

Data Path : Z:\HPCHEM1\BNA_E\Data\BE120914\
 Data File : BE089187.D
 Acq On : 9 Dec 2014 11:02
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Dec 10 00:35:11 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-ACID-BE120814.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 09 19:38:28 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	32032	5.00	ng	0.00
8) Naphthalene-d8	8.51	136	123336	5.00	ng	0.00
13) Acenaphthene-d10	10.66	164	59879	5.00	ng	0.00
19) Phenanthrene-d10	12.72	188	86284	5.00	ng	0.00
22) Chrysene-d12	18.07	240	58883	5.00	ng	0.00
23) Perylene-d12	21.12	264	52947	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	5.27	112	8565	1.04	ng	0.00
3) Phenol-d6	6.56	99	11413	1.05	ng	0.00
14) 2,4,6-Tribromophenol	11.67	330	2039	1.04	ng	0.00
Target Compounds						Qvalue
4) 2-Chlorophenol	6.69	128	9402	1.05	ng	98
5) Phenol	6.57	94	13727	1.06	ng	81
6) 2-Methylphenol	7.31	107	10822	1.05	ng	# 78
7) 3+4-Methylphenols	7.51	107	9358	1.07	ng	# 87
9) 2-Nitrophenol	8.06	139	4418	1.01	ng	# 90
10) 2,4-Dimethylphenol	8.15	122	7902	1.03	ng	91
11) 2,4-Dichlorophenol	8.37	162	6938	1.03	ng	96
12) 4-Chloro-3-methylphenol	9.25	107	7221	1.03	ng	# 90
15) 2,4,6-Trichlorophenol	9.75	196	4178	1.03	ng	# 51
16) 2,4,5-Trichlorophenol	9.81	196	4719	1.04	ng	# 73
17) 2,4-Dinitrophenol	10.79	184	962	0.98	ng	85
18) 4-Nitrophenol	10.91	139	9989	1.08	ng	# 28
20) 4,6-Dinitro-2-methylphenol	11.45	198	1515	0.94	ng	94
21) Pentachlorophenol	12.44	266	1231	0.94	ng	97

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

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