

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001430.D
 Acq On : 28 May 2015 14:58
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.1

Quant Time: May 28 15:31:49 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM052815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 28 15:25:05 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	10690	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	32648	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	20534	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	46032	5.00	ng	0.00
17) Chrysene-d12	21.03	240	43196	5.00	ng	0.00
21) Perylene-d12	23.12	264	37982	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.99	112	156	0.09	ng	0.00
3) Phenol-d6	6.59	99	158	0.09	ng	0.00
7) Nitrobenzene-d5	8.53	82	149	0.08	ng	0.00
10) 2,4,6-Tribromophenol	15.57	330	43	0.07	ng	0.00
11) 2-Fluorobiphenyl	12.69	172	530	0.09	ng	0.00
18) Terphenyl-d14	19.48	244	523	0.10	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	182	0.09	ng	# 56
5) n-Nitroso-di-n-propylamine	8.20	70	95	0.09	ng	# 57
8) bis(2-Chloroethoxy)methane	9.61	93	94	0.05	ng	# 20
12) 2,6-Dinitrotoluene	13.67	165	65	0.08	ng	# 16
13) 2,4-Dinitrotoluene	14.48	165	49	0.05	ng	# 79
14) 4-Chlorophenyl-phenylether	15.12	204	289	0.09	ng	# 78
16) 4-Bromophenyl-phenylether	16.02	248	166	0.09	ng	# 82
19) Benzo(a)anthracene	21.01	228	1084	0.41	ng	# 90
20) 3,3'-Dichlorobenzidine	20.97	252	178	0.07	ng	# 90

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

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