Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000814.D
Acq On : 02 Apr 2015 19:45
Operator : TP/IZ
Sample : SSTDICC0.8
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC0.8

Quant Time: Apr 03 18:15:21 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	26258	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	122374	5.00	ng		0.00
19) Acenaphthene-d10	14.28	164	62184	5.00	ng		0.00
26) Phenanthrene-d10	17.00	188	127471	5.00	ng		0.00
30) Chrysene-d12	21.21	240	118883	5.00	ng		0.00
34) Perylene-d12	23.40	264	107192	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	5704	0.79	ng		0.00
5) Phenol-d6	6.77	99	7965	0.79	ng		0.00
12) Nitrobenzene-d5	8.77	82	4367	0.76	ng		0.00
20) 2,4,6-Tribromophenol	15.77	330	1428	0.77	ng		0.00
23) 2-Fluorobiphenyl	12.87	172	15044	0.79	ng		0.00
32) Terphenyl-d14	19.65	244	11263	0.79	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	3026	0.76	ng		95
3) n-Nitrosodimethylamine	3.44	42	2545	0.76	ng		91
6) 2-Chlorophenol	7.17	128	5459	0.76	ng		95
7) Phenol	6.81	94	8669	0.80	ng		99
8) bis(2-Chloroethyl)ether	7.02	93	5889m	0.78	ng		
9) 2-Methylphenol	8.06	107	6325	0.82	ng		99
10) 3+4-Methylphenols	8.38	107	6823	0.83	ng		98
13) Nitrobenzene	8.80	77	5199	0.78	ng		95
14) 2-Nitrophenol	9.52	139	2646	0.87	ng		96
15) 2,4-Dimethylphenol	9.58	122	5505	0.79	ng		97
16) 2,4-Dichlorophenol	10.04	162	4605	0.75	ng		94
17) Hexachlorobutadiene	10.75	225	3850	0.81	ng		100
18) 4-Chloro-3-methylphenol	11.69	107	4761	0.78	ng		99
21) 2,4,6-Trichlorophenol	12.69	196	2811	0.84	ng		99
22) 2,4,5-Trichlorophenol	12.77	196	3210m	0.78	ng		
24) 2,4-Dinitrophenol	14.42	184	119	0.19	ng		90
25) 4-Nitrophenol	14.49	139	777	0.64	ng		88
27) 4,6-Dinitro-2-methylphenol	15.41	198	898	0.79	ng	#	40
28) Hexachlorobenzene	16.33	284	5868	0.82	ng		98
29) Pentachlorophenol	16.67	266	693	0.56	ng		97
31) Benzidine	19.28	184	2317	0.48	ng	#	60
33) 3,3'-Dichlorobenzidine	21.13	252	5814	0.73	ng		96

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

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