

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM053115\
 Data File : BM001494.D
 Acq On : 31 May 2015 03:54
 Operator : TP/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICCC001

Quant Time: May 31 05:14:04 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	13029	5.00	ng	0.00
6) Naphthalene-d8	10.15	136	49894	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	25979	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	52013	5.00	ng	0.00
17) Chrysene-d12	21.02	240	40526	5.00	ng	0.00
21) Perylene-d12	23.11	264	35538	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	2613	1.21	ng	0.00
3) Phenol-d6	6.57	99	3527	1.53	ng	0.00
7) Nitrobenzene-d5	8.53	82	2541	0.94	ng	0.00
10) 2,4,6-Tribromophenol	15.57	330	886	1.10	ng	0.00
11) 2-Fluorobiphenyl	12.67	172	7535	1.03	ng	0.00
18) Terphenyl-d14	19.47	244	5464	1.08	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.80	93	2748	1.14	ng	100
5) n-Nitroso-di-n-propylamine	8.18	70	2015	1.40	ng	100
8) bis(2-Chloroethoxy)methane	9.58	93	3217	1.18	ng	100
12) 2,6-Dinitrotoluene	13.65	165	1181	1.06	ng	100
13) 2,4-Dinitrotoluene	14.45	165	1300	1.01	ng	100
14) 4-Chlorophenyl-phenylether	15.11	204	3659	0.94	ng	100
16) 4-Bromophenyl-phenylether	16.02	248	2100	1.03	ng	100
19) Benzo(a)anthracene	21.01	228	10670	4.19	ng	100
20) 3,3'-Dichlorobenzidine	20.95	252	3401	1.26	ng	100

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

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