

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE120814\
 Data File : BE089179.D
 Acq On : 8 Dec 2014 22:54
 Operator : TP/JJ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.5

Manual Integrations
 APPROVED

mohammad
 12/10/2014 4:30:22 PM

Quant Time: Dec 09 04:11:37 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	27786	5.00	ng	0.00
8) Naphthalene-d8	8.51	136	106506	5.00	ng	0.00
13) Acenaphthene-d10	10.66	164	51075	5.00	ng	0.00
19) Phenanthrene-d10	12.72	188	78274	5.00	ng	0.00
22) Chrysene-d12	18.07	240	59438	5.00	ng	0.00
23) Perylene-d12	21.13	264	54860	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	5.28	112	3591	0.50	ng	0.00
3) Phenol-d6	6.56	99	5021	0.55	ng	0.00
14) 2,4,6-Tribromophenol	11.67	330	860	0.55	ng	0.00
Target Compounds						Qvalue
4) 2-Chlorophenol	6.69	128	3937	0.58	ng	83
5) Phenol	6.58	94	5788	0.58	ng	90
6) 2-Methylphenol	7.31	107	4795	0.25	ng	93
7) 3+4-Methylphenols	7.52	107	3825	0.16	ng	92
9) 2-Nitrophenol	8.06	139	1952	0.85	ng	# 86
10) 2,4-Dimethylphenol	8.15	122	3351	0.58	ng	95
11) 2,4-Dichlorophenol	8.38	162	2965	0.63	ng	90
12) 4-Chloro-3-methylphenol	9.26	107	3099	0.57	ng	94
15) 2,4,6-Trichlorophenol	9.75	196	1768	0.59	ng	82
16) 2,4,5-Trichlorophenol	9.81	196	2065m	0.64	ng	
17) 2,4-Dinitrophenol	10.80	184	312	0.99	ng	# 82
18) 4-Nitrophenol	10.91	139	4223	1.65	ng	86
20) 4,6-Dinitro-2-methylphenol	11.45	198	517	0.85	ng	# 64
21) Pentachlorophenol	12.45	266	427	0.57	ng	90

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

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