

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
 Data File : BM000849.D
 Acq On : 04 Apr 2015 00:46
 Operator : TP/IZ
 Sample : MDL-01-W
 Misc : MDL-WATER-0.2/0.4
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MDL-01-W

Quant Time: Apr 04 04:18:12 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 04 00:49:41 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	19548	5.00	ng	0.00
11) Naphthalene-d8	10.39	136	98108	5.00	ng	0.00
19) Acenaphthene-d10	14.26	164	51442	5.00	ng	0.00
26) Phenanthrene-d10	17.00	188	119765	5.00	ng	0.00
30) Chrysene-d12	21.21	240	98647	5.00	ng	0.00
34) Perylene-d12	23.39	264	91124	5.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.21	112	75205	13.97	ng	0.00
5) Phenol-d6	6.77	99	133226	12.86	ng	0.00
12) Nitrobenzene-d5	8.77	82	34656	7.57	ng	0.00
20) 2,4,6-Tribromophenol	15.77	330	19625	9.01	ng	0.00
23) 2-Fluorobiphenyl	12.87	172	154679	9.76	ng	0.00
32) Terphenyl-d14	19.65	244	132545	11.16	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.12	88	873	0.21	ng	# 62
3) n-Nitrosodimethylamine	3.46	42	316	0.13	ng	# 1
6) 2-Chlorophenol	7.17	128	1210	0.37	ng	86
7) Phenol	6.81	94	1276	0.16	ng	79
8) bis(2-Chloroethyl)ether	7.02	93	1030m	0.35	ng	
9) 2-Methylphenol	8.06	107	900	0.29	ng	# 86
10) 3+4-Methylphenols	8.38	107	889	0.24	ng	# 82
13) Nitrobenzene	8.80	77	928	0.17	ng	# 71
14) 2-Nitrophenol	9.52	139	333	0.25	ng	# 37
15) 2,4-Dimethylphenol	9.58	122	632	0.11	ng	88
16) 2,4-Dichlorophenol	10.07	162	641	0.32	ng	85
17) Hexachlorobutadiene	10.72	225	601	0.16	ng	# 68
18) 4-Chloro-3-methylphenol	11.69	107	712	0.14	ng	94
21) 2,4,6-Trichlorophenol	12.70	196	310	0.20	ng	# 55
22) 2,4,5-Trichlorophenol	12.77	196	406	0.16	ng	# 82
25) 4-Nitrophenol	14.54	139	89	0.29	ng	92
27) 4,6-Dinitro-2-methylphenol	15.41	198	64	0.38	ng	# 80
28) Hexachlorobenzene	16.33	284	910	0.19	ng	# 82
29) Pentachlorophenol	16.69	266	41	0.31	ng	# 69
31) Benzidine	19.29	184	1245	0.57	ng	# 73
33) 3,3'-Dichlorobenzidine	21.13	252	995	0.30	ng	# 86

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

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