

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
 Data File : BM000811.D
 Acq On : 02 Apr 2015 17:57
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.1

Quant Time: Apr 03 18:11:07 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 17:33:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	29650	5.00	ng	0.00
11) Naphthalene-d8	10.39	136	136382	5.00	ng	0.00
19) Acenaphthene-d10	14.28	164	66317	5.00	ng	0.00
26) Phenanthrene-d10	17.00	188	133836	5.00	ng	0.00
30) Chrysene-d12	21.21	240	126845	5.00	ng	0.00
34) Perylene-d12	23.40	264	117903	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.21	112	781	0.10	ng	0.00
5) Phenol-d6	6.77	99	787	0.07	ng	0.00
12) Nitrobenzene-d5	8.77	82	701	0.11	ng	0.00
20) 2,4,6-Tribromophenol	15.77	330	155	0.08	ng	0.00
23) 2-Fluorobiphenyl	12.89	172	2093	0.10	ng	0.02
32) Terphenyl-d14	19.65	244	1465	0.10	ng	0.00

Target Compounds						Qvalue
3) n-Nitrosodimethylamine	3.46	42	351	0.09	ng	# 1
6) 2-Chlorophenol	7.17	128	690	0.09	ng	# 50
7) Phenol	6.81	94	1097	0.09	ng	# 61
8) bis(2-Chloroethyl)ether	7.02	93	776m	0.09	ng	
9) 2-Methylphenol	8.06	107	708	0.08	ng	# 70
10) 3+4-Methylphenols	8.41	107	688	0.07	ng	# 51
13) Nitrobenzene	8.80	77	695	0.09	ng	# 67
14) 2-Nitrophenol	9.52	139	575	0.17	ng	# 22
15) 2,4-Dimethylphenol	9.58	122	675	0.09	ng	# 81
16) 2,4-Dichlorophenol	10.07	162	554	0.08	ng	# 85
17) Hexachlorobutadiene	10.75	225	520	0.10	ng	# 93
18) 4-Chloro-3-methylphenol	11.69	107	617	0.09	ng	# 90
21) 2,4,6-Trichlorophenol	12.70	196	270	0.08	ng	# 63
22) 2,4,5-Trichlorophenol	12.77	196	362	0.08	ng	# 83
28) Hexachlorobenzene	16.33	284	809	0.11	ng	# 75
31) Benzidine	19.30	184	965	0.19	ng	# 80
33) 3,3'-Dichlorobenzidine	21.14	252	758	0.09	ng	# 75

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

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