

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112116\
 Data File : BF091254.D
 Acq On : 21 Nov 2016 20:25
 Operator : UM/SJ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Quant Time: Nov 22 00:22:08 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 22 00:03:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.57	152	219349	20.00	ng	0.00
21) Naphthalene-d8	7.86	136	848143	20.00	ng	0.00
38) Acenaphthene-d10	9.61	164	452378	20.00	ng	0.00
63) Phenanthrene-d10	11.08	188	737084	20.00	ng	0.00
75) Chrysene-d12	13.71	240	657624	20.00	ng	0.00
86) Perylene-d12	15.07	264	577528	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.15	112	243708	18.56	ng	0.01
7) Phenol-d6	6.22	99	356936	20.29	ng	-0.01
23) Nitrobenzene-d5	7.14	82	326646	20.91	ng	-0.01
41) 2,4,6-Tribromophenol	10.40	330	86041	21.01	ng	0.00
44) 2-Fluorobiphenyl	8.93	172	622726	23.54	ng	-0.01
78) Terphenyl-d14	12.67	244	574031	21.67	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.00	88	62064	8.44	ng	97
3) Pyridine	2.65	79	132618	7.57	ng	93
4) n-Nitrosodimethylamine	2.59	42	65277	9.00	ng	95
6) Aniline	6.24	93	194996	9.44	ng	# 73
8) 2-Chlorophenol	6.35	128	149394	9.66	ng	# 79
9) Benzaldehyde	6.12	77	91850	9.63	ng	88
10) Phenol	6.24	94	203553	10.29	ng	# 67
11) bis(2-Chloroethyl)ether	6.30	93	148160	9.05	ng	94
12) 1,3-Dichlorobenzene	6.51	146	154384	9.75	ng	100
13) 1,4-Dichlorobenzene	6.59	146	166699	9.81	ng	100
14) 1,2-Dichlorobenzene	6.74	146	154911	10.08	ng	97
15) Benzyl Alcohol	6.74	79	125310	9.91	ng	97
16) 2,2'-oxybis(1-Chloropropan	6.86	45	305855	12.50	ng	93
17) 2-Methylphenol	7.01	107	150057	11.77	ng	96
18) Hexachloroethane	7.08	117	59794	9.19	ng	92
19) n-Nitroso-di-n-propylamine	6.99	70	120090	9.88	ng	96
20) 3+4-Methylphenols	7.01	107	150057	10.12	ng	96
22) Acetophenone	6.99	105	222334	11.37	ng	# 91
24) Nitrobenzene	7.16	77	176292	10.47	ng	93
25) Isophorone	7.40	82	305731	10.24	ng	97
26) 2-Nitrophenol	7.48	139	70564	9.65	ng	93
27) 2,4-Dimethylphenol	7.54	122	115288	9.74	ng	94
28) bis(2-Chloroethoxy)methane	7.63	93	158905	9.86	ng	98
29) 2,4-Dichlorophenol	7.73	162	118072	11.05	ng	96
30) 1,2,4-Trichlorobenzene	7.80	180	138167	11.68	ng	99
31) Naphthalene	7.88	128	425826	10.60	ng	98
32) Benzoic acid	7.66	122	51814	6.30	ng	94
33) 4-Chloroaniline	7.95	127	159241	9.51	ng	97
34) Hexachlorobutadiene	8.01	225	75596	11.60	ng	98
35) Caprolactam	8.29	113	33772	10.33	ng	# 79
36) 4-Chloro-3-methylphenol	8.44	107	126765	10.08	ng	90
37) 2-Methylnaphthalene	8.57	142	286495	10.98	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.74	216	140762	12.35	ng	98
40) Hexachlorocyclopentadiene	8.73	237	14078	3.87	ng	96

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112116\
 Data File : BF091254.D
 Acq On : 21 Nov 2016 20:25
 Operator : UM/SJ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Quant Time: Nov 22 00:22:08 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 22 00:03:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.86	196	75138	10.16	nq	98
43) 2,4,5-Trichlorophenol	8.91	196	87380	11.29	nq	96
45) 1,1'-Biphenyl	9.04	154	389393	11.95	nq	96
46) 2-Chloronaphthalene	9.06	162	280883	11.33	nq	93
47) 2-Nitroaniline	9.16	65	94923	10.35	nq	# 57
48) Acenaphthylene	9.47	152	414498	10.68	nq	98
49) Dimethylphthalate	9.34	163	332068	11.18	nq	100
50) 2,6-Dinitrotoluene	9.40	165	66318	10.34	nq	# 79
51) Acenaphthene	9.64	154	257270	10.50	nq	99
52) 3-Nitroaniline	9.58	138	71834	9.60	nq	93
53) 2,4-Dinitrophenol	9.71	184	23623	7.74	nq	# 76
54) Dibenzofuran	9.81	168	371065	11.25	nq	94
55) 4-Nitrophenol	9.79	139	10250m	2.64	nq	
56) 2,4-Dinitrotoluene	9.81	165	93230	11.86	nq	# 40
57) Fluorene	10.16	166	300743	11.18	nq	100
58) 2,3,4,6-Tetrachlorophenol	9.94	232	71605	11.75	nq	97
59) Diethylphthalate	10.04	149	318464	10.74	nq	96
60) 4-Chlorophenyl-phenylether	10.14	204	137829	11.44	nq	# 73
61) 4-Nitroaniline	10.19	138	61165	8.14	nq	92
62) Azobenzene	10.30	77	365661	10.36	nq	84
64) 4,6-Dinitro-2-methylphenol	10.22	198	38496	8.48	nq	# 64
65) n-Nitrosodiphenylamine	10.27	169	250952	10.76	nq	99
66) 4-Bromophenyl-phenylether	10.64	248	89059	11.52	nq	# 88
67) Hexachlorobenzene	10.69	284	90433	10.75	nq	# 66
68) Atrazine	10.80	200	75513	11.61	nq	96
69) Pentachlorophenol	10.91	266	29230	7.94	nq	97
70) Phenanthrene	11.10	178	446253	11.66	nq	99
71) Anthracene	11.16	178	434903	10.44	nq	98
72) Carbazole	11.32	167	415726	11.12	nq	97
73) Di-n-butylphthalate	11.66	149	459356	9.93	nq	# 98
74) Fluoranthene	12.29	202	454404	11.02	nq	97
76) Benzidine	12.43	184	137925	9.35	nq	97
77) Pyrene	12.51	202	497077	11.19	nq	99
79) Butylbenzylphthalate	13.15	149	193515	9.28	nq	87
80) Benzo(a)anthracene	13.70	228	397975	10.63	nq	99
81) 3,3'-Dichlorobenzidine	13.68	252	141789	10.65	nq	99
82) Chrysene	13.73	228	409797	11.00	nq	99
83) Bis(2-ethylhexyl)phthalate	13.71	149	268036	9.72	nq	98
84) Di-n-octyl phthalate	14.32	149	451806	9.94	nq	99
85) Indeno(1,2,3-cd)pyrene	16.31	276	320110	10.42	nq	98
87) Benzo(b)fluoranthene	14.71	252	371439m	9.89	nq	
88) Benzo(k)fluoranthene	14.73	252	351403m	10.33	nq	
89) Benzo(a)pyrene	15.01	252	336242	10.13	nq	99
90) Dibenzo(a,h)anthracene	16.31	278	266534	9.26	nq	96
91) Benzo(g,h,i)perylene	16.66	276	262779	9.93	nq	# 95

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

