

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101315\
 Data File : BM002893.D
 Acq On : 13 Oct 2015 19:10
 Operator : UM/IZ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.5

Manual Integrations
 APPROVED

MMdadoda
 10/14/2015 6:25:55 PM

Quant Time: Oct 14 17:32:53 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM0101315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 14 17:20:35 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.66	152	89530	5.00	ng	0.00
8) Naphthalene-d8	11.50	136	375220	5.00	ng	0.00
14) Acenaphthene-d10	15.25	164	203751	5.00	ng	0.00
21) Phenanthrene-d10	17.95	188	460181	5.00	ng	0.00
24) Chrysene-d12	22.04	240	422838	5.00	ng	0.00
26) Perylene-d12	24.61	264	397980	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	6.15	112	8414m	0.52	ng	0.02
3) Phenol-d6	7.80	99	11528	0.47	ng	0.02
9) Nitrobenzene-d5	9.87	82	10768	0.46	ng	0.01
15) 2,4,6-Tribromophenol	16.74	330	3217	0.43	ng	0.02
18) 2-Fluorobiphenyl	13.88	172	33095	0.50	ng	0.01
25) Terphenyl-d14	20.48	244	31827	0.49	ng	0.00
Target Compounds						
4) 2-Chlorophenol	8.22	128	10938	0.48	ng	Qvalue 74
5) Phenol	7.83	94	11419	0.44	ng	89
6) 2-Methylphenol	9.11	107	8794	0.48	ng	# 77
7) 3+4-Methylphenols	9.46	107	10625	0.46	ng	# 35
10) 2-Nitrophenol	10.66	139	4279	0.45	ng	# 71
11) 2,4-Dimethylphenol	10.66	122	9646	0.47	ng	92
12) 2,4-Dichlorophenol	11.21	162	7489	0.46	ng	86
13) 4-Chloro-3-methylphenol	12.77	107	9304	0.46	ng	# 74
16) 2,4,6-Trichlorophenol	13.73	196	5922	0.45	ng	# 68
17) 2,4,5-Trichlorophenol	13.84	196	6328	0.45	ng	# 96
19) 2,4-Dinitrophenol	15.56	184	71	0.07	ng	# 1
20) 4-Nitrophenol	15.54	139	1727m	0.05	ng	
22) 4,6-Dinitro-2-methylphenol	16.43	198	1006	0.31	ng	# 24
23) Pentachlorophenol	17.65	266	1014	0.29	ng	95

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

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