

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
 Data File : BM000818.D
 Acq On : 02 Apr 2015 22:08
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Quant Time: Apr 03 18:19:31 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 03 17:58:44 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	25888	5.00	ng	0.00
11) Naphthalene-d8	10.39	136	128297	5.00	ng	0.00
19) Acenaphthene-d10	14.28	164	64587	5.00	ng	0.00
26) Phenanthrene-d10	17.00	188	128358	5.00	ng	0.00
30) Chrysene-d12	21.21	240	122731	5.00	ng	0.00
34) Perylene-d12	23.40	264	109730	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.21	112	81561	11.44	ng	0.00
5) Phenol-d6	6.77	99	140222	14.10	ng	0.00
12) Nitrobenzene-d5	8.77	82	65011	10.86	ng	0.00
20) 2,4,6-Tribromophenol	0.00	330	0d	0.00	ng	0.00
23) 2-Fluorobiphenyl	12.87	172	205272	10.32	ng	0.00
32) Terphenyl-d14	19.65	244	159704	10.81	ng	0.00
Target Compounds						
3) n-Nitrosodimethylamine	3.41	42	42069	12.79	ng	Qvalue # 8
6) 2-Chlorophenol	7.17	128	91330	12.92	ng	85
7) Phenol	6.81	94	127072	11.90	ng	87
8) bis(2-Chloroethyl)ether	7.02	93	95741	12.89	ng	# 71
9) 2-Methylphenol	8.06	107	94686	12.43	ng	# 91
13) Nitrobenzene	8.80	77	78847	11.28	ng	# 65
14) 2-Nitrophenol	9.52	139	40105	12.57	ng	# 80
15) 2,4-Dimethylphenol	9.58	122	88027	12.11	ng	94
16) 2,4-Dichlorophenol	10.04	162	86609	13.44	ng	91
17) Hexachlorobutadiene	10.75	225	51861	10.46	ng	99
18) 4-Chloro-3-methylphenol	11.67	107	81159	12.60	ng	# 80
24) 2,4-Dinitrophenol	14.40	184	13705	21.28	ng	# 36
28) Hexachlorobenzene	16.33	284	80173	11.06	ng	# 95

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

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