

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050615\
 Data File : BM001228.D
 Acq On : 06 May 2015 21:53
 Operator : TP/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICCC001

Quant Time: May 07 06:39:49 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	15885	5.00	ng	0.00
9) Naphthalene-d8	10.26	136	46857	5.00	ng	0.00
16) Acenaphthene-d10	14.15	164	24831	5.00	ng	0.00
23) Phenanthrene-d10	16.90	188	44214	5.00	ng	0.00
27) Chrysene-d12	21.09	240	50117	5.00	ng	0.00
29) Perylene-d12	23.23	264	45801	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	5.06	112	2621	1.01	ng	0.00
4) Phenol-d6	6.65	99	3602	1.03	ng	0.00
10) Nitrobenzene-d5	8.61	82	2687	1.01	ng	0.00
17) 2,4,6-Tribromophenol	15.64	330	1022	0.99	ng	0.00
20) 2-Fluorobiphenyl	12.76	172	7968	1.03	ng	0.00
28) Terphenyl-d14	19.55	244	6769	0.01	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.94	88	1020	0.99	ng	100
5) 2-Chlorophenol	7.05	128	2248	0.99	ng	100
6) Phenol	6.68	94	3569	1.01	ng	100
7) 2-Methylphenol	7.94	107	1879	1.01	ng	100
8) 3+4-Methylphenols	8.27	107	2463	1.02	ng	100
11) 2-Nitrophenol	9.39	139	1037	0.99	ng	100
12) 2,4-Dimethylphenol	9.46	122	2281	1.01	ng	100
13) 2,4-Dichlorophenol	9.92	162	2335	1.01	ng	100
14) 4-Chloro-3-methylphenol	11.55	107	2390	1.00	ng	100
15) 2-Methylnaphthalene	11.92	142	6902	1.03	ng	100
18) 2,4,6-Trichlorophenol	12.56	196	1822	1.00	ng	100
19) 2,4,5-Trichlorophenol	12.64	196	2070	1.01	ng	100
21) 2,4-Dinitrophenol	14.29	184	272	0.83	ng	100
22) 4-Nitrophenol	14.40	139	509	0.48	ng	100
24) 4,6-Dinitro-2-methylphenol	15.29	198	683	0.88	ng	100
25) Hexachlorobenzene	16.21	284	3420	1.05	ng	100
26) Pentachlorophenol	16.56	266	692	0.96	ng	100

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

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