

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101415\
 Data File : BM002904.D
 Acq On : 14 Oct 2015 19:28
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001EC

Manual Integrations
 APPROVED

MMdadoda
 10/15/2015 5:29:49 PM

Quant Time: Oct 15 01:15:47 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM0101315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 01:02:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.66	152	83235	5.00	ng	0.01
8) Naphthalene-d8	11.50	136	345451	5.00	ng	0.00
14) Acenaphthene-d10	15.25	164	182289	5.00	ng	0.00
21) Phenanthrene-d10	17.95	188	420393	5.00	ng	0.00
24) Chrysene-d12	22.04	240	398620	5.00	ng	0.00
26) Perylene-d12	24.61	264	368537	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	6.15	112	14740m	0.92	ng	0.00
3) Phenol-d6	7.80	99	19778	0.87	ng	0.00
9) Nitrobenzene-d5	9.86	82	17625	0.93	ng	0.00
15) 2,4,6-Tribromophenol	16.71	330	5675	0.85	ng	0.00
18) 2-Fluorobiphenyl	13.88	172	55416	0.93	ng	0.01
25) Terphenyl-d14	20.48	244	56184	0.93	ng	0.00
Target Compounds						
4) 2-Chlorophenol	8.21	128	18856	0.89	ng	Qvalue # 70
5) Phenol	7.82	94	20277	0.90	ng	83
6) 2-Methylphenol	9.11	107	15489	0.90	ng	# 86
7) 3+4-Methylphenols	9.45	107	18794	0.88	ng	# 35
10) 2-Nitrophenol	10.64	139	7172	0.83	ng	# 64
11) 2,4-Dimethylphenol	10.66	122	16848	0.88	ng	100
12) 2,4-Dichlorophenol	11.19	162	13766	0.93	ng	# 82
13) 4-Chloro-3-methylphenol	12.76	107	16175	0.87	ng	# 84
16) 2,4,6-Trichlorophenol	13.72	196	10289	0.87	ng	# 77
17) 2,4,5-Trichlorophenol	13.83	196	11206	0.89	ng	# 94
19) 2,4-Dinitrophenol	15.51	184	354	0.90	ng	# 1
20) 4-Nitrophenol	15.51	139	3387	0.89	ng	# 75
22) 4,6-Dinitro-2-methylphenol	16.40	198	2025	0.85	ng	80
23) Pentachlorophenol	17.63	266	2265	1.03	ng	94

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

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