Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040715\
Data File : BM000881.D
Acq On : 07 Apr 2015 15:50
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040715

Quant Time: Apr 07 16:37:48 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 15:59:39 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	22421	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	115612	5.00	ng		0.00
19) Acenaphthene-d10	14.26	164	61914	5.00	ng		0.00
26) Phenanthrene-d10	17.00	188	152641	5.00	ng		0.00
30) Chrysene-d12	21.21	240	117777	5.00	ng		0.00
34) Perylene-d12	23.39	264	108175	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	7941	1.06	ng		0.00
5) Phenol-d6	6.77	99	12966	1.03	ng		0.00
12) Nitrobenzene-d5	8.73	82	5832	1.01	ng		-0.04
20) 2,4,6-Tribromophenol	15.77	330	1804	0.97	ng		0.02
23) 2-Fluorobiphenyl	12.87	172	20118	1.05	ng		0.00
32) Terphenyl-d14	19.65	244	15146	1.04	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	3717	1.04	ng		96
3) n-Nitrosodimethylamine	3.43	42	3650	1.03	ng	#	36
6) 2-Chlorophenol	7.17	128	9968	1.00	ng		96
7) Phenol	6.81	94	9742	1.04	ng		95
8) bis(2-Chloroethyl)ether	7.02	93	9991	1.00	ng		89
9) 2-Methylphenol	8.06	107	7204	1.05	ng		97
10) 3+4-Methylphenols	8.38	107	9606	1.01	ng		96
13) Nitrobenzene	8.80	77	5960	1.03	ng		85
14) 2-Nitrophenol	9.52	139	3312	1.01	ng	#	91
15) 2,4-Dimethylphenol	9.58	122	6873	1.04	ng		96
16) 2,4-Dichlorophenol	10.04	162	6927	1.00	ng		95
17) Hexachlorobutadiene	10.72	225	5469	1.04	ng		99
18) 4-Chloro-3-methylphenol	11.67	107	6485	1.00	ng		99
21) 2,4,6-Trichlorophenol	12.69	196	3740	0.96	ng		97
22) 2,4,5-Trichlorophenol	12.75	196	4789	0.98	ng		93
25) 4-Nitrophenol	14.49	139	1206	1.17	ng		96
27) 4,6-Dinitro-2-methylphenol	15.39	198	1450	0.99	ng	#	72
28) Hexachlorobenzene	16.33	284	6387	0.94	ng	#	52
29) Pentachlorophenol	16.67	266	1134	1.09	ng		91
31) Benzidine	19.28	184	6849	1.20	ng	#	95
33) 3,3'-Dichlorobenzidine	21.13	252	9646	1.16	ng	#	82

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

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