

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM053115\
 Data File : BM001491.D
 Acq On : 31 May 2015 02:05
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Quant Time: May 31 05:13:26 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	13996	5.00	ng	0.00
6) Naphthalene-d8	10.15	136	53207	5.00	ng	0.00
9) Acenaphthene-d10	14.06	164	28408	5.00	ng	0.00
15) Phenanthrene-d10	16.80	188	56270	5.00	ng	0.00
17) Chrysene-d12	21.03	240	46471	5.00	ng	0.00
21) Perylene-d12	23.11	264	42420	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	26998	11.62	ng	0.00
3) Phenol-d6	6.57	99	36336	14.65	ng	0.00
7) Nitrobenzene-d5	8.53	82	30416	10.57	ng	0.00
10) 2,4,6-Tribromophenol	15.56	330	11980	13.58	ng	0.00
11) 2-Fluorobiphenyl	12.67	172	80119	9.99	ng	0.00
18) Terphenyl-d14	19.47	244	61632	10.61	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.80	93	26365	10.17	ng	82
5) n-Nitroso-di-n-propylamine	8.18	70	25683	16.67	ng	# 94
8) bis(2-Chloroethoxy)methane	9.58	93	37520	12.94	ng	# 95
12) 2,6-Dinitrotoluene	13.66	165	17750	14.59	ng	# 73
13) 2,4-Dinitrotoluene	14.45	165	22936	16.34	ng	# 88
14) 4-Chlorophenyl-phenylether	15.12	204	41814	9.77	ng	# 75
16) 4-Bromophenyl-phenylether	16.01	248	24552	11.13	ng	# 75
19) Benzo(a)anthracene	21.00	228	117541	40.22	ng	# 81
20) 3,3'-Dichlorobenzidine	20.95	252	46766	15.16	ng	# 92

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

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