

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120814\
 Data File : BE089181.D
 Acq On : 8 Dec 2014 23:59
 Operator : TP/JJ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICCC001

Quant Time: Dec 09 03:49:20 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 03:43:23 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	28601	5.00	ng	0.00
8) Naphthalene-d8	8.51	136	110080	5.00	ng	0.00
13) Acenaphthene-d10	10.66	164	53194	5.00	ng	0.00
19) Phenanthrene-d10	12.72	188	78741	5.00	ng	0.00
22) Chrysene-d12	18.07	240	56035	5.00	ng	0.00
23) Perylene-d12	21.12	264	51034	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	5.27	112	7997	1.07	ng	0.00
3) Phenol-d6	6.56	99	10272	1.09	ng	0.00
14) 2,4,6-Tribromophenol	11.67	330	1847	0.94	ng	0.00
Target Compounds						Qvalue
4) 2-Chlorophenol	6.69	128	8425	1.20	ng	100
5) Phenol	6.57	94	12262	1.19	ng	100
6) 2-Methylphenol	7.31	107	9991	0.51	ng	100
7) 3+4-Methylphenols	7.52	107	8249	0.34	ng	100
9) 2-Nitrophenol	8.06	139	4067	1.50	ng	100
10) 2,4-Dimethylphenol	8.15	122	7083	1.19	ng	100
11) 2,4-Dichlorophenol	8.38	162	6095	1.18	ng	100
12) 4-Chloro-3-methylphenol	9.26	107	6378	1.16	ng	100
15) 2,4,6-Trichlorophenol	9.75	196	3665	1.07	ng	100
16) 2,4,5-Trichlorophenol	9.81	196	4259	1.21	ng	100
17) 2,4-Dinitrophenol	10.79	184	817	2.17	ng	100
18) 4-Nitrophenol	10.91	139	8824	3.34	ng	100
20) 4,6-Dinitro-2-methylphenol	11.45	198	1323	1.81	ng	100
21) Pentachlorophenol	12.44	266	1119	0.89	ng	100

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

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