

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050615\
 Data File : BM001229.D
 Acq On : 06 May 2015 22:29
 Operator : TP/IZ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.8

Quant Time: May 07 06:39:59 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	16521	5.00	ng	0.00
9) Naphthalene-d8	10.26	136	48792	5.00	ng	0.00
16) Acenaphthene-d10	14.15	164	25964	5.00	ng	0.00
23) Phenanthrene-d10	16.90	188	45943	5.00	ng	0.00
27) Chrysene-d12	21.09	240	54535	5.00	ng	0.00
29) Perylene-d12	23.23	264	50466	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	5.06	112	2099	0.78	ng	0.00
4) Phenol-d6	6.65	99	2742	0.75	ng	0.00
10) Nitrobenzene-d5	8.61	82	2105	0.76	ng	0.00
17) 2,4,6-Tribromophenol	15.64	330	766	0.71	ng	0.00
20) 2-Fluorobiphenyl	12.76	172	6344	0.78	ng	0.00
28) Terphenyl-d14	19.55	244	5384	0.00	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.94	88	819	0.76	ng	99
5) 2-Chlorophenol	7.05	128	1802	0.77	ng	98
6) Phenol	6.68	94	2733	0.75	ng	98
7) 2-Methylphenol	7.94	107	1494	0.77	ng	98
8) 3+4-Methylphenols	8.27	107	1907	0.76	ng	91
11) 2-Nitrophenol	9.39	139	809	0.74	ng	94
12) 2,4-Dimethylphenol	9.46	122	1728	0.73	ng	98
13) 2,4-Dichlorophenol	9.92	162	1811	0.76	ng	98
14) 4-Chloro-3-methylphenol	11.55	107	1788	0.72	ng	98
15) 2-Methylnaphthalene	11.92	142	5378	0.77	ng	100
18) 2,4,6-Trichlorophenol	12.56	196	1383	0.73	ng	98
19) 2,4,5-Trichlorophenol	12.64	196	1604	0.75	ng	97
21) 2,4-Dinitrophenol	14.29	184	204	0.60	ng	# 83
22) 4-Nitrophenol	14.40	139	363	0.33	ng	96
24) 4,6-Dinitro-2-methylphenol	15.29	198	511	0.64	ng	96
25) Hexachlorobenzene	16.21	284	2686	0.79	ng	99
26) Pentachlorophenol	16.56	266	503	0.67	ng	96

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

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