Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000854.D
Acq On : 04 Apr 2015 04:21
Operator : TP/IZ
Sample : MDL-06-W
Misc : MDL-WATER-0.2/0.4
ALS Vial : 43 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : MDL-06-W

Quant Time: Apr 04 05:14:34 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	18775	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	95266	5.00	ng		0.00
19) Acenaphthene-d10	14.26	164	48950	5.00	ng		0.00
26) Phenanthrene-d10	17.00	188	109087	5.00	ng		0.00
30) Chrysene-d12	21.21	240	89217	5.00	ng		0.00
34) Perylene-d12	23.39	264	83263	5.00	ng		-0.02
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	78113	15.10	ng		0.00
5) Phenol-d6	6.77	99	134511	13.50	ng		0.00
12) Nitrobenzene-d5	8.77	82	34315	7.72	ng		0.00
20) 2,4,6-Tribromophenol	15.77	330	17975	8.68	ng		0.00
23) 2-Fluorobiphenyl	12.87	172	156315	10.37	ng		0.00
32) Terphenyl-d14	19.65	244	118713	11.05	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	849	0.21	ng	#	63
3) n-Nitrosodimethylamine	3.46	42	305	0.13	ng	#	1
6) 2-Chlorophenol	7.17	128	1257	0.39	ng		85
7) Phenol	6.81	94	1334	0.17	ng		79
8) bis(2-Chloroethyl)ether	7.02	93	1088m	0.37	ng		
9) 2-Methylphenol	8.06	107	951	0.30	ng	#	88
10) 3+4-Methylphenols	8.38	107	932	0.25	ng	#	83
13) Nitrobenzene	8.80	77	1022	0.20	ng	#	73
14) 2-Nitrophenol	9.52	139	353	0.26	ng	#	37
15) 2,4-Dimethylphenol	9.58	122	701	0.13	ng		83
16) 2,4-Dichlorophenol	10.07	162	673	0.33	ng		85
17) Hexachlorobutadiene	10.72	225	670	0.18	ng	#	71
18) 4-Chloro-3-methylphenol	11.69	107	711	0.15	ng		93
21) 2,4,6-Trichlorophenol	12.70	196	299	0.20	ng	#	55
22) 2,4,5-Trichlorophenol	12.77	196	393	0.16	ng	#	79
25) 4-Nitrophenol	14.54	139	85	0.29	ng		87
27) 4,6-Dinitro-2-methylphenol	15.41	198	69	0.39	ng		92
28) Hexachlorobenzene	16.33	284	970	0.21	ng	#	83
29) Pentachlorophenol	16.69	266	31	0.31	ng	#	67
31) Benzidine	19.29	184	1134	0.58	ng	#	72
33) 3,3'-Dichlorobenzidine	21.13	252	906	0.30	ng	#	87

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

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