

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001427.D
 Acq On : 28 May 2015 12:32
 Operator : TP/IZ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.8

Quant Time: May 28 15:27:06 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	11890	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	37743	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	22665	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	52431	5.00	ng	0.00
17) Chrysene-d12	21.03	240	47230	5.00	ng	0.00
21) Perylene-d12	23.12	264	40773	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.98	112	1442	0.76	ng	0.00
3) Phenol-d6	6.57	99	1477	0.76	ng	-0.03
7) Nitrobenzene-d5	8.53	82	1583	0.77	ng	0.00
10) 2,4,6-Tribromophenol	15.58	330	520	0.75	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	5160	0.81	ng	0.00
18) Terphenyl-d14	19.48	244	4767	0.82	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	1592	0.74	ng	100
5) n-Nitroso-di-n-propylamine	8.18	70	1002	0.81	ng	95
8) bis(2-Chloroethoxy)methane	9.60	93	1623	0.80	ng	96
12) 2,6-Dinitrotoluene	13.66	165	693	0.73	ng	# 61
13) 2,4-Dinitrotoluene	14.46	165	780	0.70	ng	92
14) 4-Chlorophenyl-phenylether	15.13	204	2772	0.80	ng	89
16) 4-Bromophenyl-phenylether	16.02	248	1625	0.79	ng	# 80
19) Benzo(a)anthracene	21.01	228	9116	3.17	ng	100
20) 3,3'-Dichlorobenzidine	20.96	252	2091	0.70	ng	95

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

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