

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120814\
 Data File : BE089178.D
 Acq On : 8 Dec 2014 22:21
 Operator : TP/JJ
 Sample : SSTDICC0.2
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.2

Manual Integrations
 APPROVED

mohammad
 12/10/2014 4:30:18 PM

Quant Time: Dec 09 09:01:43 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	28928	5.00	ng	0.00
8) Naphthalene-d8	8.51	136	111610	5.00	ng	0.00
13) Acenaphthene-d10	10.66	164	54223	5.00	ng	0.00
19) Phenanthrene-d10	12.72	188	81680	5.00	ng	0.00
22) Chrysene-d12	18.07	240	58342	5.00	ng	0.00
23) Perylene-d12	21.13	264	53160	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	5.28	112	1459	0.19	ng	0.00
3) Phenol-d6	6.57	99	2106	0.22	ng	0.01
14) 2,4,6-Tribromophenol	11.67	330	376	0.33	ng	0.00
Target Compounds						Qvalue
4) 2-Chlorophenol	6.70	128	1780	0.25	ng	89
5) Phenol	6.58	94	2534	0.24	ng	96
6) 2-Methylphenol	7.32	107	2088	0.11	ng	# 87
7) 3+4-Methylphenols	7.53	107	1651	0.07	ng	92
9) 2-Nitrophenol	8.06	139	787	0.46	ng	# 62
10) 2,4-Dimethylphenol	8.15	122	1504	0.25	ng	88
11) 2,4-Dichlorophenol	8.38	162	1404	0.32	ng	91
12) 4-Chloro-3-methylphenol	9.27	107	1382	0.23	ng	# 94
15) 2,4,6-Trichlorophenol	9.75	196	776m	0.30	ng	
16) 2,4,5-Trichlorophenol	9.82	196	873	0.29	ng	87
17) 2,4-Dinitrophenol	10.81	184	96	0.16	ng	# 46
18) 4-Nitrophenol	10.91	139	1858	0.66	ng	92
20) 4,6-Dinitro-2-methylphenol	11.45	198	180	0.38	ng	# 1
21) Pentachlorophenol	12.45	266	173m	0.45	ng	

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

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