Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000816.D
Acq On : 02 Apr 2015 20:57
Operator : TP/IZ
Sample : SSTDICC002
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC002

Quant Time: Apr 03 18:16:15 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	27053	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	123709	5.00	ng		0.00
19) Acenaphthene-d10	14.28	164	60881	5.00	ng		0.00
26) Phenanthrene-d10	17.00	188	123846	5.00	ng		0.00
30) Chrysene-d12	21.21	240	116391	5.00	ng		0.00
34) Perylene-d12	23.40	264	104671	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	15502	2.08	ng		0.00
5) Phenol-d6	6.77	99	23587	2.27	ng		0.00
12) Nitrobenzene-d5	8.77	82	11485	1.99	ng		0.00
20) 2,4,6-Tribromophenol	15.77	330	4486	2.47	ng		0.00
23) 2-Fluorobiphenyl	12.87	172	39382	2.10	ng		0.00
32) Terphenyl-d14	19.65	244	29327	2.09	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	6683	1.63	ng		91
3) n-Nitrosodimethylamine	3.43	42	7369	2.14	ng	#	38
6) 2-Chlorophenol	7.17	128	16141	2.18	ng		93
7) Phenol	6.81	94	23410	2.10	ng		93
8) bis(2-Chloroethyl)ether	7.02	93	16331	2.10	ng		89
9) 2-Methylphenol	8.06	107	16972	2.13	ng		95
10) 3+4-Methylphenols	8.38	107	19705	2.34	ng		95
13) Nitrobenzene	8.80	77	14106	2.09	ng		90
14) 2-Nitrophenol	9.52	139	6589	2.14	ng	#	90
15) 2,4-Dimethylphenol	9.58	122	14886	2.12	ng		97
16) 2,4-Dichlorophenol	10.04	162	13702	2.21	ng		95
17) Hexachlorobutadiene	10.75	225	9899	2.07	ng		100
18) 4-Chloro-3-methylphenol	11.69	107	13210	2.13	ng		98
21) 2,4,6-Trichlorophenol	12.69	196	8912	2.71	ng		99
22) 2,4,5-Trichlorophenol	12.75	196	10641	2.64	ng		95
24) 2,4-Dinitrophenol	14.40	184	904	1.49	ng	#	64
25) 4-Nitrophenol	14.49	139	3191	2.67	ng		87
27) 4,6-Dinitro-2-methylphenol	15.39	198	2628	2.38	ng	#	80
28) Hexachlorobenzene	16.33	284	15153	2.17	ng		97
29) Pentachlorophenol	16.67	266	3553	2.94	ng		95
31) Benzidine	19.28	184	10468	2.23	ng	#	50
33) 3,3'-Dichlorobenzidine	21.13	252	18760	2.40	ng		99

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

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