Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000841.D
Acq On : 03 Apr 2015 18:48
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001

Quant Time: Apr 04 07:18:45 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 03 18:16:53 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	22443	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	113138	5.00	ng		0.00
19) Acenaphthene-d10	14.26	164	58071	5.00	ng		-0.02
26) Phenanthrene-d10	17.00	188	132696	5.00	ng		0.00
30) Chrysene-d12	21.21	240	110496	5.00	ng		0.00
34) Perylene-d12	23.40	264	100150	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	7238	1.17	ng		0.00
5) Phenol-d6	6.77	99	10983	1.15	ng		0.00
12) Nitrobenzene-d5	8.77	82	5235	0.99	ng		0.00
20) 2,4,6-Tribromophenol	15.77	330	1812	0.90	ng		0.00
23) 2-Fluorobiphenyl	12.87	172	18322	1.02	ng		0.00
32) Terphenyl-d14	19.65	244	13741	1.03	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	3470	1.17	ng		95
3) n-Nitrosodimethylamine	3.43	42	3193	1.12	ng	#	32
6) 2-Chlorophenol	7.17	128	8123	1.23	ng		85
7) Phenol	6.81	94	9841	1.06	ng		96
8) bis(2-Chloroethyl)ether	7.02	93	7964	1.17	ng	#	76
9) 2-Methylphenol	8.06	107	7435	1.07	ng		96
10) 3+4-Methylphenols	8.38	107	8895	1.10	ng		95
13) Nitrobenzene	8.80	77	5806	0.94	ng	#	83
14) 2-Nitrophenol	9.52	139	3132	1.04	ng		98
15) 2,4-Dimethylphenol	9.58	122	6729	1.05	ng		95
16) 2,4-Dichlorophenol	10.04	162	6273	1.04	ng		95
17) Hexachlorobutadiene	10.72	225	4739	1.08	ng	#	67
18) 4-Chloro-3-methylphenol	11.69	107	6099	1.07	ng		99
21) 2,4,6-Trichlorophenol	12.69	196	3559	0.93	ng		93
22) 2,4,5-Trichlorophenol	12.75	196	4483	1.05	ng		95
25) 4-Nitrophenol	14.49	139	1027	0.93	ng		96
27) 4,6-Dinitro-2-methylphenol	15.39	198	1215	0.93	ng		93
28) Hexachlorobenzene	16.33	284	6471	0.85	ng		96
29) Pentachlorophenol	16.67	266	1040	0.84	ng		98
31) Benzidine	19.28	184	7233	1.79	ng	#	51
33) 3,3'-Dichlorobenzidine	21.13	252	8410	1.00	ng		98

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

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