

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM022015\
 Data File : BM000560.D
 Acq On : 20 Feb 2015 14:57
 Operator : TP/IZ
 Sample : MDL-01-W
 Misc : MDL-WATER-3PPM-SOM0
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-01-W

Quant Time: Feb 21 03:05:50 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	108016	20.00	ng/ul	0.00
16) Naphthalene-d8	10.46	136	477686	20.00	ng/ul	0.00
33) Acenaphthene-d10	14.33	164	319561	20.00	ng/ul	0.00
59) Phenanthrene-d10	17.07	188	699101	20.00	ng/ul	0.00
75) Chrysene-d12	21.26	240	524988	20.00	ng/ul	0.00
83) Perylene-d12	23.47	264	390379	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) Phenol-d5	6.85	99	251052	31.93	ng/ul	0.00
5) Bis-(2-Chloroethyl)ether-d	7.01	67	164117	35.32	ng/ul	0.00
7) 2-Chlorophenol-d4	7.20	132	220589	32.98	ng/ul	0.00
11) 4-Methylphenol-d8	8.39	113	221241	33.19	ng/ul	0.00
17) Nitrobenzene-d5	8.83	128	112277	35.03	ng/ul	0.00
20) 2-Nitrophenol-d4	9.55	143	132357	35.67	ng/ul	0.00
24) 2,4-Dichlorophenol-d3	10.08	165	243979	31.03	ng/ul	0.00
27) 4-Chloroaniline-d4	10.60	131	345773	49.31	ng/ul	0.00
41) Dimethylphthalate-d6	13.74	166	866136	37.45	ng/ul	0.00
44) Acenaphthylene-d8	14.02	160	1059479	34.19	ng/ul	0.00
49) 4-Nitrophenol-d4	14.53	143	119647	32.35	ng/ul	0.00
55) Fluorene-d10	15.32	176	782674	34.63	ng/ul	0.00
60) 4,6-Dinitro-2-methylphenol	15.44	200	107422	26.44	ng/ul	0.00
68) Anthracene-d10	17.17	188	1168355	35.33	ng/ul	0.00
76) Pyrene-d10	19.47	212	1139741	37.30	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.33	264	667150	36.45	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
70) 1,2,3,4-Tetrachlorobenzene	13.12	216	20383	1.97	ng/uL	99
71) Pentachlorobenzene	14.65	250	40519	4.08	ng/uL	97

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

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