

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101315\
 Data File : BM002894.D
 Acq On : 13 Oct 2015 19:46
 Operator : UM/IZ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.8

Quant Time: Oct 14 17:33:11 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	85725	5.00	ng	0.00
8) Naphthalene-d8	11.50	136	361500	5.00	ng	0.00
14) Acenaphthene-d10	15.25	164	194712	5.00	ng	0.00
21) Phenanthrene-d10	17.95	188	451393	5.00	ng	0.00
24) Chrysene-d12	22.04	240	427995	5.00	ng	0.00
26) Perylene-d12	24.61	264	401099	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	6.15	112	11868	0.77	ng	0.02
3) Phenol-d6	7.80	99	17425	0.74	ng	0.02
9) Nitrobenzene-d5	9.86	82	15816	0.70	ng	0.00
15) 2,4,6-Tribromophenol	16.71	330	5239	0.74	ng	0.00
18) 2-Fluorobiphenyl	13.88	172	49823	0.78	ng	0.01
25) Terphenyl-d14	20.48	244	49986	0.77	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) 2-Chlorophenol	8.21	128	16689	0.76	ng	# 69
5) Phenol	7.83	94	17658	0.71	ng	94
6) 2-Methylphenol	9.11	107	13310	0.75	ng	# 85
7) 3+4-Methylphenols	9.46	107	16521	0.75	ng	# 36
10) 2-Nitrophenol	10.64	139	6664	0.73	ng	# 59
11) 2,4-Dimethylphenol	10.66	122	14963	0.75	ng	99
12) 2,4-Dichlorophenol	11.21	162	11926	0.77	ng	93
13) 4-Chloro-3-methylphenol	12.76	107	14613	0.74	ng	88
16) 2,4,6-Trichlorophenol	13.73	196	9329	0.73	ng	# 61
17) 2,4,5-Trichlorophenol	13.83	196	10286	0.76	ng	# 98
19) 2,4-Dinitrophenol	15.51	184	252	0.25	ng	# 1
20) 4-Nitrophenol	15.51	139	3101	0.09	ng	# 75
22) 4,6-Dinitro-2-methylphenol	16.40	198	2032	0.63	ng	78
23) Pentachlorophenol	17.63	266	1836	0.54	ng	88

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

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