

Data Path : Z:\HPCHEM1\BNA_M\Data\BM010215\
 Data File : BM000102.D
 Acq On : 02 Jan 2015 10:16
 Operator : TP/IZ
 Sample : BLK-S-MDL-ML
 Misc : MDL-SOIL-BLK-ML
 ALS Vial : 11 Sample Multiplier: 1

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

Quant Time: Jan 03 04:42:24 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM122914.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 31 02:57:11 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	325926	20.00	ng/ul	0.00
18) Naphthalene-d8	10.63	136	1387775	20.00	ng/ul	-0.01
35) Acenaphthene-d10	14.47	164	992537	20.00	ng/ul	-0.01
61) Phenanthrene-d10	17.21	188	1951220	20.00	ng/ul	-0.01
75) Chrysene-d12	21.40	240	1661673	20.00	ng/ul	-0.01
83) Perylene-d12	23.69	264	1443676	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.32	96	38464	6.84	ng/uL	-0.01
5) Phenol-d5	7.00	99	1082418	38.34	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.17	67	601039	34.82	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.37	132	771419	38.37	ng/ul	0.00
13) 4-Methylphenol-d8	8.54	113	830709	37.24	ng/ul	0.00
19) Nitrobenzene-d5	9.00	128	364908	38.02	ng/ul	0.00
22) 2-Nitrophenol-d4	9.72	143	390266	39.89	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.24	165	795370	36.60	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.77	131	1001945	40.81	ng/ul	0.00
43) Dimethylphthalate-d6	13.89	166	2605721	35.60	ng/ul	0.00
46) Acenaphthylene-d8	14.16	160	2995486	34.16	ng/ul	-0.01
51) 4-Nitrophenol-d4	14.67	143	372803	31.53	ng/ul	0.00
57) Fluorene-d10	15.47	176	2107654	33.53	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.58	200	168943	17.46	ng/ul	0.00
70) Anthracene-d10	17.32	188	3177566	35.22	ng/ul	0.00
76) Pyrene-d10	19.61	212	3160695	41.23	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.56	264	2337016	35.87	ng/ul	0.00

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

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