

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050615\
 Data File : BM001232.D
 Acq On : 07 May 2015 00:16
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.1

Quant Time: May 07 07:16:02 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 06:25:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	14980	5.00	ng	0.00
9) Naphthalene-d8	10.26	136	42930	5.00	ng	0.00
16) Acenaphthene-d10	14.15	164	23021	5.00	ng	0.00
23) Phenanthrene-d10	16.90	188	43394	5.00	ng	0.00
27) Chrysene-d12	21.09	240	47338	5.00	ng	0.00
29) Perylene-d12	23.23	264	44422	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	5.06	112	235	0.10	ng	0.00
4) Phenol-d6	6.65	99	259	0.08	ng	0.00
10) Nitrobenzene-d5	8.64	82	160	0.07	ng	0.03
17) 2,4,6-Tribromophenol	15.64	330	71	0.07	ng	0.00
20) 2-Fluorobiphenyl	12.76	172	706	0.10	ng	0.00
28) Terphenyl-d14	19.55	244	598	0.00	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.95	88	114	0.12	ng	87
5) 2-Chlorophenol	7.05	128	213	0.10	ng	# 60
6) Phenol	6.68	94	310	0.09	ng	# 49
7) 2-Methylphenol	7.94	107	150	0.09	ng	# 63
8) 3+4-Methylphenols	8.27	107	163	0.07	ng	# 59
11) 2-Nitrophenol	9.39	139	69	0.07	ng	# 1
12) 2,4-Dimethylphenol	9.46	122	182	0.09	ng	# 75
13) 2,4-Dichlorophenol	9.95	162	169	0.08	ng	# 63
14) 4-Chloro-3-methylphenol	11.57	107	181	0.08	ng	# 55
15) 2-Methylnaphthalene	11.92	142	566	0.09	ng	98
18) 2,4,6-Trichlorophenol	12.58	196	133	0.08	ng	# 47
19) 2,4,5-Trichlorophenol	12.66	196	138	0.07	ng	87
25) Hexachlorobenzene	16.21	284	295	0.09	ng	97

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

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