANALYSIS RESULTS:

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ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.1400, 0.0017, 0.1383, 0.0000, 0.0033, 0.2795, 0.0002,  
30, 2-methylnaphthalene, 0.2800, 0.0033, 0.2767, 0.0000, 0.0066, 0.5590, 0.0004,  
32, acenaphthene, 0.0980, 0.0021, 0.0959, 0.0000, 0.0042, 0.1937, 0.0019,  
31, acenaphthylene, 0.2000, 0.0078, 0.1922, 0.0000, 0.0155, 0.3882, 0.0062,  
35, anthracene, 0.2800, 0.0020, 0.2780, 0.0000, 0.0040, 0.5616, 0.1404,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 0.5800, 0.0000, 0.5800, 0.0000, 0.0001, 1.1717, 0.0235,  
40, benzo(a)pyrene, 0.6300, 0.0000, 0.6300, 0.0000, 0.0000, 1.2727, 0.0304,  
50, benzo(b)fluoranthene, 1.5000, 0.0003, 1.4997, 0.0000, 0.0006, 3.0298, 0.0620,  
38, chrysene, 0.6400, 0.0003, 0.6397, 0.0000, 0.0006, 1.2923, 0.5654,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 1.1000, 0.0029, 1.0971, 0.0000, 0.0058, 2.2163, 0.0446,  
33, 9h-fluorene, 0.1500, 0.0021, 0.1479, 0.0000, 0.0041, 0.2989, 0.0038,  
28, naphthalene, 1.7000, 0.1237, 1.5763, 0.0000, 0.2450, 3.1845, 0.0122,  
34, phenanthrene, 0.5300, 0.0038, 0.5262, 0.0000, 0.0076, 1.0630, 0.0120,  
37, pyrene, 0.7900, 0.0021, 0.7879, 0.0000, 0.0042, 1.5917, 0.0500,  
43, toluene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
51, benzo(k)fluoranthene, 0.3500, 0.0001, 0.3499, 0.0000, 0.0001, 0.7070, 0.0471,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.270000,  
Water Volume Frac.(l/l), 0.730000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.730000,

Bulk Density (kg/l), 1.4455,

Dilution Factor (Vol. fac.), 0.5751,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
 29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
 30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
 32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
 31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
 35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
 42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
 39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
 40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
 50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
 38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
 44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
 36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
 33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
 28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
 34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
 37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
 43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
 51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:32:00 PM

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Sample Name Identification, BH-SED-13C-00  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.7800,

NAPLANAL ANALYSIS RESULTS:

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~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.0700, 0.0011, 0.0689, 0.0000, 0.0019, 0.1611, 0.0002,  
30, 2-methylnaphthalene, 0.1400, 0.0022, 0.1378, 0.0000, 0.0038, 0.3222, 0.0004,  
32, acenaphthene, 0.0590, 0.0017, 0.0573, 0.0000, 0.0029, 0.1340, 0.0022,  
31, acenaphthylene, 0.1100, 0.0056, 0.1044, 0.0000, 0.0098, 0.2441, 0.0063,  
35, anthracene, 0.1600, 0.0015, 0.1585, 0.0000, 0.0026, 0.3705, 0.1493,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 0.3100, 0.0000, 0.3100, 0.0000, 0.0001, 0.7247, 0.0234,  
40, benzo(a)pyrene, 0.3200, 0.0000, 0.3200, 0.0000, 0.0000, 0.7481, 0.0288,  
50, benzo(b)fluoranthene, 1.4000, 0.0003, 1.3997, 0.0000, 0.0006, 3.2723, 0.1079,  
38, chrysene, 0.3100, 0.0002, 0.3098, 0.0000, 0.0004, 0.7243, 0.5108,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 0.5900, 0.0021, 0.5879, 0.0000, 0.0036, 1.3745, 0.0446,  
33, 9h-fluorene, 0.1000, 0.0018, 0.0982, 0.0000, 0.0031, 0.2296, 0.0047,  
28, naphthalene, 0.7700, 0.0719, 0.6981, 0.0000, 0.1256, 1.6322, 0.0101,  
34, phenanthrene, 0.3100, 0.0029, 0.3071, 0.0000, 0.0051, 0.7179, 0.0130,  
37, pyrene, 0.4300, 0.0015, 0.4285, 0.0000, 0.0026, 1.0018, 0.0507,  
43, toluene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
51, benzo(k)fluoranthene, 0.1900, 0.0000, 0.1900, 0.0000, 0.0001, 0.4441, 0.0477,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.220000,  
Water Volume Frac.(l/l), 0.780000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.780000,

Bulk Density (kg/l), 1.3630,

Dilution Factor (Vol. fac.), 1.5390,

No NAPL (i.e., NAPL Saturation Equals Zero)



The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:35:20 PM

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Sample Name Identification, BH-SED-13C-6  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.4200,

NAPLANAL ANALYSIS RESULTS:

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~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.3400, 0.0011, 0.3389, 0.0000, 0.0051, 0.4315, 0.0003,  
30, 2-methylnaphthalene, 0.1800, 0.0006, 0.1794, 0.0000, 0.0027, 0.2285, 0.0002,  
32, acenaphthene, 1.1000, 0.0065, 1.0935, 0.0000, 0.0303, 1.3923, 0.0149,  
31, acenaphthylene, 0.1400, 0.0015, 0.1385, 0.0000, 0.0071, 0.1763, 0.0030,  
35, anthracene, 0.5700, 0.0011, 0.5689, 0.0000, 0.0052, 0.7243, 0.1933,  
42, benzene, 0.0640, 0.0159, 0.0481, 0.0000, 0.0739, 0.0613, 0.0001,  
39, benz(a)anthracene, 1.0000, 0.0000, 1.0000, 0.0000, 0.0001, 1.2732, 0.0272,  
40, benzo(a)pyrene, 0.7200, 0.0000, 0.7200, 0.0000, 0.0000, 0.9167, 0.0233,  
50, benzo(b)fluoranthene, 0.7800, 0.0000, 0.7800, 0.0000, 0.0002, 0.9931, 0.0217,  
38, chrysene, 0.7700, 0.0001, 0.7699, 0.0000, 0.0005, 0.9803, 0.4577,  
44, ethylbenzene, 0.0044, 0.0001, 0.0043, 0.0000, 0.0005, 0.0055, 0.0000,  
36, fluoranthene, 2.7000, 0.0019, 2.6981, 0.0000, 0.0090, 3.4353, 0.0738,  
33, 9h-fluorene, 0.5600, 0.0021, 0.5579, 0.0000, 0.0097, 0.7104, 0.0097,  
28, naphthalene, 8.4000, 0.1729, 8.2271, 0.0000, 0.8058, 10.4752, 0.0427,  
34, phenanthrene, 1.3000, 0.0025, 1.2975, 0.0000, 0.0118, 1.6520, 0.0198,  
37, pyrene, 1.7000, 0.0012, 1.6988, 0.0000, 0.0057, 2.1630, 0.0725,  
43, toluene, 0.0072, 0.0006, 0.0066, 0.0000, 0.0028, 0.0084, 0.0000,  
51, benzo(k)fluoranthene, 0.4400, 0.0000, 0.4400, 0.0000, 0.0001, 0.5602, 0.0398,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.580000,  
Water Volume Frac.(l/l), 0.420000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.420000,

Bulk Density (kg/l), 1.9570,

Dilution Factor (Vol. fac.), 0.6809,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
 29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
 30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
 32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
 31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
 35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
 42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
 39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
 40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
 50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
 38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
 44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
 36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 3800.0000
 33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
 28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
 34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 1400.0000
 37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 3800.0000
 43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
 51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time:

8/3/2009 1:37:09 PM

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Sample Name Identification, BH-SED-14-00  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.7600,

NAPLANAL ANALYSIS RESULTS:

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~~~~~  
ID#, Name, Total, Mass, Mass, Mass, Conc., Sorbed, Mole fraction,  
, , mass, in water, in soil, in NAPL, in water, in soil, in NAPL,  
, , (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/kg)\*, (mg/L), (mg/kg)^, ,  
====, =====, =====, =====, =====, =====, =====, =====, =====,  
29, 1-methylnaphthalene, 0.1700, 0.0024, 0.1676, 0.0000, 0.0043, 0.3680, 0.0002,  
30, 2-methylnaphthalene, 0.3200, 0.0044, 0.3156, 0.0000, 0.0081, 0.6927, 0.0003,  
32, acenaphthene, 0.2400, 0.0061, 0.2339, 0.0000, 0.0112, 0.5135, 0.0035,  
31, acenaphthylene, 0.2000, 0.0091, 0.1909, 0.0000, 0.0168, 0.4190, 0.0046,  
35, anthracene, 0.3900, 0.0033, 0.3867, 0.0000, 0.0061, 0.8488, 0.1455,  
42, benzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
39, benz(a)anthracene, 0.7700, 0.0001, 0.7699, 0.0000, 0.0001, 1.6900, 0.0232,  
40, benzo(a)pyrene, 0.7300, 0.0000, 0.7300, 0.0000, 0.0000, 1.6023, 0.0262,  
50, benzo(b)fluoranthene, 1.7000, 0.0004, 1.6996, 0.0000, 0.0007, 3.7306, 0.0523,  
38, chrysene, 0.8800, 0.0005, 0.8795, 0.0000, 0.0010, 1.9304, 0.5789,  
44, ethylbenzene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
36, fluoranthene, 1.7000, 0.0053, 1.6947, 0.0000, 0.0098, 3.7197, 0.0513,  
33, 9h-fluorene, 0.3200, 0.0052, 0.3148, 0.0000, 0.0095, 0.6911, 0.0060,  
28, naphthalene, 1.5000, 0.1263, 1.3737, 0.0000, 0.2319, 3.0153, 0.0079,  
34, phenanthrene, 0.8600, 0.0073, 0.8527, 0.0000, 0.0134, 1.8717, 0.0144,  
37, pyrene, 1.2000, 0.0038, 1.1962, 0.0000, 0.0069, 2.6257, 0.0565,  
43, toluene, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000, 0.0000,  
51, benzo(k)fluoranthene, 0.2900, 0.0001, 0.2899, 0.0000, 0.0001, 0.6364, 0.0291,

