

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101315\
 Data File : BM002896.D
 Acq On : 13 Oct 2015 20:59
 Operator : UM/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC002

Quant Time: Oct 14 17:34:05 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	80044	5.00	ng	0.00
8) Naphthalene-d8	11.50	136	343434	5.00	ng	0.00
14) Acenaphthene-d10	15.25	164	187622	5.00	ng	0.00
21) Phenanthrene-d10	17.95	188	424144	5.00	ng	0.00
24) Chrysene-d12	22.04	240	403790	5.00	ng	0.00
26) Perylene-d12	24.61	264	380723	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	6.13	112	30054	2.09	ng	0.00
3) Phenol-d6	7.79	99	41162	1.87	ng	0.00
9) Nitrobenzene-d5	9.85	82	36664	1.71	ng	-0.01
15) 2,4,6-Tribromophenol	16.71	330	14249	2.08	ng	0.00
18) 2-Fluorobiphenyl	13.87	172	116963	1.90	ng	0.00
25) Terphenyl-d14	20.48	244	117273	1.91	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) 2-Chlorophenol	8.21	128	39913	1.96	ng	# 75
5) Phenol	7.82	94	43033	1.84	ng	# 95
6) 2-Methylphenol	9.10	107	29354	1.78	ng	# 84
7) 3+4-Methylphenols	9.44	107	41269	2.02	ng	# 34
10) 2-Nitrophenol	10.62	139	17017	1.97	ng	# 70
11) 2,4-Dimethylphenol	10.64	122	37876	2.00	ng	# 77
12) 2,4-Dichlorophenol	11.17	162	31427	2.13	ng	# 82
13) 4-Chloro-3-methylphenol	12.75	107	37919	2.03	ng	# 81
16) 2,4,6-Trichlorophenol	13.72	196	24628	2.01	ng	# 67
17) 2,4,5-Trichlorophenol	13.81	196	27278	2.09	ng	# 92
19) 2,4-Dinitrophenol	15.45	184	1675	1.74	ng	# 82
20) 4-Nitrophenol	15.49	139	11246	0.35	ng	# 65
22) 4,6-Dinitro-2-methylphenol	16.38	198	6837	2.25	ng	# 72
23) Pentachlorophenol	17.62	266	6825	2.12	ng	# 85

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

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