

Data Path : Z:\HPCHEM1\BNA M\DATA\BM050615\
 Data File : BM001233.D
 Acq On : 07 May 2015 00:52
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM050615

Quant Time: May 07 07:18:55 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	19960	5.00	ng	0.00
9) Naphthalene-d8	10.26	136	60832	5.00	ng	0.00
16) Acenaphthene-d10	14.15	164	32465	5.00	ng	0.00
23) Phenanthrene-d10	16.90	188	59230	5.00	ng	0.00
27) Chrysene-d12	21.09	240	68246	5.00	ng	0.00
29) Perylene-d12	23.23	264	63425	5.00	ng	0.00

System Monitoring Compounds						
3) 2-Fluorophenol	5.06	112	3416	1.05	ng	0.00
4) Phenol-d6	6.65	99	4755	1.02	ng	0.00
10) Nitrobenzene-d5	8.61	82	3659	1.00	ng	0.00
17) 2,4,6-Tribromophenol	15.64	330	1335	0.95	ng	0.00
20) 2-Fluorobiphenyl	12.76	172	10439	1.03	ng	0.00
28) Terphenyl-d14	19.55	244	9019	1.00	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.94	88	1326	1.02	ng	99
5) 2-Chlorophenol	7.05	128	2888	1.02	ng	99
6) Phenol	6.68	94	4387	0.99	ng	94
7) 2-Methylphenol	7.94	107	2342	1.00	ng	97
8) 3+4-Methylphenols	8.24	107	3353	1.03	ng	# 40
11) 2-Nitrophenol	9.39	139	1344	0.95	ng	94
12) 2,4-Dimethylphenol	9.46	122	2952	1.01	ng	97
13) 2,4-Dichlorophenol	9.92	162	3011	1.01	ng	97
14) 4-Chloro-3-methylphenol	11.55	107	3075	0.96	ng	91
15) 2-Methylnaphthalene	11.92	142	8947	1.03	ng	100
18) 2,4,6-Trichlorophenol	12.56	196	2383	0.97	ng	96
19) 2,4,5-Trichlorophenol	12.64	196	2712	0.96	ng	93
21) 2,4-Dinitrophenol	14.29	184	363	0.95	ng	91
22) 4-Nitrophenol	14.38	139	737	1.09	ng	# 85
24) 4,6-Dinitro-2-methylphenol	15.29	198	894	0.96	ng	# 72
25) Hexachlorobenzene	16.21	284	4503	1.03	ng	98
26) Pentachlorophenol	16.56	266	870	1.02	ng	98

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

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