

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001426.D
 Acq On : 28 May 2015 11:56
 Operator : TP/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICCC001

Quant Time: May 28 15:26:56 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	11929	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	38490	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	22871	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	51197	5.00	ng	0.00
17) Chrysene-d12	21.03	240	48143	5.00	ng	0.00
21) Perylene-d12	23.12	264	41281	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.98	112	1862	0.97	ng	0.00
3) Phenol-d6	6.59	99	1904	0.98	ng	0.00
7) Nitrobenzene-d5	8.53	82	2048	0.98	ng	0.00
10) 2,4,6-Tribromophenol	15.57	330	691	0.99	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	6537	1.01	ng	0.00
18) Terphenyl-d14	19.48	244	6032	1.01	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	2060	0.96	ng	100
5) n-Nitroso-di-n-propylamine	8.18	70	1268	1.02	ng	100
8) bis(2-Chloroethoxy)methane	9.60	93	2156	1.05	ng	100
12) 2,6-Dinitrotoluene	13.67	165	919	0.95	ng	100
13) 2,4-Dinitrotoluene	14.45	165	1051	0.94	ng	100
14) 4-Chlorophenyl-phenylether	15.12	204	3663	1.05	ng	100
16) 4-Bromophenyl-phenylether	16.02	248	2090	1.04	ng	100
19) Benzo(a)anthracene	21.01	228	11800	4.03	ng	100
20) 3,3'-Dichlorobenzidine	20.97	252	2788	0.92	ng	100

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

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