

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120814\
 Data File : BE089185.D
 Acq On : 9 Dec 2014 8:50
 Operator : TP/JJ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 ICVBE120814

Quant Time: Dec 09 09:29:28 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	29817	5.00	ng	0.00
8) Naphthalene-d8	8.51	136	115133	5.00	ng	0.00
13) Acenaphthene-d10	10.66	164	55348	5.00	ng	0.00
19) Phenanthrene-d10	12.72	188	81793	5.00	ng	0.00
22) Chrysene-d12	18.07	240	57211	5.00	ng	0.00
23) Perylene-d12	21.12	264	51689	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	5.27	112	8307	1.09	ng	0.00
3) Phenol-d6	6.56	99	10592	1.05	ng	0.00
14) 2,4,6-Tribromophenol	11.67	330	1879	1.03	ng	0.00
Target Compounds						Qvalue
4) 2-Chlorophenol	6.69	128	8703	1.04	ng	96
5) Phenol	6.57	94	12831	1.07	ng	86
6) 2-Methylphenol	7.31	107	10054	1.04	ng	# 78
7) 3+4-Methylphenols	7.52	107	8703	1.07	ng	# 87
9) 2-Nitrophenol	8.06	139	4248	1.04	ng	# 88
10) 2,4-Dimethylphenol	8.15	122	7418	1.04	ng	92
11) 2,4-Dichlorophenol	8.38	162	6365	1.02	ng	95
12) 4-Chloro-3-methylphenol	9.26	107	6616	1.02	ng	# 88
15) 2,4,6-Trichlorophenol	9.75	196	3787	1.01	ng	# 42
16) 2,4,5-Trichlorophenol	9.81	196	4364	1.04	ng	# 70
17) 2,4-Dinitrophenol	10.78	184	913	0.92	ng	88
18) 4-Nitrophenol	10.91	139	9185	1.08	ng	# 28
20) 4,6-Dinitro-2-methylphenol	11.45	198	1457	0.95	ng	93
21) Pentachlorophenol	12.44	266	1112	0.90	ng	95

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

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