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.240000,  
Water Volume Frac.(l/l), 0.760000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.760000,

Bulk Density (kg/l), 1.3960,

Dilution Factor (Vol. fac.), 0.0796,

No NAPL (i.e., NAPL Saturation Equals Zero)

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

```

=====
ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 38000.0000
33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 14000.0000
37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 38000.0000
43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:38:51 PM

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Sample Name Identification, BH-SED-14-8

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.7000,

NAPLANAL ANALYSIS RESULTS:  
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| ID#, | Name,                 | Total,<br>mass,<br>(mg/kg)*, | Mass,<br>in water,<br>(mg/kg)*, | Mass in,<br>in soil,<br>(mg/kg)*, | Mass,<br>in NAPL,<br>(mg/kg)*, | Conc.,<br>in water,<br>(mg/L), | Sorbed,<br>in soil,<br>(mg/kg)^, | Conc.,<br>in NAPL,<br>(kg/L), | Mole fraction,<br>in NAPL, |
|------|-----------------------|------------------------------|---------------------------------|-----------------------------------|--------------------------------|--------------------------------|----------------------------------|-------------------------------|----------------------------|
| 29,  | 1-methylnaphthalene,  | 0.6700,                      | 0.0069,                         | 0.6617,                           | 0.0015,                        | 0.0146,                        | 1.2443,                          | 0.0003,                       | 0.0006,                    |
| 30,  | 2-methylnaphthalene,  | 1.2000,                      | 0.0123,                         | 1.1850,                           | 0.0027,                        | 0.0262,                        | 2.2284,                          | 0.0006,                       | 0.0010,                    |
| 32,  | acenaphthene,         | 0.7700,                      | 0.0140,                         | 0.7311,                           | 0.0249,                        | 0.0299,                        | 1.3748,                          | 0.0057,                       | 0.0087,                    |
| 31,  | acenaphthylene,       | 0.3800,                      | 0.0123,                         | 0.3489,                           | 0.0188,                        | 0.0262,                        | 0.6561,                          | 0.0043,                       | 0.0067,                    |
| 35,  | anthracene,           | 0.9900,                      | 0.0031,                         | 0.4974,                           | 0.4895,                        | 0.0067,                        | 0.9353,                          | 0.1122,                       | 0.1485,                    |
| 42,  | benzene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 39,  | benz(a)anthracene,    | 2.2000,                      | 0.0001,                         | 1.9981,                           | 0.2017,                        | 0.0003,                        | 3.7575,                          | 0.0463,                       | 0.0478,                    |
| 40,  | benzo(a)pyrene,       | 2.1000,                      | 0.0000,                         | 1.8535,                           | 0.2465,                        | 0.0001,                        | 3.4855,                          | 0.0565,                       | 0.0528,                    |
| 50,  | benzo(b)fluoranthene, | 2.3000,                      | 0.0003,                         | 2.0643,                           | 0.2353,                        | 0.0007,                        | 3.8821,                          | 0.0540,                       | 0.0504,                    |
| 38,  | chrysene,             | 2.1000,                      | 0.0003,                         | 0.6549,                           | 1.4448,                        | 0.0006,                        | 1.2316,                          | 0.3313,                       | 0.3421,                    |
| 44,  | ethylbenzene,         | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 36,  | fluoranthene,         | 5.2000,                      | 0.0110,                         | 4.7610,                           | 0.4279,                        | 0.0236,                        | 8.9533,                          | 0.0981,                       | 0.1144,                    |
| 33,  | 9h-fluorene,          | 0.8100,                      | 0.0092,                         | 0.7649,                           | 0.0359,                        | 0.0197,                        | 1.4385,                          | 0.0082,                       | 0.0117,                    |
| 28,  | naphthalene,          | 6.1000,                      | 0.3831,                         | 5.6557,                           | 0.0612,                        | 0.8181,                        | 10.6358,                         | 0.0140,                       | 0.0258,                    |
| 34,  | phenanthrene,         | 2.7000,                      | 0.0162,                         | 2.5700,                           | 0.1138,                        | 0.0345,                        | 4.8330,                          | 0.0261,                       | 0.0345,                    |
| 37,  | pyrene,               | 3.4000,                      | 0.0069,                         | 2.9757,                           | 0.4174,                        | 0.0147,                        | 5.5959,                          | 0.0957,                       | 0.1116,                    |
| 43,  | toluene,              | 0.0000,                      | 0.0000,                         | 0.0000,                           | 0.0000,                        | 0.0000,                        | 0.0000,                          | 0.0000,                       | 0.0000,                    |
| 51,  | benzo(k)fluoranthene, | 0.7500,                      | 0.0001,                         | 0.5469,                           | 0.2030,                        | 0.0002,                        | 1.0285,                          | 0.0465,                       | 0.0435,                    |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.700000,  
NAPL Volume Frac.(l/l), 0.000007,  
Soil Volume Frac.(l/l), 0.299993,  
Porosity (Volume Frac.), 0.700007,

Bulk Density (kg/l), 1.4950,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0009,

Numerical Accuracy Information

The solution converged in 6 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenapthene          | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenapthylene        | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 550000.0000 |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 200000.0000 |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 550000.0000 |

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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:41:02 PM

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Sample Name Identification, BH-SED-15-2  
Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.4000,  
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NAPLANAL ANALYSIS RESULTS:

| ID# | Name                  | Total,<br>mass,<br>(mg/kg)* | Mass,<br>in water,<br>(mg/kg)* | Mass,<br>in soil,<br>(mg/kg)* | Mass,<br>in NAPL,<br>(mg/kg)* | Conc.,<br>in water,<br>(mg/L) | Sorbed,<br>in soil,<br>(mg/kg)^ | Mole fraction,<br>in NAPL, |
|-----|-----------------------|-----------------------------|--------------------------------|-------------------------------|-------------------------------|-------------------------------|---------------------------------|----------------------------|
| 29, | 1-methylnaphthalene,  | 0.0970,                     | 0.0003,                        | 0.0967,                       | 0.0000,                       | 0.0014,                       | 0.1210,                         | 0.0001,                    |
| 30, | 2-methylnaphthalene,  | 0.2400,                     | 0.0007,                        | 0.2393,                       | 0.0000,                       | 0.0035,                       | 0.2995,                         | 0.0002,                    |
| 32, | acenapthene,          | 0.0990,                     | 0.0005,                        | 0.0985,                       | 0.0000,                       | 0.0027,                       | 0.1232,                         | 0.0010,                    |
| 31, | acenapthylene,        | 0.3400,                     | 0.0034,                        | 0.3366,                       | 0.0000,                       | 0.0169,                       | 0.4213,                         | 0.0055,                    |
| 35, | anthracene,           | 0.4600,                     | 0.0008,                        | 0.4592,                       | 0.0000,                       | 0.0041,                       | 0.5747,                         | 0.1168,                    |
| 42, | benzene,              | 0.0000,                     | 0.0000,                        | 0.0000,                       | 0.0000,                       | 0.0000,                       | 0.0000,                         | 0.0000,                    |
| 39, | benz(a)anthracene,    | 1.1000,                     | 0.0000,                        | 1.1000,                       | 0.0000,                       | 0.0001,                       | 1.3767,                         | 0.0224,                    |
