Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000819.D
Acq On : 02 Apr 2015 22:44
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040315

Quant Time: Apr 03 18:28:18 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	27611	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	127534	5.00	ng		0.00
19) Acenaphthene-d10	14.28	164	62543	5.00	ng		0.00
26) Phenanthrene-d10	17.00	188	124983	5.00	ng		0.00
30) Chrysene-d12	21.21	240	120032	5.00	ng		0.00
34) Perylene-d12	23.40	264	107252	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	7924	1.04	ng		0.00
5) Phenol-d6	6.77	99	11263	1.00	ng		0.00
12) Nitrobenzene-d5	8.77	82	5912	0.99	ng		0.00
20) 2,4,6-Tribromophenol	15.77	330	2071	0.95	ng		0.00
23) 2-Fluorobiphenyl	12.87	172	20091	1.04	ng		0.00
32) Terphenyl-d14	19.65	244	15045	1.04	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	3670	0.98	ng		94
3) n-Nitrosodimethylamine	3.43	42	3666	1.04	ng	#	33
6) 2-Chlorophenol	7.17	128	7843	1.00	ng		98
7) Phenol	6.81	94	11776	1.03	ng		97
8) bis(2-Chloroethyl)ether	7.02	93	7842	0.98	ng		100
9) 2-Methylphenol	8.06	107	8625	1.02	ng		99
10) 3+4-Methylphenols	8.38	107	9470	0.97	ng		99
13) Nitrobenzene	8.80	77	7221	1.04	ng		99
14) 2-Nitrophenol	9.52	139	3494	1.03	ng		98
15) 2,4-Dimethylphenol	9.58	122	7607	1.05	ng		100
16) 2,4-Dichlorophenol	10.04	162	6701	1.00	ng		99
17) Hexachlorobutadiene	10.75	225	5043	1.02	ng		100
18) 4-Chloro-3-methylphenol	11.69	107	6435	1.01	ng		98
21) 2,4,6-Trichlorophenol	12.69	196	4030	0.97	ng		100
22) 2,4,5-Trichlorophenol	12.75	196	4965	1.07	ng		99
24) 2,4-Dinitrophenol	14.40	184	277	0.45	ng		99
25) 4-Nitrophenol	14.49	139	1299	1.04	ng		98
27) 4,6-Dinitro-2-methylphenol	15.41	198	1063	0.89	ng	#	30
28) Hexachlorobenzene	16.33	284	7600	1.05	ng		100
29) Pentachlorophenol	16.67	266	1342	1.03	ng		97
31) Benzidine	19.28	184	4102	1.18	ng	#	53
33) 3,3'-Dichlorobenzidine	21.13	252	8656	0.96	ng		97

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

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