

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050615\
 Data File : BM001227.D
 Acq On : 06 May 2015 21:17
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC002

Quant Time: May 07 06:39:40 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	18625	5.00	ng	0.00
9) Naphthalene-d8	10.26	136	53711	5.00	ng	0.00
16) Acenaphthene-d10	14.15	164	28827	5.00	ng	0.00
23) Phenanthrene-d10	16.90	188	49851	5.00	ng	0.00
27) Chrysene-d12	21.09	240	58948	5.00	ng	0.00
29) Perylene-d12	23.23	264	54009	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	5.06	112	6039	1.99	ng	0.00
4) Phenol-d6	6.65	99	8737	2.12	ng	0.00
10) Nitrobenzene-d5	8.61	82	6573	2.15	ng	0.00
17) 2,4,6-Tribromophenol	15.64	330	2584	2.15	ng	0.00
20) 2-Fluorobiphenyl	12.76	172	17887	1.99	ng	0.00
28) Terphenyl-d14	19.55	244	15770	0.01	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.94	88	2356	1.95	ng	98
5) 2-Chlorophenol	7.05	128	5235	1.98	ng	97
6) Phenol	6.68	94	8291	2.01	ng	93
7) 2-Methylphenol	7.94	107	4357	2.00	ng	96
8) 3+4-Methylphenols	8.24	107	6130	2.16	ng	# 39
11) 2-Nitrophenol	9.39	139	2585	2.15	ng	# 84
12) 2,4-Dimethylphenol	9.46	122	5498	2.12	ng	94
13) 2,4-Dichlorophenol	9.92	162	5647	2.14	ng	96
14) 4-Chloro-3-methylphenol	11.55	107	5813	2.12	ng	94
15) 2-Methylnaphthalene	11.92	142	15787	2.06	ng	100
18) 2,4,6-Trichlorophenol	12.56	196	4466	2.11	ng	99
19) 2,4,5-Trichlorophenol	12.64	196	5214	2.19	ng	96
21) 2,4-Dinitrophenol	14.29	184	950	2.50	ng	# 70
22) 4-Nitrophenol	14.38	139	1749	1.43	ng	91
24) 4,6-Dinitro-2-methylphenol	15.29	198	1990	2.29	ng	# 72
25) Hexachlorobenzene	16.21	284	7622	2.08	ng	99
26) Pentachlorophenol	16.56	266	1870	2.29	ng	96

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

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