| 40, | benzo(a)pyrene,       | 1.5000,                     | 0.0000,                        | 1.5000,                       | 0.0000,                       | 0.0000,                       | 1.8773,                         | 0.0364,                    |
| 50, | benzo(b)fluoranthene, | 1.5000,                     | 0.0001,                        | 1.4999,                       | 0.0000,                       | 0.0003,                       | 1.8773,                         | 0.0312,                    |
| 38, | chrysene,             | 1.3000,                     | 0.0002,                        | 1.2998,                       | 0.0000,                       | 0.0008,                       | 1.6268,                         | 0.5784,                    |
| 44, | ethylbenzene,         | 0.0000,                     | 0.0000,                        | 0.0000,                       | 0.0000,                       | 0.0000,                       | 0.0000,                         | 0.0000,                    |
| 36, | fluoranthene,         | 3.0000,                     | 0.0020,                        | 2.9980,                       | 0.0000,                       | 0.0099,                       | 3.7522,                         | 0.0614,                    |
| 33, | 9h-fluorene,          | 0.2200,                     | 0.0008,                        | 0.2192,                       | 0.0000,                       | 0.0038,                       | 0.2744,                         | 0.0028,                    |
| 28, | naphthalene,          | 3.2000,                     | 0.0607,                        | 3.1393,                       | 0.0000,                       | 0.3022,                       | 3.9290,                         | 0.0122,                    |
| 34, | phenanthrene,         | 0.6600,                     | 0.0012,                        | 0.6588,                       | 0.0000,                       | 0.0059,                       | 0.8246,                         | 0.0075,                    |
| 37, | pyrene,               | 2.0000,                     | 0.0013,                        | 1.9987,                       | 0.0000,                       | 0.0066,                       | 2.5015,                         | 0.0638,                    |
| 43, | toluene,              | 0.0000,                     | 0.0000,                        | 0.0000,                       | 0.0000,                       | 0.0000,                       | 0.0000,                         | 0.0000,                    |
| 51, | benzo(k)fluoranthene, | 0.8900,                     | 0.0000,                        | 0.8900,                       | 0.0000,                       | 0.0002,                       | 1.1138,                         | 0.0603,                    |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Soil Volume Frac.(l/l), 0.600000,  
Water Volume Frac.(l/l), 0.400000,  
NAPL Volume Frac.(l/l), 0.000000,  
Porosity (Volume Frac.), 0.400000,

Bulk Density (kg/l), 1.9900,

Dilution Factor (Vol. fac.), 0.2800,

No NAPL (i.e., NAPL Saturation Equals Zero)



The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

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ID# CAS# Name Molecular Weight Density Henry's Constant Water Solubility Koc
 (g/mol) (kg/l) (atm-L/mol) (mg/L) (ml/g)
=====
29 90-12-0 1-methylnaphthalene 142.2000 0.9000 0.2600 25.8000 8500.0000
30 91-57-6 2-methylnaphthalene 142.2000 0.9000 0.5180 25.4000 8500.0000
32 83-32-9 acenaphthene 154.2100 0.9000 0.0920 3.4200 4600.0000
31 208-96-8 acenaphthylene 152.2000 0.9000 1.4800 3.9300 2500.0000
35 120-12-7 anthracene 178.2400 0.9000 1.0200 0.0450 14000.0000
42 71-43-2 benzene 78.1100 0.8740 5.5900 1750.0000 83.0000
39 56-55-3 benz(a)anthracene 228.3000 0.9000 0.0012 0.0057 138000.0000
40 50-32-8 benzo(a)pyrene 252.3200 0.9000 0.0016 0.0012 550000.0000
50 205-99-2 benzo(b)fluoranthene 252.3200 0.9000 0.0119 0.0140 55000.0000
38 218-01-9 chrysene 228.3000 0.9000 0.0011 0.0018 20000.0000
44 100-41-4 ethylbenzene 106.1700 0.8670 6.4300 152.0000 1100.0000
36 206-44-0 fluoranthene 202.2600 0.9000 0.0065 0.2060 3800.0000
33 86-73-7 9h-fluorene 166.2200 0.9000 0.0642 1.6900 7300.0000
28 91-20-3 naphthalene 128.1800 0.9000 1.1500 31.7000 1300.0000
34 85-01-8 phenanthrene 178.2400 0.9000 0.1590 1.0000 1400.0000
37 129-00-0 pyrene 202.2600 0.9000 0.0050 0.1320 3800.0000
43 108-88-3 toluene 92.1400 0.8620 6.3700 535.0000 300.0000
51 207-08-9 benzo(k)fluoranthene 252.3200 0.9000 0.0394 0.0043 55000.0000
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END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 7/28/2009 2:10:43 PM

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Sample Name Identification, BH-SED-16-0

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5500,

NAPLANAL ANALYSIS RESULTS:  
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| ID# | Name                 | Total mass, (mg/kg)* | Mass in water, (mg/kg)* | Mass in soil, (mg/kg)* | Mass in NAPL, (mg/kg)* | Conc. in water, (mg/L) | Sorbed in soil, (mg/kg)^ | Conc. in NAPL, (kg/L) | Mole fraction in NAPL |
|-----|----------------------|----------------------|-------------------------|------------------------|------------------------|------------------------|--------------------------|-----------------------|-----------------------|
| 29  | 1-methylnaphthalene  | 0.1400               | 0.0008                  | 0.1389                 | 0.0003                 | 0.0024                 | 0.2030                   | 0.0001                | 0.0001                |
| 30  | 2-methylnaphthalene  | 0.3400               | 0.0018                  | 0.3373                 | 0.0008                 | 0.0058                 | 0.4929                   | 0.0001                | 0.0002                |
| 32  | acenaphthene         | 0.0980               | 0.0009                  | 0.0936                 | 0.0035                 | 0.0030                 | 0.1367                   | 0.0005                | 0.0009                |
| 31  | acenaphthylene       | 0.8600               | 0.0147                  | 0.7981                 | 0.0472                 | 0.0466                 | 1.1662                   | 0.0073                | 0.0119                |
| 35  | anthracene           | 0.8000               | 0.0013                  | 0.3840                 | 0.4148                 | 0.0040                 | 0.5611                   | 0.0643                | 0.0891                |
| 42  | benzene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                 | 0.0000                   | 0.0000                | 0.0000                |
| 39  | benz(a)anthracene    | 3.4000               | 0.0001                  | 3.0607                 | 0.3392                 | 0.0003                 | 4.4725                   | 0.0526                | 0.0569                |
| 40  | benzo(a)pyrene       | 5.5000               | 0.0000                  | 4.7994                 | 0.7005                 | 0.0001                 | 7.0131                   | 0.1087                | 0.1063                |
| 50  | benzo(b)fluoranthene | 8.3000               | 0.0006                  | 7.3765                 | 0.9229                 | 0.0020                 | 10.7788                  | 0.1432                | 0.1400                |
| 38  | chrysene             | 3.1000               | 0.0002                  | 0.9061                 | 2.1937                 | 0.0007                 | 1.3240                   | 0.3404                | 0.3678                |
| 44  | ethylbenzene         | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                 | 0.0000                   | 0.0000                | 0.0000                |
| 36  | fluoranthene         | 4.5000               | 0.0050                  | 4.0914                 | 0.4036                 | 0.0157                 | 5.9785                   | 0.0626                | 0.0764                |
| 33  | 9h-fluorene          | 0.2900               | 0.0017                  | 0.2742                 | 0.0141                 | 0.0055                 | 0.4006                   | 0.0022                | 0.0032                |
| 28  | naphthalene          | 1.9000               | 0.0644                  | 1.8141                 | 0.0215                 | 0.2039                 | 2.6508                   | 0.0033                | 0.0064                |
| 34  | phenanthrene         | 0.9900               | 0.0031                  | 0.9412                 | 0.0457                 | 0.0098                 | 1.3752                   | 0.0071                | 0.0098                |
| 37  | pyrene               | 5.2000               | 0.0055                  | 4.5015                 | 0.6930                 | 0.0173                 | 6.5778                   | 0.1075                | 0.1311                |
| 43  | toluene              | 0.0000               | 0.0000                  | 0.0000                 | 0.0000                 | 0.0000                 | 0.0000                   | 0.0000                | 0.0000                |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.550000,  
NAPL Volume Frac.(l/l), 0.000011,  
Soil Volume Frac.(l/l), 0.449989,  
Porosity (Volume Frac.), 0.550011,

Bulk Density (kg/l), 1.7425,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%), 0.0020,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
 Data for chlorinated compounds are from database of Pankow and Johnson with  
 CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
 Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
