

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM022015\
 Data File : BM000583.D
 Acq On : 21 Feb 2015 07:32
 Operator : TP/IZ
 Sample : MDL-06-S-ML
 Misc : MDL-SOIL-3PPM-SOM0-ML
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-06-S-ML

Quant Time: Feb 23 15:52:20 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-MA-BM022015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Feb 21 00:39:15 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	100481	20.00	ng/ul	0.00
16) Naphthalene-d8	10.46	136	427585	20.00	ng/ul	0.00
33) Acenaphthene-d10	14.32	164	272383	20.00	ng/ul	0.00
59) Phenanthrene-d10	17.06	188	570782	20.00	ng/ul	-0.01
75) Chrysene-d12	21.26	240	481487	20.00	ng/ul	0.00
83) Perylene-d12	23.47	264	402173	20.00	ng/ul	0.00

System Monitoring Compounds

3) Phenol-d5	6.85	99	191644	26.20	ng/ul	0.00
5) Bis-(2-Chloroethyl)ether-d	7.00	67	122548	28.35	ng/ul	0.00
7) 2-Chlorophenol-d4	7.20	132	171361	27.54	ng/ul	-0.01
11) 4-Methylphenol-d8	8.38	113	165630	26.71	ng/ul	0.00
17) Nitrobenzene-d5	8.83	128	80172	27.94	ng/ul	0.00
20) 2-Nitrophenol-d4	9.55	143	95820	28.85	ng/ul	0.00
24) 2,4-Dichlorophenol-d3	10.08	165	177253	25.19	ng/ul	0.00
27) 4-Chloroaniline-d4	10.60	131	251151	40.01	ng/ul	0.00
41) Dimethylphthalate-d6	13.73	166	578463	29.34	ng/ul	0.00
44) Acenaphthylene-d8	14.01	160	742283	28.10	ng/ul	0.00
49) 4-Nitrophenol-d4	14.53	143	69256	21.97	ng/ul	-0.01
55) Fluorene-d10	15.32	176	538748	27.96	ng/ul	0.00
60) 4,6-Dinitro-2-methylphenol	15.44	200	63277	19.07	ng/ul	0.00
68) Anthracene-d10	17.16	188	776136	28.75	ng/ul	0.00
76) Pyrene-d10	19.46	212	766105	27.34	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.33	264	545906	28.95	ng/ul	0.00

Target Compounds

					Qvalue
70) 1,2,3,4-Tetrachlorobenzene	13.12	216	16421	1.94 ng/uL	96
71) Pentachlorobenzene	14.64	250	31205	3.85 ng/uL	98

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

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