Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000828.D
Acq On : 03 Apr 2015 04:08
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001

Quant Time: Apr 03 18:39:34 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	21784	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	105336	5.00	ng		0.00
19) Acenaphthene-d10	14.28	164	53478	5.00	ng		0.00
26) Phenanthrene-d10	17.01	188	136562	5.00	ng		0.00
30) Chrysene-d12	21.20	240	105218	5.00	ng		0.00
34) Perylene-d12	23.40	264	94983	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	6467	1.08	ng		0.00
5) Phenol-d6	6.77	99	9094	1.02	ng		0.00
12) Nitrobenzene-d5	8.77	82	4700	0.96	ng		0.00
20) 2,4,6-Tribromophenol	15.75	330	1542	0.85	ng		-0.02
23) 2-Fluorobiphenyl	12.87	172	16635	1.01	ng		0.00
32) Terphenyl-d14	19.66	244	12991	1.03	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	3157	1.08	ng		97
3) n-Nitrosodimethylamine	3.44	42	2905	1.05	ng		93
6) 2-Chlorophenol	7.17	128	6438	1.04	ng		96
7) Phenol	6.81	94	9196	1.02	ng		98
8) bis(2-Chloroethyl)ether	7.02	93	6515	1.02	ng		92
9) 2-Methylphenol	8.06	107	6697	1.01	ng		98
10) 3+4-Methylphenols	8.38	107	7633	0.99	ng		99
13) Nitrobenzene	8.80	77	5404	0.94	ng		97
14) 2-Nitrophenol	9.52	139	2726	0.98	ng		99
15) 2,4-Dimethylphenol	9.58	122	5979	1.00	ng		99
16) 2,4-Dichlorophenol	10.04	162	5268	0.96	ng		99
17) Hexachlorobutadiene	10.72	225	4219	1.04	ng	#	67
18) 4-Chloro-3-methylphenol	11.69	107	5301	1.00	ng		100
21) 2,4,6-Trichlorophenol	12.69	196	3062	0.88	ng		99
22) 2,4,5-Trichlorophenol	12.75	196	3935	1.01	ng		99
24) 2,4-Dinitrophenol	14.41	184	135	0.26	ng		89
25) 4-Nitrophenol	14.51	139	880	0.88	ng		91
27) 4,6-Dinitro-2-methylphenol	15.40	198	1002	0.82	ng	#	44
28) Hexachlorobenzene	16.32	284	5543	0.72	ng	#	1
29) Pentachlorophenol	16.67	266	813	0.72	ng		90
31) Benzidine	19.27	184	3675	1.20	ng	#	54
33) 3,3'-Dichlorobenzidine	21.14	252	7004	0.90	ng	#	76

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

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