 are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
 densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
 accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
 estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
 for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 1380000.0000  |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 5500000.0000  |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 550000.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 200000.0000   |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |

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 END OF RECORDS

NAPLANAL Version 1.0.0

Date and Time: 8/3/2009 1:44:53 PM

Sample Name Identification, BH-SED-17-00

Model used: Liquid saturated & water moisture content known,  
Water moisture content (Volume Frac.), 0.5700,

NAPLANAL ANALYSIS RESULTS:

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| ID# | Name                 | Total mass, (mg/kg)* | Mass, in water, (mg/kg)* | Mass in, in soil, (mg/kg)* | Mass, in NAPL, (mg/kg)* | Conc., in water, (mg/L) | Sorbed, in soil, (mg/kg)^ | Conc., in NAPL, (kg/L) | Mole fraction, in NAPL, |
|-----|----------------------|----------------------|--------------------------|----------------------------|-------------------------|-------------------------|---------------------------|------------------------|-------------------------|
| 29  | 1-methylnaphthalene  | 0.1800               | 0.0010                   | 0.1781                     | 0.0009                  | 0.0031                  | 0.2672                    | 0.0001                 | 0.0001                  |
| 30  | 2-methylnaphthalene  | 0.4100               | 0.0024                   | 0.4056                     | 0.0020                  | 0.0072                  | 0.6085                    | 0.0002                 | 0.0003                  |
| 32  | acenaphthene         | 0.1500               | 0.0015                   | 0.1382                     | 0.0103                  | 0.0045                  | 0.2073                    | 0.0008                 | 0.0013                  |
| 31  | acenaphthylene       | 1.1000               | 0.0193                   | 0.9663                     | 0.1144                  | 0.0580                  | 1.4497                    | 0.0093                 | 0.0148                  |
| 35  | anthracene           | 1.4000               | 0.0016                   | 0.4423                     | 0.9561                  | 0.0047                  | 0.6636                    | 0.0776                 | 0.1053                  |
| 42  | benzene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 39  | benz(a)anthracene    | 4.2000               | 0.0001                   | 3.4376                     | 0.7623                  | 0.0004                  | 5.1573                    | 0.0619                 | 0.0656                  |
| 40  | benzo(a)pyrene       | 5.0000               | 0.0000                   | 3.8697                     | 1.1303                  | 0.0001                  | 5.8056                    | 0.0918                 | 0.0880                  |
| 50  | benzo(b)fluoranthene | 6.0000               | 0.0004                   | 4.7983                     | 1.2013                  | 0.0013                  | 7.1987                    | 0.0975                 | 0.0935                  |
| 38  | chrysene             | 3.6000               | 0.0002                   | 0.6159                     | 2.9840                  | 0.0005                  | 0.9240                    | 0.2422                 | 0.2567                  |
| 44  | ethylbenzene         | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 36  | fluoranthene         | 7.5000               | 0.0082                   | 6.2567                     | 1.2351                  | 0.0247                  | 9.3867                    | 0.1003                 | 0.1199                  |
| 33  | 9h-fluorene          | 0.5100               | 0.0031                   | 0.4595                     | 0.0473                  | 0.0094                  | 0.6894                    | 0.0038                 | 0.0056                  |
| 28  | naphthalene          | 2.4000               | 0.0869                   | 2.2594                     | 0.0537                  | 0.2607                  | 3.3896                    | 0.0044                 | 0.0082                  |
| 34  | phenanthrene         | 2.8000               | 0.0091                   | 2.5435                     | 0.2474                  | 0.0273                  | 3.8159                    | 0.0201                 | 0.0273                  |
| 37  | pyrene               | 6.9000               | 0.0069                   | 5.2697                     | 1.6234                  | 0.0208                  | 7.9059                    | 0.1318                 | 0.1576                  |
| 43  | toluene              | 0.0000               | 0.0000                   | 0.0000                     | 0.0000                  | 0.0000                  | 0.0000                    | 0.0000                 | 0.0000                  |
| 51  | benzo(k)fluoranthene | 1.6000               | 0.0001                   | 0.8814                     | 0.7185                  | 0.0002                  | 1.3224                    | 0.0583                 | 0.0559                  |

(mg/kg)\* --- mg per kg of soil sample (wet soil)

(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.570000,  
NAPL Volume Frac.(l/l), 0.000021,  
Soil Volume Frac.(l/l), 0.429979,  
Porosity (Volume Frac.), 0.570021,

Bulk Density (kg/l), 1.7095,  
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

0.0037,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight (g/mol) | Density (kg/l) | Henry's Constant (atm-L/mol) | Water Solubility (mg/L) | Koc (ml/g)  |
|-----|----------|----------------------|--------------------------|----------------|------------------------------|-------------------------|-------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                 | 0.9000         | 0.2600                       | 25.8000                 | 8500.0000   |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                 | 0.9000         | 0.5180                       | 25.4000                 | 8500.0000   |
| 32  | 83-32-9  | acenaphthene         | 154.2100                 | 0.9000         | 0.0920                       | 3.4200                  | 4600.0000   |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                 | 0.9000         | 1.4800                       | 3.9300                  | 2500.0000   |
| 35  | 120-12-7 | anthracene           | 178.2400                 | 0.9000         | 1.0200                       | 0.0450                  | 14000.0000  |
| 42  | 71-43-2  | benzene              | 78.1100                  | 0.8740         | 5.5900                       | 1750.0000               | 83.0000     |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                 | 0.9000         | 0.0012                       | 0.0057                  | 138000.0000 |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                 | 0.9000         | 0.0016                       | 0.0012                  | 550000.0000 |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                 | 0.9000         | 0.0119                       | 0.0140                  | 55000.0000  |
| 38  | 218-01-9 | chrysene             | 228.3000                 | 0.9000         | 0.0011                       | 0.0018                  | 20000.0000  |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                 | 0.8670         | 6.4300                       | 152.0000                | 1100.0000   |
| 36  | 206-44-0 | fluoranthene         | 202.2600                 | 0.9000         | 0.0065                       | 0.2060                  | 38000.0000  |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                 | 0.9000         | 0.0642                       | 1.6900                  | 7300.0000   |
| 28  | 91-20-3  | naphthalene          | 128.1800                 | 0.9000         | 1.1500                       | 31.7000                 | 1300.0000   |
| 34  | 85-01-8  | phenanthrene         | 178.2400                 | 0.9000         | 0.1590                       | 1.0000                  | 14000.0000  |
| 37  | 129-00-0 | pyrene               | 202.2600                 | 0.9000         | 0.0050                       | 0.1320                  | 38000.0000  |
| 43  | 108-88-3 | toluene              | 92.1400                  | 0.8620         | 6.3700                       | 535.0000                | 300.0000    |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                 | 0.9000         | 0.0394                       | 0.0043                  | 55000.0000  |

\*\*\*\*\*  
END OF RECORDS

Sample Name Identification, BH-SED-18-0

Model used: Liquid saturated & water moisture content known,
Water moisture content (Volume Frac.), 0.7500,

NAPLANAL ANALYSIS RESULTS:

Table with 10 columns: ID#, Name, Total mass, Mass in water, Mass in soil, Mass in NAPL, Conc. in water, Sorbed in soil, Conc. in NAPL, Mole fraction in NAPL. Lists various chemical compounds and their concentrations.

(mg/kg)\* --- mg per kg of soil sample (wet soil)
(mg/kg)^ --- mg per kg of solid (dry soil)

Water Volume Frac.(l/l), 0.750000,
NAPL Volume Frac.(l/l), 0.000061,
Soil Volume Frac.(l/l), 0.249939,
Porosity (Volume Frac.), 0.750061,

Bulk Density (kg/l), 1.4124,
NAPL Density (kg/l), 0.9000,

NAPL Saturation (%),

0.0081,

