

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM053115\
 Data File : BM001496.D
 Acq On : 31 May 2015 05:06
 Operator : TP/IZ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.5

Manual Integrations
 APPROVED

apatel
 6/1/2015 8:44:14 PM

Quant Time: May 31 07:34:16 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	12408	5.00	ng	0.00
6) Naphthalene-d8	10.15	136	46206	5.00	ng	0.00
9) Acenaphthene-d10	14.06	164	24320	5.00	ng	0.00
15) Phenanthrene-d10	16.80	188	49971	5.00	ng	0.00
17) Chrysene-d12	21.03	240	40885	5.00	ng	0.00
21) Perylene-d12	23.11	264	36214	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	1262	0.61	ng	0.00
3) Phenol-d6	6.57	99	1611	0.73	ng	0.00
7) Nitrobenzene-d5	8.53	82	1249	0.50	ng	0.00
10) 2,4,6-Tribromophenol	15.56	330	417	0.55	ng	0.00
11) 2-Fluorobiphenyl	12.67	172	3555	0.52	ng	0.00
18) Terphenyl-d14	19.47	244	2757	0.54	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.80	93	1255	0.55	ng	90
5) n-Nitroso-di-n-propylamine	8.18	70	972	0.71	ng	# 91
8) bis(2-Chloroethoxy)methane	9.58	93	1448	0.58	ng	# 95
12) 2,6-Dinitrotoluene	13.65	165	515	0.49	ng	# 43
13) 2,4-Dinitrotoluene	14.46	165	549	0.46	ng	# 94
14) 4-Chlorophenyl-phenylether	15.12	204	1757	0.48	ng	# 74
16) 4-Bromophenyl-phenylether	16.01	248	1025	0.52	ng	# 79
19) Benzo(a)anthracene	20.99	228	5709	2.22	ng	# 87
20) 3,3'-Dichlorobenzidine	20.95	252	1861m	0.69	ng	

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

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