Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000813.D
Acq On : 02 Apr 2015 19:09
Operator : TP/IZ
Sample : SSTDICC0.5
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC0.5

Quant Time: Apr 03 18:14:19 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 17:58:44 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	24580	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	116308	5.00	ng		0.00
19) Acenaphthene-d10	14.28	164	56297	5.00	ng		0.00
26) Phenanthrene-d10	17.00	188	114194	5.00	ng		0.00
30) Chrysene-d12	21.21	240	104545	5.00	ng		0.00
34) Perylene-d12	23.40	264	95716	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	3365	0.50	ng		0.00
5) Phenol-d6	6.77	99	4333	0.46	ng		0.00
12) Nitrobenzene-d5	8.77	82	2783	0.51	ng		0.00
20) 2,4,6-Tribromophenol	15.77	330	750	0.45	ng		0.00
23) 2-Fluorobiphenyl	12.87	172	8878	0.51	ng		0.00
32) Terphenyl-d14	19.65	244	6445	0.51	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	1972	0.53	ng		88
3) n-Nitrosodimethylamine	3.44	42	1421	0.45	ng	#	57
6) 2-Chlorophenol	7.17	128	3153	0.47	ng		88
7) Phenol	6.81	94	5087	0.50	ng		93
8) bis(2-Chloroethyl)ether	7.02	93	3226m	0.46	ng		
9) 2-Methylphenol	8.06	107	3578	0.49	ng	#	95
10) 3+4-Methylphenols	8.38	107	3747	0.49	ng		93
13) Nitrobenzene	8.80	77	3168	0.50	ng	#	84
14) 2-Nitrophenol	9.52	139	1629	0.56	ng	#	82
15) 2,4-Dimethylphenol	9.58	122	3132	0.48	ng		93
16) 2,4-Dichlorophenol	10.04	162	2572	0.44	ng		87
17) Hexachlorobutadiene	10.75	225	2299	0.51	ng		98
18) 4-Chloro-3-methylphenol	11.69	107	2641	0.45	ng		97
21) 2,4,6-Trichlorophenol	12.69	196	1384	0.46	ng		97
22) 2,4,5-Trichlorophenol	12.77	196	1647m	0.44	ng		
25) 4-Nitrophenol	14.51	139	333	0.30	ng		94
27) 4,6-Dinitro-2-methylphenol	15.41	198	354	0.35	ng	#	54
28) Hexachlorobenzene	16.33	284	3363	0.52	ng		96
29) Pentachlorophenol	16.67	266	233	0.21	ng	#	84
31) Benzidine	19.29	184	1736	0.41	ng	#	67
33) 3,3'-Dichlorobenzidine	21.13	252	3367	0.48	ng	#	96

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

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