Numerical Accuracy Information

The solution converged in 2 iterations with residual less than 1.0E-6.

The Thermodynamic Physical Properties of the Chemical Components Used in This Set of Soil Data Analysis

Database Source of Reference:

Combo.dbs (combination of P&jcalc.dbs and aromatic.dbs)  
Data for chlorinated compounds are from database of Pankow and Johnson with  
CALCULATED solubilities as presented in: Dense Chlorinated Solvents... by  
Pankow and Cherry, Waterloo Press, 1996. Data for (poly)aromatic hydrocarbons  
are from EPA/600/8-90/003, March 1990, Basics of Pump-and-Treat Tech. Liquid  
densities of solid phase compounds at 20-25'C are set at 0.9 g/mL though more  
accurate values can be obtained. Exceptions: Kocs for #49, 52, 53, and 54 are  
estimated from Eqtn 4-6, W.J. Lyman et al., Handbook... (ACS, 1990); properties  
for #52-54 are from Howard and Meylan, Handbook...(Lewis, 1997). #30 estimated.

| ID# | CAS#     | Name                 | Molecular Weight<br>(g/mol) | Density<br>(kg/l) | Henry's Constant<br>(atm-L/mol) | Water Solubility<br>(mg/L) | Koc<br>(ml/g) |
|-----|----------|----------------------|-----------------------------|-------------------|---------------------------------|----------------------------|---------------|
| 29  | 90-12-0  | 1-methylnaphthalene  | 142.2000                    | 0.9000            | 0.2600                          | 25.8000                    | 8500.0000     |
| 30  | 91-57-6  | 2-methylnaphthalene  | 142.2000                    | 0.9000            | 0.5180                          | 25.4000                    | 8500.0000     |
| 32  | 83-32-9  | acenaphthene         | 154.2100                    | 0.9000            | 0.0920                          | 3.4200                     | 4600.0000     |
| 31  | 208-96-8 | acenaphthylene       | 152.2000                    | 0.9000            | 1.4800                          | 3.9300                     | 2500.0000     |
| 35  | 120-12-7 | anthracene           | 178.2400                    | 0.9000            | 1.0200                          | 0.0450                     | 14000.0000    |
| 42  | 71-43-2  | benzene              | 78.1100                     | 0.8740            | 5.5900                          | 1750.0000                  | 83.0000       |
| 39  | 56-55-3  | benz(a)anthracene    | 228.3000                    | 0.9000            | 0.0012                          | 0.0057                     | 138000.0000   |
| 40  | 50-32-8  | benzo(a)pyrene       | 252.3200                    | 0.9000            | 0.0016                          | 0.0012                     | 550000.0000   |
| 50  | 205-99-2 | benzo(b)fluoranthene | 252.3200                    | 0.9000            | 0.0119                          | 0.0140                     | 550000.0000   |
| 38  | 218-01-9 | chrysene             | 228.3000                    | 0.9000            | 0.0011                          | 0.0018                     | 200000.0000   |
| 44  | 100-41-4 | ethylbenzene         | 106.1700                    | 0.8670            | 6.4300                          | 152.0000                   | 1100.0000     |
| 36  | 206-44-0 | fluoranthene         | 202.2600                    | 0.9000            | 0.0065                          | 0.2060                     | 38000.0000    |
| 33  | 86-73-7  | 9h-fluorene          | 166.2200                    | 0.9000            | 0.0642                          | 1.6900                     | 7300.0000     |
| 28  | 91-20-3  | naphthalene          | 128.1800                    | 0.9000            | 1.1500                          | 31.7000                    | 1300.0000     |
| 34  | 85-01-8  | phenanthrene         | 178.2400                    | 0.9000            | 0.1590                          | 1.0000                     | 14000.0000    |
| 37  | 129-00-0 | pyrene               | 202.2600                    | 0.9000            | 0.0050                          | 0.1320                     | 38000.0000    |
| 43  | 108-88-3 | toluene              | 92.1400                     | 0.8620            | 6.3700                          | 535.0000                   | 300.0000      |
| 51  | 207-08-9 | benzo(k)fluoranthene | 252.3200                    | 0.9000            | 0.0394                          | 0.0043                     | 550000.0000   |

END OF RECORDS

# **MASS DISTRIBUTION CALCULATIONS**



# **CALCULATION OF DISSOLVED HYDROCARBON MASS IN GROUNDWATER COKE POINT PENINSULA, BALTIMORE, MARYLAND**

## **OBJECTIVE**

The objective of the calculation was to determine mass of hydrocarbons in the groundwater in the onshore investigation area.

## **APPROACH**

1. Use isocontour maps of benzene and naphthalene concentrations in the shallow and intermediate aquifers to identify highly contaminated regions for each aquifer (Figures C-1 to C-4)
2. Determine the area of each region using a geographic information system (GIS).
3. Calculate geometric mean concentrations for benzene, toluene, and naphthalene from the sample concentrations measured for the applicable aquifer within each region.
4. Calculate the volume of groundwater present in each aquifer in each region, based on the area of the region and the average porosity and thickness of the aquifer.
5. Calculate the mass of benzene, toluene, and naphthalene present in each region, using the estimated concentrations and the calculated water volumes.
6. Concentration values are as listed on isocontour maps in URS 2005a, Tables D-9 and D-13.

## **CONCENTRATION CALCULATIONS**

- Calculate mass of benzene, toluene and naphthalene in the shallow and intermediate aquifers using table below (for regions, see Figures C-1 and C-2).

**CALCULATION OF DISSOLVED HYDROCARBON MASS IN GROUNDWATER  
COKE POINT PENINSULA, BALTIMORE, MARYLAND**

**SHALLOW AQUIFER**

| Section C-1-A         |                |                | Section C-1-B         |                |                |
|-----------------------|----------------|----------------|-----------------------|----------------|----------------|
| Sample                | Benzene (µg/L) | Toluene (µg/L) | Sample                | Benzene (µg/L) | Toluene (µg/L) |
| CO02PZM006            | 790,000        | 48,000         | CP07-PZM006           | 1,800          | 97             |
| CO03-PZM005           | 1              | 1              | CP08-PZM008           | 12,000         | 3,700          |
| CO07-PZM008           | 42,000         | 4,900          |                       |                |                |
| CO08-PZM005           | 5,500          | 4,500          |                       |                |                |
| CO15-PZM005           | 98,000         | 14,000         |                       |                |                |
| CO16-PZM006           | 200,000        | 15,000         |                       |                |                |
| CO17-PZM005           | 28,000         | 840            |                       |                |                |
| CO18-PZM006           | 1,100,000      | 71,000         |                       |                |                |
| CO19-PZM004           | 12,000         | 400            |                       |                |                |
| CO27-PZM012           | 25,000         | 5,800          |                       |                |                |
| CO28-PZM010           | 2,000          | 130            |                       |                |                |
| CO29-PZM010           | 25             | 8              |                       |                |                |
| CO30-PZM015           | 54,000         | 6,300          |                       |                |                |
| SW17-PZM007           | 20,000         | 2,500          |                       |                |                |
| <b>Geometric Mean</b> | <b>12,286</b>  | <b>1,565</b>   | <b>Geometric Mean</b> | <b>4,648</b>   | <b>599</b>     |

| Section C-1-C             |                |                | Section C-1-D             |                |                |
|---------------------------|----------------|----------------|---------------------------|----------------|----------------|
| Sample                    | Benzene (µg/L) | Toluene (µg/L) | Sample                    | Benzene (µg/L) | Toluene (µg/L) |
| CO25-PZM008               | 3,000          | 1,900          | CO05PZM006                | 2,100          | 1,700          |
| 1000 ug/L Benzene Contour | 1,000          | N/A            | 1000 ug/L Benzene Contour | 1,000          | N/A            |
| <b>Geometric Mean</b>     | <b>1,732</b>   | <b>1,900</b>   | <b>Geometric Mean</b>     | <b>1,449</b>   | <b>1,700</b>   |

**CALCULATION OF DISSOLVED HYDROCARBON MASS IN GROUNDWATER  
COKE POINT PENINSULA, BALTIMORE, MARYLAND**

| Section C-2-A         |                    | Section C-2-B         |                    |
|-----------------------|--------------------|-----------------------|--------------------|
| Sample                | Naphthalene (µg/L) | Sample                | Naphthalene (µg/L) |
| CO15-PZM005           | 1,200              | CO04-PZM004           | 6,700              |
| SW17-PZM007           | 1,200              | CO05-PZM006           | 8,100              |
| CO30-PZM015           | 2,100              | CO08-PZM005           | 1,400              |
|                       |                    | CO09-PZM007           | 2,000              |
|                       |                    | CO10-PZM006           | 1,100              |
|                       |                    | CO12-PZM008           | 2,300              |
