

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010715\
 Data File : BM000138.D
 Acq On : 07 Jan 2015 21:46
 Operator : TP/IZ
 Sample : MDL-03-S
 Misc : MDL-SOIL-0.07PPM-SIM
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-03-S

Quant Time: Jan 08 12:05:23 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM010715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 08 10:41:47 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	2960	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	9505	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5576	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	10488	0.40	ng/ul	0.02
16) Chrysene-d12	21.41	240	8958	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	8174	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.22	152	4805	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	8534	0.38	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.68	128	1417	0.06	ng/ul#	87
5) 2-Methylnaphthalene	12.29	142	885	0.05	ng/ul	97
7) Acenaphthylene	14.20	152	1415	0.05	ng/ul#	88
8) Acenaphthene	14.54	153	956	0.05	ng/ul	95
9) Fluorene	15.52	166	1060	0.05	ng/ul#	85
11) Pentachlorophenol	16.90	266	15	0.01	ng/ul#	44
12) Phenanthrene	17.26	178	1780	0.05	ng/ul#	89
13) Anthracene	17.35	178	1674	0.06	ng/ul#	88
15) Fluoranthene	19.27	202	2061	0.06	ng/ul	91
17) Pyrene	19.64	202	2114	0.06	ng/ul#	93
18) Benzo(a)anthracene	21.39	228	1690	0.06	ng/ul	92
19) Chrysene	21.44	228	1821	0.06	ng/ul	96
21) Benzo(b)fluoranthene	23.01	252	1717	0.05	ng/ul	65
22) Benzo(k)fluoranthene	23.05	252	1700	0.06	ng/ul#	69
23) Benzo(a)pyrene	23.60	252	1707	0.06	ng/ul#	57
24) Indeno(1,2,3-cd)pyrene	26.05	276	1830	0.06	ng/ul#	73
25) Dibenzo(a,h)anthracene	26.07	278	1432	0.05	ng/ul#	59
26) Benzo(g,h,i)perylene	26.75	276	1666	0.06	ng/ul#	77

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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