Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000815.D
Acq On : 02 Apr 2015 20:21
Operator : TP/IZ
Sample : SSTDICCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICCC001

Quant Time: Apr 03 18:15:39 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	23726	5.00	ng	0.00
11) Naphthalene-d8	10.39	136	107784	5.00	ng	0.00
19) Acenaphthene-d10	14.28	164	53068	5.00	ng	0.00
26) Phenanthrene-d10	17.00	188	112695	5.00	ng	0.00
30) Chrysene-d12	21.21	240	103677	5.00	ng	0.00
34) Perylene-d12	23.40	264	93220	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.21	112	6583	1.01	ng	0.00
5) Phenol-d6	6.77	99	9220	1.01	ng	0.00
12) Nitrobenzene-d5	8.77	82	4840	0.96	ng	0.00
20) 2,4,6-Tribromophenol	15.77	330	1741	1.10	ng	0.00
23) 2-Fluorobiphenyl	12.87	172	16716	1.02	ng	0.00
32) Terphenyl-d14	19.65	244	12775	1.02	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.12	88	3293	0.92	ng	100
3) n-Nitrosodimethylamine	3.44	42	3022	1.00	ng	100
6) 2-Chlorophenol	7.17	128	6464	1.00	ng	100
7) Phenol	6.81	94	9587	0.98	ng	100
8) bis(2-Chloroethyl)ether	7.02	93	6441	0.95	ng	100
9) 2-Methylphenol	8.06	107	7022	1.01	ng	100
10) 3+4-Methylphenols	8.38	107	8029	1.09	ng	100
13) Nitrobenzene	8.80	77	5729	0.98	ng	100
14) 2-Nitrophenol	9.52	139	2874	1.07	ng	100
15) 2,4-Dimethylphenol	9.58	122	6250	1.02	ng	100
16) 2,4-Dichlorophenol	10.04	162	5540	1.02	ng	100
17) Hexachlorobutadiene	10.75	225	4260	1.02	ng	100
18) 4-Chloro-3-methylphenol	11.69	107	5360	0.99	ng	100
21) 2,4,6-Trichlorophenol	12.69	196	3460	1.21	ng	100
22) 2,4,5-Trichlorophenol	12.75	196	4179	1.19	ng	100
24) 2,4-Dinitrophenol	14.40	184	240	0.45	ng	100
25) 4-Nitrophenol	14.49	139	1009	0.97	ng	100
27) 4,6-Dinitro-2-methylphenol	15.39	198	1043	1.04	ng	100
28) Hexachlorobenzene	16.33	284	6485	1.02	ng	100
29) Pentachlorophenol	16.67	266	1178	1.07	ng	100
31) Benzidine	19.30	184	528	0.13	ng	100
33) 3,3'-Dichlorobenzidine	21.13	252	5845	0.84	ng	100

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

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