|                       |                    | CO13-PZM008           | 22,000             |
|                       |                    | CO20-PZM004           | 1,700              |
|                       |                    | CO21-PZM005           | 2                  |
|                       |                    | CO22-PZM005           | 2,600              |
|                       |                    | CO23-PZM008           | 1,300              |
|                       |                    | CO24-PZM007           | 2,500              |
|                       |                    | CO25-PZM008           | 12,000             |
|                       |                    | CO26-PZM007           | 5,700              |
|                       |                    | TS08-PPM007           | 340                |
| <b>Geometric Mean</b> | <b>1,446</b>       | <b>Geometric Mean</b> | <b>1,861</b>       |

**INTERMEDIATE AQUIFER**

| Section C-3-A         |                |                |
|-----------------------|----------------|----------------|
| Sample                | Benzene (µg/L) | Toluene (µg/L) |
| CO28-PZM048           | 350,000        | 28,000         |
| CO27-PZM046           | 390,000        | 49,000         |
| CO02-PZM041           | 32,000         | 30             |
| CO08-PZM036           | 12,000         | 10,000         |
| SW17-PZM038           | 54,000         | 330            |
| <b>Geometric Mean</b> | <b>25,456</b>  | <b>1,817</b>   |

| Section C-4-A         |                    | Section C-4-B         |                    | Section C-4-C         |                    |
|-----------------------|--------------------|-----------------------|--------------------|-----------------------|--------------------|
| Sample                | Naphthalene (µg/L) | Sample                | Naphthalene (µg/L) | Sample                | Naphthalene (µg/L) |
| CO28-PZM048           | 1,900              | CO08-PZM036           | 3,500              | CP05-PZM028           | 1,300              |
| 1000 ug/L contour     | 1,000              | CO13-PZM030           | 3,100              | 1000 ug/L contour     | 1,000              |
|                       |                    | CO04-PZM048           | 1,100              |                       |                    |
| <b>Geometric Mean</b> | <b>1,378</b>       | <b>Geometric Mean</b> | <b>2,285</b>       | <b>Geometric Mean</b> | <b>1,140</b>       |

**CALCULATION OF DISSOLVED HYDROCARBON MASS IN GROUNDWATER  
COKE POINT PENINSULA, BALTIMORE, MARYLAND**

**AQUIFER VOLUME CALCULATIONS**

- Calculate the volume of each contaminated aquifer section

| <b>Section</b> | <b>Aquifer</b> | <b>Area (ft<sup>2</sup>)</b> | <b>Average<br/>Thickness (ft)</b> | <b>Porosity</b> | <b>Volume (ft<sup>3</sup>)</b> |
|----------------|----------------|------------------------------|-----------------------------------|-----------------|--------------------------------|
| C-1-A          | Shallow        | 3,600,000                    | 30                                | 0.25            | 6,180,000                      |
| C-1-B          | Shallow        | 349,000                      | 30                                | 0.25            | 2,617,500                      |
| C-1-C          | Shallow        | 112,000                      | 30                                | 0.25            | 840,000                        |
| C-1-D          | Shallow        | 92,600                       | 30                                | 0.25            | 694,500                        |
| C-2-A          | Shallow        | 824,000                      | 30                                | 0.25            | 61,800                         |
| C-2-B          | Shallow        | 2,610,000                    | 30                                | 0.25            | 19,575,000                     |
| C-3-A          | Intermediate   | 2,350,000                    | 40                                | 0.3             | 28,200,000                     |
| C-4-A          | Intermediate   | 322,000                      | 40                                | 0.3             | 3,864,000                      |
| C-4-B          | Intermediate   | 1,180,000                    | 40                                | 0.3             | 14,160,000                     |
| C-4-C          | Intermediate   | 108,000                      | 40                                | 0.3             | 1,296,000                      |

**CALCULATION OF DISSOLVED HYDROCARBON MASS IN GROUNDWATER  
COKE POINT PENINSULA, BALTIMORE, MARYLAND**

**CONTAMINANT MASS CALCULATIONS**

- Calculate the mass of the contaminants in each aquifer section, and the total mass in all sections

| Section            | Aquifer      | Volume (ft <sup>3</sup> ) | Mean Concentration (µg/L) | Contaminant Mass (µg) | Contaminant Mass (lb) |
|--------------------|--------------|---------------------------|---------------------------|-----------------------|-----------------------|
| <b>Benzene</b>     |              |                           |                           |                       |                       |
| C-1-A              | Shallow      | 6,180,000                 | 12,286                    | 2.15E+12              | 4.74E+03              |
| C-1-B              | Shallow      | 6,180,000                 | 4,648                     | 8.13E+11              | 1.79E+03              |
| C-1-C              | Shallow      | 6,180,000                 | 1,732                     | 3.03E+11              | 6.68E+02              |
| C-1-D              | Shallow      | 6,180,000                 | 1,449                     | 2.54E+11              | 5.59E+02              |
| C-3-A              | Intermediate | 28,200,000                | 25,456                    | 2.03E+13              | 4.48E+04              |
|                    |              |                           | <b>Total</b>              | <b>2.38E+13</b>       | <b>5.26E+04</b>       |
| <b>Toluene</b>     |              |                           |                           |                       |                       |
| C-1-A              | Shallow      | 6,180,000                 | 1,565                     | 2.74E+11              | 6.04E+02              |
| C-1-B              | Shallow      | 6,180,000                 | 599                       | 1.05E+11              | 2.31E+02              |
| C-1-C              | Shallow      | 6,180,000                 | 1,900                     | 3.32E+11              | 7.33E+02              |
| C-1-D              | Shallow      | 6,180,000                 | 1,700                     | 2.97E+11              | 6.56E+02              |
| C-3-A              | Intermediate | 28,200,000                | 1,817                     | 1.45E+12              | 3.20E+03              |
|                    |              |                           | <b>Total</b>              | <b>2.46E+12</b>       | <b>5.42E+03</b>       |
| <b>Naphthalene</b> |              |                           |                           |                       |                       |
| C-2-A              | Shallow      | 61,800                    | 1,446                     | 2.53E+09              | 5.58E+00              |
| C-2-B              | Shallow      | 19,575,000                | 1,861                     | 1.03E+12              | 2.27E+03              |
| C-4-A              | Intermediate | 3,864,000                 | 1,378                     | 1.51E+11              | 3.33E+02              |
| C-4-B              | Intermediate | 14,160,000                | 2,285                     | 9.16E+11              | 2.02E+03              |
| C-4-C              | Intermediate | 1,296,000                 | 1,140                     | 4.18E+10              | 9.23E+01              |
|                    |              |                           | <b>Total</b>              | <b>2.14E+12</b>       | <b>4.73E+03</b>       |

|              |          |
|--------------|----------|
| Total pounds | 6.27E+04 |
|--------------|----------|

**CONCLUSION**

The highly contaminated groundwater areas contain approximately 53,000 lb. of benzene, 5,400 lb. of toluene, and 4,700 lb. of naphthalene.

Site Assessment for Proposed Coke Point Dredged Material Containment Facility at Sparrows Point



Legend

- Sampling Locations ●
- Area of Concern ■

Sources  
 ESRI, i-cubed, GeoEye, 2009  
 Tele Atlas North America Inc., ESRI, 2006



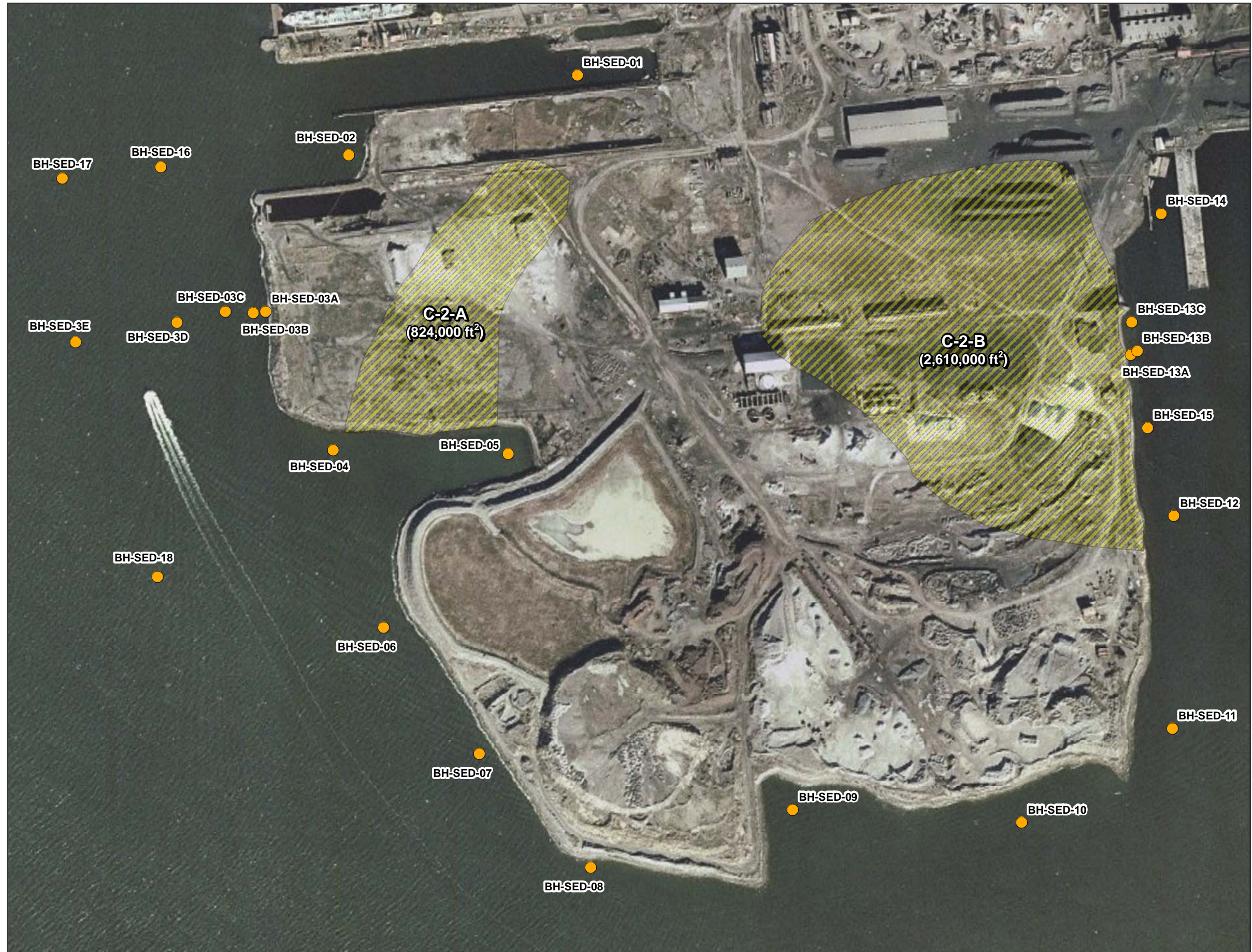
**EA** EA ENGINEERING, SCIENCE, AND TECHNOLOGY, INC.



Figure C-1. Shallow Benzene, Coke Point Peninsula, Baltimore Maryland

H:\projects\1453406\WXD\2009\_Report\Figure C-1

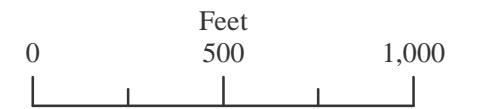
Site Assessment for Proposed Coke Point Dredged Material Containment Facility at Sparrows Point



Legend

- Sampling Locations ●
- Area of Concern

Sources  
 ESRI, i-cubed, GeoEye, 2009  
 Tele Atlas North America Inc., ESRI, 2006



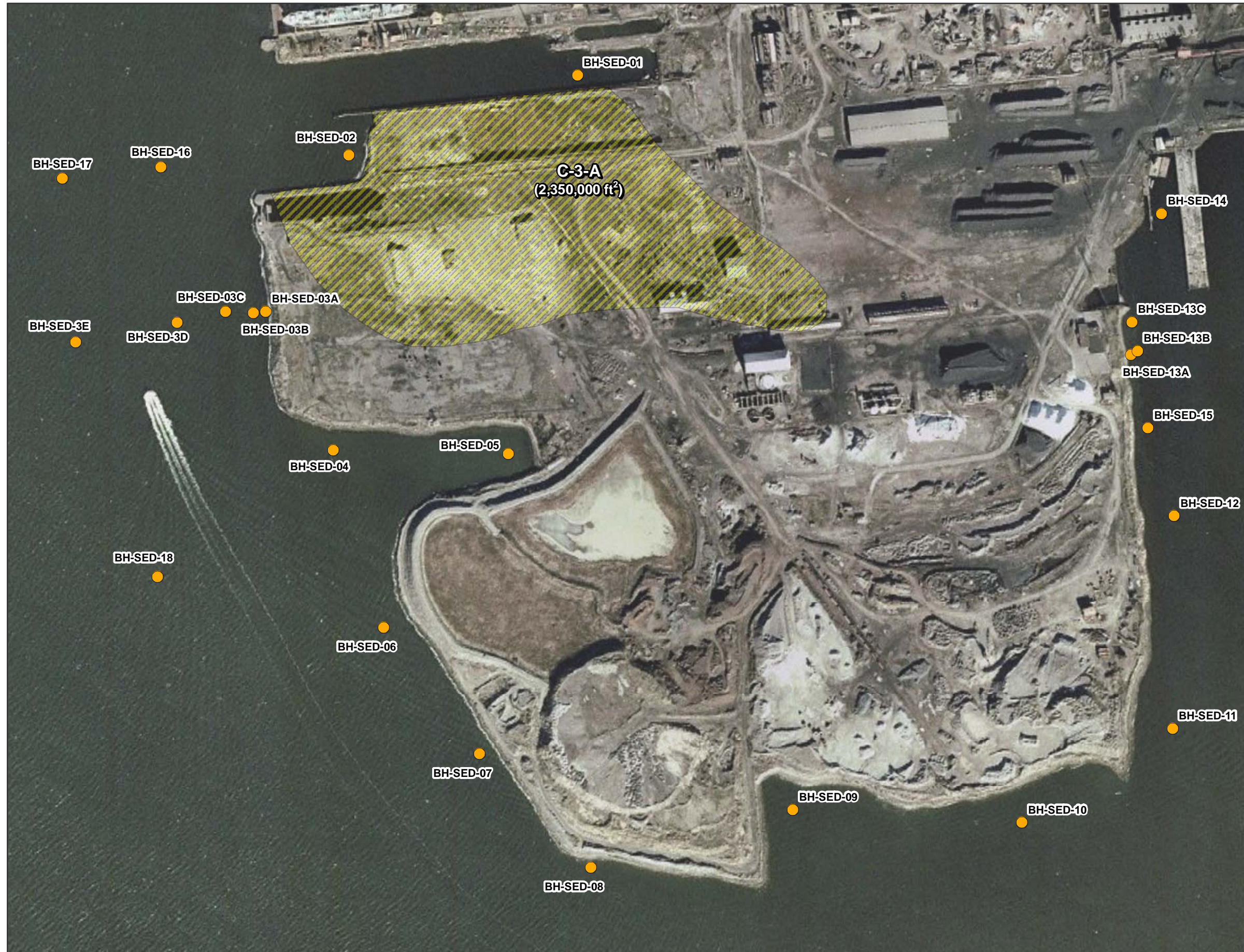
**EA** EA ENGINEERING, SCIENCE, AND TECHNOLOGY, INC.



Figure C-2. Shallow Naphthalene, Coke Point Peninsula, Baltimore Maryland

H:\projects\1453406\WXD\2009\_Report\Figure C-2

Site Assessment for Proposed Coke Point Dredged Material Containment Facility at Sparrows Point



Legend

- Sampling Locations ●
- Area of Concern

Sources  
 ESRI, i-cubed, GeoEye, 2009  
 Tele Atlas North America Inc., ESRI, 2006



**EA** EA ENGINEERING, SCIENCE, AND TECHNOLOGY, INC.



H:\projects\1453406\WXD\2009\_Report\Figure C-3

Figure C-3. Intermediate Benzene, Coke Point Peninsula, Baltimore Maryland



