

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM053115\
 Data File : BM001495.D
 Acq On : 31 May 2015 04:30
 Operator : TP/IZ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.8

Manual Integrations
 APPROVED

apatel
 6/1/2015 8:44:09 PM

Quant Time: May 31 05:20:39 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	13661	5.00	ng	0.00
6) Naphthalene-d8	10.15	136	51919	5.00	ng	0.00
9) Acenaphthene-d10	14.06	164	26726	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	51735	5.00	ng	0.00
17) Chrysene-d12	21.01	240	42099	5.00	ng	-0.01
21) Perylene-d12	23.11	264	37137	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	2190	0.97	ng	0.00
3) Phenol-d6	6.57	99	2915	1.20	ng	0.00
7) Nitrobenzene-d5	8.53	82	2228	0.79	ng	0.00
10) 2,4,6-Tribromophenol	15.56	330	727	0.88	ng	-0.01
11) 2-Fluorobiphenyl	12.67	172	6381	0.85	ng	0.00
18) Terphenyl-d14	19.46	244	4577	0.87	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.80	93	2304	0.91	ng	98
5) n-Nitroso-di-n-propylamine	8.18	70	1717	1.14	ng	99
8) bis(2-Chloroethoxy)methane	9.58	93	2708	0.96	ng	99
12) 2,6-Dinitrotoluene	13.65	165	959	0.84	ng	# 55
13) 2,4-Dinitrotoluene	14.45	165	1053	0.80	ng	# 82
14) 4-Chlorophenyl-phenylether	15.11	204	3263m	0.81	ng	
16) 4-Bromophenyl-phenylether	16.01	248	1764	0.87	ng	# 62
19) Benzo(a)anthracene	21.00	228	9292	3.51	ng	# 87
20) 3,3'-Dichlorobenzidine	20.95	252	3126m	1.12	ng	

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

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