

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120814\
 Data File : BE089177.D
 Acq On : 8 Dec 2014 21:45
 Operator : TP/JJ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.1

Manual Integrations
 APPROVED

mohammad
 12/10/2014 4:24:12 PM

Quant Time: Dec 09 04:10:25 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 03:43:18 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	31372	5.00	ng	0.00
8) Naphthalene-d8	8.51	136	120250	5.00	ng	0.00
13) Acenaphthene-d10	10.66	164	58616	5.00	ng	0.00
19) Phenanthrene-d10	12.72	188	89602	5.00	ng	0.00
22) Chrysene-d12	18.07	240	68180	5.00	ng	0.00
23) Perylene-d12	21.13	264	62502	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	5.29	112	604	0.07	ng	0.01
3) Phenol-d6	6.57	99	970	0.09	ng	0.01
14) 2,4,6-Tribromophenol	11.68	330	175	0.25	ng	0.01
Target Compounds						Qvalue
4) 2-Chlorophenol	6.70	128	856	0.11	ng	92
5) Phenol	6.59	94	1205	0.11	ng	91
6) 2-Methylphenol	7.32	107	1046m	0.05	ng	
7) 3+4-Methylphenols	7.54	107	761	0.03	ng	# 84
9) 2-Nitrophenol	8.06	139	352	0.32	ng	# 27
10) 2,4-Dimethylphenol	8.16	122	699	0.11	ng	90
11) 2,4-Dichlorophenol	8.40	162	650	0.18	ng	# 84
12) 4-Chloro-3-methylphenol	9.27	107	664	0.09	ng	# 93
15) 2,4,6-Trichlorophenol	9.75	196	359	0.18	ng	# 63
16) 2,4,5-Trichlorophenol	9.82	196	432	0.16	ng	89
18) 4-Nitrophenol	10.91	139	646	0.19	ng	94
20) 4,6-Dinitro-2-methylphenol	11.46	198	71	0.22	ng	# 1
21) Pentachlorophenol	12.46	266	101m	0.42	ng	

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

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