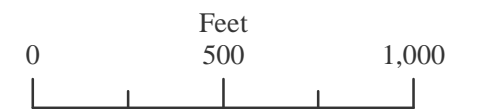
Site Assessment for Proposed Coke Point Dredged Material Containment Facility at Sparrows Point



**Legend**

- Sampling Locations ●
- Area of Concern

Sources  
 ESRI, i-cubed, GeoEye, 2009  
 Tele Atlas North America Inc., ESRI, 2006



**EA** EA ENGINEERING, SCIENCE, AND TECHNOLOGY, INC.



Figure C-4. Intermediate Naphthalene, Coke Point Peninsula, Baltimore Maryland

H:\projects\1453406\WXD\2009\_Report\Figure C-4

**CALCULATION OF HYDROCARBON MASS IN ONSHORE SOILS  
COKE POINT PENINSULA, BALTIMORE, MARYLAND**

**OBJECTIVE**

The objective of the calculation was to determine volume and mass of hydrocarbons sorbed within the soil matrix for the onshore investigation areas.

**APPROACH**

1. Estimate an "equivalent thickness" of sorbed hydrocarbon mass using porosity, saturation percent (expressed as a fraction), and thickness of sample interval
2. Calculate volume using geometric mean of equivalent thickness multiplied by the square footage of each area of concern
  - Assumes sorbed mass is log-normally distributed over each area of concern, which were bounded based on groundwater impacts that should correlate with impacted fill
3. Calculate mass using NAPL density (from Table 3-10)
4. Calculations use rounded numbers and conservative assumptions

**CALCULATION (BENZOL PROCESSING AREA)**

- Calculate equivalent thickness of sorbed hydrocarbons in Benzol Processing Area using table below

**Equivalent Thickness of Sorbed Hydrocarbons from NAPL Saturation  
Benzol Processing Area**

| Borehole ID                                     | Sample ID    | Estimated Thickness of Impacted Soil <sup>1</sup> (ft) | Hydrocarbon Pore Saturation <sup>2</sup> | Total Porosity | Equivalent Thickness (ft) |
|-------------------------------------------------|--------------|--------------------------------------------------------|------------------------------------------|----------------|---------------------------|
| BP-B01                                          | BP-SO-B01-8  | 1                                                      | 0.0523                                   | 0.22           | 0.0115                    |
|                                                 | BP-SO-B01-14 | 4                                                      | 0.0377                                   | 0.25           | 0.0377                    |
|                                                 | BP-SO-B01-20 | 2                                                      | 0.0009                                   | 0.37           | 0.0007                    |
| BP-B02                                          | BP-SO-B02-8  | 6                                                      | 0.0002                                   | 0.30           | 0.0004                    |
|                                                 | BP-SO-B02-20 | 2                                                      | 0.0013                                   | 0.33           | 0.0009                    |
| BP-B04                                          | BP-SO-B04-24 | 2                                                      | 0.0001                                   | 0.25           | 0.0001                    |
| BP-B06                                          | BP-SO-B06-8  | 4                                                      | 0.0093                                   | 0.19           | 0.0071                    |
|                                                 | BP-SO-B06-12 | 4                                                      | 0.0181                                   | 0.26           | 0.0188                    |
| BP-B07                                          | BP-SO-B07-12 | 5                                                      | 0.0078                                   | 0.16           | 0.0062                    |
| BP-B09                                          | BP-SO-B09-14 | 3                                                      | 0.1073                                   | 0.18           | 0.0579                    |
|                                                 | BP-SO-B09-18 | 3                                                      | 0.0582                                   | 0.33           | 0.0576                    |
| <b>Geometric Mean Equivalent Thickness (ft)</b> |              |                                                        |                                          |                | <b>0.0046</b>             |

<sup>1</sup> Based on interpretations from boring log descriptions

<sup>2</sup> Values from saturation percent in Table 3-13 expressed as a fraction.  
Values less than 0.0001 excluded from calculations.

**CALCULATION OF HYDROCARBON MASS IN ONSHORE SOILS  
COKE POINT PENINSULA, BALTIMORE, MARYLAND**

- Benzol Processing area is 424,000 ft<sup>2</sup> (Figure C-5)  
This is based on the previously defined impacted area (440,000 ft<sup>2</sup>), with overlapping mobile NAPL area (16,000 ft<sup>2</sup>) subtracted.
- Calculate approximate volume of sorbed hydrocarbons (as equivalent NAPL)  
(424,000 ft<sup>2</sup>) \* 0.0046 ft = 1,950 ft<sup>3</sup>  
(1,950 ft<sup>3</sup>) \* (7.48 gal/ft<sup>3</sup>) = 14,600 gal equivalent NAPL
- Convert to pounds of equivalent NAPL  
NAPL specific gravity in Benzol Processing Area is 0.8906 at 70 °F  
One gallon of water weighs 8.329 lb at 70 °F  
Therefore NAPL weight is (0.8906) \* (8.329 lb/gal) = 7.42 lb/gal  
(14,600 gal) \* (7.42 lbs/gal) = 108,000 lbs equivalent NAPL

**CALCULATION (COAL TAR STORAGE AREA)**

- Calculate equivalent thickness of sorbed hydrocarbons in Coal Tar Storage Area using table below

**Equivalent Thickness of Sorbed Hydrocarbons from NAPL Saturation  
Coal Tar Storage Area**

| Borehole ID                                     | Sample ID    | Estimated Thickness of Impacted Soil <sup>1</sup> (ft) | Hydrocarbon Pore Saturation <sup>2</sup> | Total Porosity | Equivalent Thickness (ft) |
|-------------------------------------------------|--------------|--------------------------------------------------------|------------------------------------------|----------------|---------------------------|
| CT-B01                                          | CT-SO-B01-18 | 2                                                      | 0.0351                                   | 0.24           | 0.0169                    |
| CT-B02                                          | CT-SO-B02-16 | 2                                                      | 0.0006                                   | 0.14           | 0.0002                    |
|                                                 | CT-SO-B02-20 | 2                                                      | 0.0103                                   | 0.20           | 0.0041                    |
| CT-B03                                          | CT-SO-B03-20 | 2                                                      | 0.0065                                   | 0.16           | 0.0021                    |
| CT-B04                                          | CT-SO-B04-18 | 2                                                      | 0.0077                                   | 0.24           | 0.0037                    |
| CT-B05                                          | CT-SO-B05-20 | 2                                                      | 0.0101                                   | 0.20           | 0.0040                    |
| <b>Geometric Mean Equivalent Thickness (ft)</b> |              |                                                        |                                          |                | <b>0.0027</b>             |

<sup>1</sup> Based on interpretations from boring log descriptions

<sup>2</sup> Values from saturation percent in Table 3-13 expressed as a fraction.

Values less than 0.001 excluded from calculations.

- Coal Tar Storage area is 398,000 ft<sup>2</sup> (Figure C-5)
- Calculate approximate volume of sorbed hydrocarbons (as equivalent NAPL)  
(398,000 ft<sup>2</sup>) \* 0.0027 ft = 1,100 ft<sup>3</sup>  
(1,100 ft<sup>3</sup>) \* (7.48 gal/ft<sup>3</sup>) = 7,900 gal equivalent NAPL
- Convert to pounds of equivalent NAPL  
NAPL specific gravity in Coal Tar Storage Area is 1.1531 at 70 °F  
One gallon of water weighs 8.329 lb at 70 °F  
Therefore NAPL weight is (1.1531) \* (8.329 lb/gal) = 9.60 lb/gal  
(7,900 gal) \* (9.60 lbs/gal) = 76,000 lbs equivalent NAPL

**CALCULATION OF HYDROCARBON MASS IN ONSHORE SOILS  
COKE POINT PENINSULA, BALTIMORE, MARYLAND**

**CONCLUSION**

The estimated volume of equivalent NAPL within the Benzol Processing area is 14,000 gallons.  
The estimated mass of equivalent NAPL is 108,000 pounds.

The estimated volume of equivalent NAPL within the Coal Tar Storage area is 7,900 gallons.  
The estimated mass of equivalent NAPL is 76,000 pounds.

This gives a total mass equivalent of 184,000 pounds for the onshore investigation areas.

Site Assessment for Proposed Coke Point Dredged Material Containment Facility at Sparrows Point



Legend

- Sampling Locations ●
- Area of Concern ■

Sources  
 ESRI, i-cubed, GeoEye, 2009  
 Tele Atlas North America Inc., ESRI, 2006



**EA** EA ENGINEERING, SCIENCE, AND TECHNOLOGY, INC.



H:\projects\1453406\MXD\2009\_Report\Figure C-5

Figure C-5. Estimated Extent of Onshore Residual NAPL, Coke Point Peninsula, Baltimore Maryland

**CALCULATION OF NAPL VOLUME AND MASS  
BENZOL PROCESSING AREA, COKE POINT PENINSULA, BALTIMORE, MARYLAND**

**OBJECTIVE**

The objective of the calculation was to determine volume and mass of LNAPL within the Benzol Processing area based on NAPL gauging conducted on 23 August 2009.

