

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001431.D
 Acq On : 28 May 2015 15:35
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM052815

Quant Time: May 28 16:12:30 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 15:41:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	11426	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	36916	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	22352	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	52057	5.00	ng	0.00
17) Chrysene-d12	21.03	240	48257	5.00	ng	0.00
21) Perylene-d12	23.11	264	40883	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.98	112	1700	0.97	ng	0.00
3) Phenol-d6	6.57	99	1747	0.95	ng	-0.03
7) Nitrobenzene-d5	8.53	82	1887	0.96	ng	0.00
10) 2,4,6-Tribromophenol	15.58	330	648	0.91	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	6130	0.99	ng	0.00
18) Terphenyl-d14	19.48	244	5863	0.98	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	1851	0.96	ng	99
5) n-Nitroso-di-n-propylamine	8.18	70	1177	0.94	ng	97
8) bis(2-Chloroethoxy)methane	9.60	93	1954	0.90	ng	99
12) 2,6-Dinitrotoluene	13.66	165	852	0.87	ng	# 67
13) 2,4-Dinitrotoluene	14.46	165	969	0.83	ng	# 87
14) 4-Chlorophenyl-phenylether	15.13	204	3337	0.97	ng	88
16) 4-Bromophenyl-phenylether	16.02	248	2002	0.97	ng	# 81
19) Benzo(a)anthracene	21.01	228	11155	0.92	ng	100
20) 3,3'-Dichlorobenzidine	20.96	252	2642	0.90	ng	98

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

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