

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120814\
 Data File : BE089180.D
 Acq On : 8 Dec 2014 23:27
 Operator : TP/JJ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.8

Manual Integrations
 APPROVED

mohammad
 12/10/2014 4:30:25 PM

Quant Time: Dec 09 04:12:07 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	29077	5.00	ng	0.00
8) Naphthalene-d8	8.51	136	110751	5.00	ng	0.00
13) Acenaphthene-d10	10.66	164	54913	5.00	ng	0.00
19) Phenanthrene-d10	12.72	188	82253	5.00	ng	0.00
22) Chrysene-d12	18.08	240	60677	5.00	ng	0.00
23) Perylene-d12	21.12	264	54853	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	5.28	112	4825	0.64	ng	0.00
3) Phenol-d6	6.56	99	6760	0.70	ng	0.00
14) 2,4,6-Tribromophenol	11.67	330	1191	0.66	ng	0.00
Target Compounds						Qvalue
4) 2-Chlorophenol	6.69	128	5430	0.76	ng	89
5) Phenol	6.58	94	7936	0.76	ng	86
6) 2-Methylphenol	7.31	107	6479	0.33	ng	95
7) 3+4-Methylphenols	7.52	107	5308	0.22	ng	95
9) 2-Nitrophenol	8.06	139	2781	1.09	ng	95
10) 2,4-Dimethylphenol	8.15	122	4606	0.77	ng	97
11) 2,4-Dichlorophenol	8.38	162	4049	0.80	ng	94
12) 4-Chloro-3-methylphenol	9.25	107	4208	0.75	ng	94
15) 2,4,6-Trichlorophenol	9.75	196	2416	0.72	ng	# 82
16) 2,4,5-Trichlorophenol	9.81	196	2753m	0.78	ng	
17) 2,4-Dinitrophenol	10.80	184	476	1.36	ng	# 88
18) 4-Nitrophenol	10.91	139	5839	2.13	ng	92
20) 4,6-Dinitro-2-methylphenol	11.45	198	800	1.16	ng	88
21) Pentachlorophenol	12.44	266	634	0.66	ng	95

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

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