Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040715\
Data File : BM000887.D
Acq On : 07 Apr 2015 19:25
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001EC

Quant Time: Apr 08 01:55:52 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 07 15:59:39 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	16742	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	86176	5.00	ng		0.00
19) Acenaphthene-d10	14.26	164	47392	5.00	ng		0.00
26) Phenanthrene-d10	17.00	188	115632	5.00	ng		0.00
30) Chrysene-d12	21.21	240	85472	5.00	ng		0.00
34) Perylene-d12	23.39	264	77088	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	5627	1.00	ng		0.00
5) Phenol-d6	6.77	99	8916	0.96	ng		0.00
12) Nitrobenzene-d5	8.73	82	4034	0.94	ng		-0.04
20) 2,4,6-Tribromophenol	15.77	330	1192	0.87	ng		0.02
23) 2-Fluorobiphenyl	12.87	172	14625	1.00	ng		0.00
32) Terphenyl-d14	19.65	244	10562	1.00	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	2791	1.05	ng		98
3) n-Nitrosodimethylamine	3.43	42	2524	0.97	ng	#	48
6) 2-Chlorophenol	7.17	128	7105	0.96	ng		95
7) Phenol	6.81	94	6221	0.89	ng		98
8) bis(2-Chloroethyl)ether	7.02	93	7139	0.97	ng		84
9) 2-Methylphenol	8.06	107	4887	0.95	ng		96
10) 3+4-Methylphenols	8.38	107	6533	0.94	ng		97
13) Nitrobenzene	8.80	77	3950	0.91	ng		86
14) 2-Nitrophenol	9.52	139	2415	0.99	ng		97
15) 2,4-Dimethylphenol	9.58	122	4786	0.97	ng		98
16) 2,4-Dichlorophenol	10.04	162	4724	0.93	ng		96
17) Hexachlorobutadiene	10.72	225	4007	1.02	ng		99
18) 4-Chloro-3-methylphenol	11.67	107	4548	0.95	ng		99
21) 2,4,6-Trichlorophenol	12.69	196	2518	0.87	ng		97
22) 2,4,5-Trichlorophenol	12.75	196	3302	0.90	ng		94
25) 4-Nitrophenol	14.49	139	660	0.93	ng		91
27) 4,6-Dinitro-2-methylphenol	15.39	198	801	0.78	ng	#	48
28) Hexachlorobenzene	16.33	284	4549	0.88	ng	#	53
29) Pentachlorophenol	16.67	266	580	0.91	ng		94
31) Benzidine	19.28	184	3912	1.00	ng		98
33) 3,3'-Dichlorobenzidine	21.13	252	5943	1.01	ng	#	85

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

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