

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050615\
 Data File : BM001237.D
 Acq On : 07 May 2015 03:50
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001EC

Quant Time: May 07 07:29:28 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM050615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 07 07:00:24 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	14940	5.00	ng	0.00
9) Naphthalene-d8	10.26	136	44680	5.00	ng	0.00
16) Acenaphthene-d10	14.15	164	24304	5.00	ng	0.00
23) Phenanthrene-d10	16.90	188	44187	5.00	ng	0.00
27) Chrysene-d12	21.09	240	49848	5.00	ng	0.00
29) Perylene-d12	23.23	264	46081	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	5.06	112	2527	1.04	ng	0.00
4) Phenol-d6	6.65	99	3529	1.02	ng	0.00
10) Nitrobenzene-d5	8.61	82	2709	1.01	ng	0.00
17) 2,4,6-Tribromophenol	15.64	330	960	0.92	ng	0.00
20) 2-Fluorobiphenyl	12.76	172	7786	1.03	ng	0.00
28) Terphenyl-d14	19.54	244	6753	1.03	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.94	88	993	1.02	ng	100
5) 2-Chlorophenol	7.05	128	2166	1.02	ng	99
6) Phenol	6.68	94	3232	0.98	ng	97
7) 2-Methylphenol	7.94	107	1742	1.00	ng	99
8) 3+4-Methylphenols	8.24	107	2448	1.01	ng	# 39
11) 2-Nitrophenol	9.39	139	981	0.95	ng	98
12) 2,4-Dimethylphenol	9.46	122	2160	1.00	ng	99
13) 2,4-Dichlorophenol	9.92	162	2184	1.00	ng	99
14) 4-Chloro-3-methylphenol	11.55	107	2255	0.96	ng	92
15) 2-Methylnaphthalene	11.92	142	6712	1.05	ng	99
18) 2,4,6-Trichlorophenol	12.56	196	1734	0.95	ng	98
19) 2,4,5-Trichlorophenol	12.64	196	1932	0.92	ng	98
21) 2,4-Dinitrophenol	14.29	184	250	0.91	ng	97
22) 4-Nitrophenol	14.40	139	465	1.04	ng	98
24) 4,6-Dinitro-2-methylphenol	15.29	198	641	0.94	ng	# 77
25) Hexachlorobenzene	16.21	284	3381	1.04	ng	98
26) Pentachlorophenol	16.56	266	632	1.01	ng	98

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

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