

Data Path : Z:\HPCHEM1\BNA_E\Data\BE120914\
 Data File : BE089189.D
 Acq On : 9 Dec 2014 14:39
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Dec 09 15:35:35 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 08:57:21 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	31118	5.00	ng	0.00
8) Naphthalene-d8	8.51	136	122620	5.00	ng	0.00
13) Acenaphthene-d10	10.66	164	59522	5.00	ng	0.00
19) Phenanthrene-d10	12.72	188	86095	5.00	ng	0.00
22) Chrysene-d12	18.07	240	60401	5.00	ng	0.00
23) Perylene-d12	21.12	264	54209	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	5.27	112	8682	1.09	ng	0.00
3) Phenol-d6	6.56	99	11285	1.07	ng	0.00
14) 2,4,6-Tribromophenol	11.67	330	1996	1.02	ng	0.00
Target Compounds						Qvalue
4) 2-Chlorophenol	6.69	128	9183	1.05	ng	92
5) Phenol	6.57	94	13557	1.08	ng	88
6) 2-Methylphenol	7.31	107	10366	1.03	ng	# 76
7) 3+4-Methylphenols	7.52	107	9391	1.10	ng	# 88
9) 2-Nitrophenol	8.05	139	4492	1.04	ng	# 51
10) 2,4-Dimethylphenol	8.14	122	7758	1.02	ng	# 67
11) 2,4-Dichlorophenol	8.38	162	6634	0.99	ng	95
12) 4-Chloro-3-methylphenol	9.25	107	7041	1.01	ng	# 90
15) 2,4,6-Trichlorophenol	9.75	196	4050	1.01	ng	# 43
16) 2,4,5-Trichlorophenol	9.81	196	4673	1.03	ng	# 73
17) 2,4-Dinitrophenol	10.78	184	965	0.90	ng	# 83
18) 4-Nitrophenol	10.91	139	9869	1.08	ng	# 28
20) 4,6-Dinitro-2-methylphenol	11.45	198	1509	0.94	ng	93
21) Pentachlorophenol	12.44	266	1272	0.96	ng	99

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

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