**APPROACH**

1. Estimate the footprint of NAPL occurrence from gauging results
2. Obtain the area of the NAPL footprint using a GIS
3. Assume an average NAPL thickness from gauging results
4. Calculate NAPL volume using an assumed soil porosity and NAPL saturation
5. Calculate NAPL mass using gasoline density to approximate NAPL density
6. Calculations use rounded numbers and conservative assumptions

**CALCULATION**

- Area of NAPL footprint as shown on Figure C-6 is 33,000 ft<sup>2</sup>
- Calculate average thickness of NAPL from wells BP-MW-05, -08, and -10  
 $(3.59 + 4.64 + 0.62) / 3 = 3 \text{ ft}$
- Calculate approximate volume of aquifer material encompassed by NAPL  
 $(33,000 \text{ ft}^2) * 3 \text{ ft} = 99,000 \text{ ft}^3$
- Calculate approximate volume of NAPL  
Assume total porosity of 0.3  
Assume NAPL saturation of 50% of total porosity  
Therefore, NAPL volume = 0.015 times the aquifer volume, or 1,500 ft<sup>3</sup>
- Convert NAPL volume from cubic feet to gallons  
Conversion factor is 7.48 gal per ft<sup>3</sup>  
 $(1,500 \text{ ft}^3) * (7.48 \text{ gal/ft}^3) = 11,000 \text{ gal}$
- Convert gallons of NAPL into pounds of NAPL  
NAPL specific gravity is 0.8906 at 70 °F  
Water density is 8.329 lb/gal at 70 °F  
NAPL density is  $(0.8906) * (8.327 \text{ lb/gal}) = 7.42 \text{ lb/gal}$   
 $(11,000 \text{ gal}) * (7.42 \text{ lbs/gal}) = 82,000 \text{ lbs NAPL}$

**CONCLUSION**

The estimated volume of NAPL within the Benzol Processing area is 11,000 gallons.  
The estimated mass of NAPL is 82,000 pounds.

Site Assessment for Proposed Coke Point Dredged Material Containment Facility at Sparrows Point



**Legend**

- Sampling Locations ●
- Area of Concern

Sources  
 ESRI, i-cubed, GeoEye, 2009  
 Tele Atlas North America Inc., ESRI, 2006



**EA** EA ENGINEERING, SCIENCE, AND TECHNOLOGY, INC.



H:\projects\1453406\WXD\2009\_Report\Figure C-6

Figure C-6. Floating Product, Coke Point Peninsula, Baltimore Maryland

**CALCULATION OF HYDROCARBON MASS IN OFFSHORE SEDIMENTS  
BENZOL PROCESSING AREA, COKE POINT PENINSULA, BALTIMORE, MARYLAND**

**OBJECTIVE**

The objective of the calculation was to determine volume and mass of hydrocarbons sorbed within the surface and subsurface sediment for the offshore investigation area

**APPROACH**

1. Estimate an "equivalent thickness" of sediment hydrocarbon mass using porosity, saturation percent (expressed as a fraction), and thickness of sample interval
2. Calculate volume using geometric mean equivalent thickness over the approximate areas of impacted sediment (Figure C-7)
  - Assumes sorbed mass is log-normally distributed over the impacted areas
3. Calculate mass using NAPL specific gravity from Table 3-10
4. Calculations use rounded numbers and conservative assumptions

**CALCULATION (OFFSHORE INVESTIGATION AREA)**

- Calculate equivalent thickness of sorbed hydrocarbons in offshore sediment using table below

**Equivalent Thickness of Sorbed Hydrocarbons from NAPL Saturation  
Surrounding the Coke Point Peninsula**

| <b>Borehole ID</b>                              | <b>Sample ID</b> | <b>Estimated Thickness of Impacted Sediment<sup>1</sup> (ft)</b> | <b>Hydrocarbon Pore Saturation<sup>2</sup></b> | <b>Total Porosity</b> | <b>Equivalent Thickness (ft)</b> |
|-------------------------------------------------|------------------|------------------------------------------------------------------|------------------------------------------------|-----------------------|----------------------------------|
| BH-SED-02                                       | BH-SED-02-00     | 1                                                                | 0.0009                                         | 0.34                  | 0.0003                           |
|                                                 | BH-SED-02-4      | 4                                                                | 0.0017                                         | 0.15                  | 0.0010                           |
| BH-SED-03A                                      | BH-SED-3A-00     | 1                                                                | 0.0002                                         | 0.43                  | 0.0001                           |
| BH-SED-03B                                      | BH-SED-3B-00     | 1                                                                | 0.0201                                         | 0.64                  | 0.0129                           |
| BH-SED-03C                                      | BH-SED-3C-00     | 2                                                                | 0.0005                                         | 0.69                  | 0.0006                           |
|                                                 | BH-SED-3C-2      | 3                                                                | 0.0082                                         | 0.60                  | 0.0147                           |
|                                                 | BH-SED-04-8      | 6                                                                | 0.0123                                         | 0.59                  | 0.0437                           |
| BH-SED-05                                       | BH-SED-05-00     | 1                                                                | 0.0016                                         | 0.38                  | 0.0006                           |
|                                                 | BH-SED-05-4      | 5                                                                | 0.0003                                         | 0.46                  | 0.0006                           |
| BH-SED-06                                       | BH-SED-06-00     | 1                                                                | 0.0003                                         | 0.59                  | 0.0002                           |
|                                                 | BH-SED-06-6      | 5                                                                | 0.0025                                         | 0.58                  | 0.0073                           |
| BH-SED-07                                       | BH-SED-07-00     | 1                                                                | 0.0018                                         | 0.51                  | 0.0009                           |
| BH-SED-08                                       | BH-SED-08-00     | 1                                                                | 0.0001                                         | 0.64                  | 0.0001                           |
| BH-SED-09                                       | BH-SED-09-00     | 1                                                                | 0.0002                                         | 0.49                  | 0.0001                           |
| BH-SED-10                                       | BH-SED-10-00     | 1                                                                | 0.0001                                         | 0.59                  | 0.0001                           |
| BH-SED-11                                       | BH-SED-11-00     | 1                                                                | 0.0001                                         | 0.58                  | 0.0001                           |
|                                                 | BH-SED-11-02     | 2                                                                | 0.0076                                         | 0.54                  | 0.0082                           |
| <b>Geometric Mean Equivalent Thickness (ft)</b> |                  |                                                                  |                                                |                       | <b>0.0008</b>                    |



**CALCULATION OF HYDROCARBON MASS IN OFFSHORE SEDIMENTS  
BENZOL PROCESSING AREA, COKE POINT PENINSULA, BALTIMORE, MARYLAND**

<sup>1</sup> Based on interpretations from boring log descriptions

<sup>2</sup> Values from saturation percent in Table 3-13 expressed as a fraction.

Values less than 0.0001 excluded from calculations.

- offshore impacted areas are 4,004,000 ft<sup>2</sup> (Figure C-7)  
 $(4,004,000 \text{ ft}^2) * 0.0008 \text{ ft} = 3200 \text{ ft}^3$
- Calculate approximate volume of sorbed hydrocarbons (as equivalent NAPL)  
 $(3,200 \text{ ft}^3) * (7.48 \text{ gal/ft}^3) = 24,000 \text{ gal equivalent NAPL}$
- Convert to pounds of equivalent NAPL  
Primary contaminants in sediment are PAHs. Assume NAPL specific gravity  
is the same as in Coal Tar Storage area (1.1531 at 70 °F)  
One gallon of water weighs 8.329 lb at 70 °F  
Therefore NAPL weight is  $(1.1531) * (8.329 \text{ lb/gal}) = 9.60 \text{ lb/gal}$   
 $(24,000 \text{ gal}) * (9.60 \text{ lbs/gal}) = 230,000 \text{ lbs equivalent NAPL}$

**CONCLUSION**

The estimated volume of NAPL within the impacted offshore areas is 24,000 gallons.

The estimated mass of NAPL is 230,000 pounds.

Site Assessment for Proposed Coke Point Dredged Material Containment Facility at Sparrows Point



Legend

- Sampling Locations ●
- Area of Concern ■

Sources  
 ESRI, i-cubed, GeoEye, 2009  
 Tele Atlas North America Inc., ESRI, 2006



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Figure C-7. Offshore Areas of Concern, Coke Point Peninsula, Baltimore Maryland

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