

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040715\
 Data File : BM000877.D
 Acq On : 07 Apr 2015 13:27
 Operator : TP/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICCC001

Quant Time: Apr 07 14:36:11 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 07 14:10:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	15467	5.00	ng	0.00
11) Naphthalene-d8	10.39	136	77212	5.00	ng	0.00
19) Acenaphthene-d10	14.26	164	41936	5.00	ng	0.00
26) Phenanthrene-d10	17.01	188	92382	5.00	ng	0.00
30) Chrysene-d12	21.20	240	82076	5.00	ng	0.00
34) Perylene-d12	23.38	264	74028	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.21	112	5031	1.16	ng	0.00
5) Phenol-d6	6.77	99	7704	1.25	ng	0.00
12) Nitrobenzene-d5	8.77	82	3568	0.99	ng	0.00
20) 2,4,6-Tribromophenol	15.75	330	1166	0.96	ng	0.00
23) 2-Fluorobiphenyl	12.87	172	12881	1.00	ng	0.00
32) Terphenyl-d14	19.65	244	10108	1.02	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.12	88	2619	1.08	ng	100
3) n-Nitrosodimethylamine	3.44	42	2182	1.10	ng	100
6) 2-Chlorophenol	7.17	128	6037	1.36	ng	100
7) Phenol	6.81	94	6118	0.96	ng	100
8) bis(2-Chloroethyl)ether	7.02	93	5874	1.26	ng	100
9) 2-Methylphenol	8.06	107	4634	1.02	ng	100
10) 3+4-Methylphenols	8.38	107	5776	1.18	ng	100
13) Nitrobenzene	8.80	77	3752	0.90	ng	100
14) 2-Nitrophenol	9.52	139	2139	0.97	ng	100
15) 2,4-Dimethylphenol	9.58	122	4347	1.00	ng	100
16) 2,4-Dichlorophenol	10.04	162	4116	1.06	ng	100
17) Hexachlorobutadiene	10.72	225	3535	1.16	ng	100
18) 4-Chloro-3-methylphenol	11.67	107	3952	1.02	ng	100
21) 2,4,6-Trichlorophenol	12.69	196	2025	0.94	ng	100
22) 2,4,5-Trichlorophenol	12.75	196	2679	1.00	ng	100
25) 4-Nitrophenol	14.51	139	614	0.79	ng	100
27) 4,6-Dinitro-2-methylphenol	15.40	198	835	1.02	ng	100
28) Hexachlorobenzene	16.32	284	5202	1.00	ng	100
29) Pentachlorophenol	16.67	266	400	0.50	ng	100
31) Benzidine	19.27	184	3623	0.96	ng	100
33) 3,3'-Dichlorobenzidine	21.12	252	5385	0.96	ng	100

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

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040715\
Data File : BM000877.D
Acq On : 07 Apr 2015 13:27
Operator : TP/IZ
Sample : SSTDICCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SSTDICCC001

Quant Time: Apr 07 14:36:11 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 07 14:10:39 2015
Response via : Initial Calibration

