Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040715\
Data File : BM000878.D
Acq On : 07 Apr 2015 14:02
Operator : TP/IZ
Sample : SSTDICC002
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC002

Quant Time: Apr 07 15:08:38 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 14:20:59 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	23013	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	120356	5.00	ng		0.00
19) Acenaphthene-d10	14.26	164	64301	5.00	ng		0.00
26) Phenanthrene-d10	17.00	188	151640	5.00	ng		0.00
30) Chrysene-d12	21.21	240	114979	5.00	ng		0.00
34) Perylene-d12	23.39	264	106638	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	16437	2.24	ng		0.00
5) Phenol-d6	6.77	99	27636	2.60	ng		0.00
12) Nitrobenzene-d5	8.73	82	12590	2.20	ng		-0.04
20) 2,4,6-Tribromophenol	15.77	330	3981	2.57	ng		0.02
23) 2-Fluorobiphenyl	12.87	172	42009	2.14	ng		0.00
32) Terphenyl-d14	19.65	244	29627	2.15	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	6957	1.44	ng	#	85
3) n-Nitrosodimethylamine	3.43	42	7706	2.56	ng	#	53
6) 2-Chlorophenol	7.17	128	22319	2.67	ng		93
7) Phenol	6.81	94	19465	2.07	ng		90
8) bis(2-Chloroethyl)ether	7.02	93	22608	2.85	ng	#	76
9) 2-Methylphenol	8.06	107	14389	2.10	ng	#	92
10) 3+4-Methylphenols	8.38	107	20882	2.62	ng		92
13) Nitrobenzene	8.80	77	12349	2.11	ng	#	64
14) 2-Nitrophenol	9.52	139	6763	1.72	ng	#	83
15) 2,4-Dimethylphenol	9.58	122	14501	2.25	ng		92
16) 2,4-Dichlorophenol	10.04	162	15391	2.67	ng		91
17) Hexachlorobutadiene	10.72	225	11283	2.10	ng		99
18) 4-Chloro-3-methylphenol	11.67	107	14000	2.37	ng		98
21) 2,4,6-Trichlorophenol	12.69	196	8799	3.27	ng		95
22) 2,4,5-Trichlorophenol	12.75	196	10874	2.99	ng		87
25) 4-Nitrophenol	14.49	139	2886	4.43	ng		86
27) 4,6-Dinitro-2-methylphenol	15.39	198	3714	3.63	ng		91
28) Hexachlorobenzene	16.33	284	12594	1.81	ng	#	50
29) Pentachlorophenol	16.67	266	3082	6.02	ng		90
31) Benzidine	19.26	184	13530	2.84	ng	#	91
33) 3,3'-Dichlorobenzidine	21.13	252	19685	2.75	ng	#	82

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

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