

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101315\
 Data File : BM002895.D
 Acq On : 13 Oct 2015 20:23
 Operator : UM/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICCC001

Quant Time: Oct 14 17:26:20 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM0101315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 14 17:20:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.66	152	91833	5.00	ng	0.00
8) Naphthalene-d8	11.50	136	385860	5.00	ng	0.00
14) Acenaphthene-d10	15.25	164	203740	5.00	ng	0.00
21) Phenanthrene-d10	17.95	188	455898	5.00	ng	0.00
24) Chrysene-d12	22.04	240	427990	5.00	ng	0.00
26) Perylene-d12	24.61	264	399004	5.00	ng	0.00

System Monitoring Compounds						
2) 2-Fluorophenol	6.13	112	16785	1.02	ng	0.00
3) Phenol-d6	7.79	99	23851	0.95	ng	0.00
9) Nitrobenzene-d5	9.86	82	21243	0.88	ng	0.00
15) 2,4,6-Tribromophenol	16.71	330	7139	0.96	ng	0.00
18) 2-Fluorobiphenyl	13.87	172	66155	0.99	ng	0.00
25) Terphenyl-d14	20.48	244	63148	0.97	ng	0.00

Target Compounds						Qvalue
4) 2-Chlorophenol	8.21	128	22990	0.98	ng	# 72
5) Phenol	7.82	94	24337	0.91	ng	# 87
6) 2-Methylphenol	9.11	107	18234	0.96	ng	# 89
7) 3+4-Methylphenols	9.45	107	22975	0.98	ng	# 35
10) 2-Nitrophenol	10.64	139	9317	0.96	ng	# 67
11) 2,4-Dimethylphenol	10.66	122	20787	0.98	ng	# 96
12) 2,4-Dichlorophenol	11.19	162	17096	1.03	ng	# 89
13) 4-Chloro-3-methylphenol	12.76	107	19949	0.95	ng	# 72
16) 2,4,6-Trichlorophenol	13.72	196	12950	0.97	ng	# 76
17) 2,4,5-Trichlorophenol	13.83	196	14014	0.99	ng	# 93
19) 2,4-Dinitrophenol	15.49	184	475	0.45	ng	# 1
20) 4-Nitrophenol	15.52	139	4514	0.13	ng	# 72
22) 4,6-Dinitro-2-methylphenol	16.40	198	2785	0.85	ng	# 74
23) Pentachlorophenol	17.63	266	2727	0.79	ng	# 94

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

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