

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM053115\
 Data File : BM001498.D
 Acq On : 31 May 2015 06:18
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.1

Manual Integrations
 APPROVED

apatel
 6/1/2015 8:44:23 PM

Quant Time: May 31 07:36:40 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	12433	5.00	ng	0.00
6) Naphthalene-d8	10.15	136	45357	5.00	ng	0.00
9) Acenaphthene-d10	14.06	164	24213	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	48962m	5.00	ng	0.00
17) Chrysene-d12	21.01	240	40773	5.00	ng	-0.01
21) Perylene-d12	23.11	264	36194	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	238	0.12	ng	0.00
3) Phenol-d6	6.57	99	290	0.13	ng	0.00
7) Nitrobenzene-d5	8.53	82	234	0.10	ng	0.00
10) 2,4,6-Tribromophenol	15.58	330	77m	0.10	ng	0.00
11) 2-Fluorobiphenyl	12.67	172	668	0.10	ng	0.00
18) Terphenyl-d14	19.46	244	531	0.10	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.80	93	238	0.10	ng	87
5) n-Nitroso-di-n-propylamine	8.18	70	185	0.14	ng	# 52
8) bis(2-Chloroethoxy)methane	9.60	93	249	0.10	ng	# 55
12) 2,6-Dinitrotoluene	13.66	165	81	0.08	ng	# 53
13) 2,4-Dinitrotoluene	14.46	165	79	0.07	ng	# 60
14) 4-Chlorophenyl-phenylether	15.11	204	340m	0.09	ng	
16) 4-Bromophenyl-phenylether	16.01	248	194	0.10	ng	# 57
19) Benzo(a)anthracene	21.00	228	1232	0.48	ng	# 88
20) 3,3'-Dichlorobenzidine	20.95	252	311m	0.11	ng	

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

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