

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM053115\
 Data File : BM001493.D
 Acq On : 31 May 2015 03:18
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC002

Quant Time: May 31 05:13:51 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM053115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun May 31 05:09:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	14016	5.00	ng	0.00
6) Naphthalene-d8	10.15	136	52030	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	26664	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	52770	5.00	ng	0.00
17) Chrysene-d12	21.03	240	42254	5.00	ng	0.00
21) Perylene-d12	23.11	264	36741	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	5498	2.36	ng	0.00
3) Phenol-d6	6.57	99	7329	2.95	ng	0.00
7) Nitrobenzene-d5	8.53	82	5870	2.09	ng	0.00
10) 2,4,6-Tribromophenol	15.57	330	1966	2.37	ng	0.00
11) 2-Fluorobiphenyl	12.67	172	16097	2.14	ng	0.00
18) Terphenyl-d14	19.46	244	11742	2.22	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.80	93	5519	2.13	ng	86
5) n-Nitroso-di-n-propylamine	8.18	70	4710	3.05	ng	# 96
8) bis(2-Chloroethoxy)methane	9.58	93	7220	2.55	ng	97
12) 2,6-Dinitrotoluene	13.65	165	2880	2.52	ng	94
13) 2,4-Dinitrotoluene	14.45	165	3338	2.53	ng	# 82
14) 4-Chlorophenyl-phenylether	15.11	204	7809	1.94	ng	97
16) 4-Bromophenyl-phenylether	16.02	248	4478	2.16	ng	97
19) Benzo(a)anthracene	21.00	228	23230	8.74	ng	# 85
20) 3,3'-Dichlorobenzidine	20.95	252	8039	2.87	ng	# 85

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

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