

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120616\
 Data File : BF091457.D
 Acq On : 6 Dec 2016 14:35
 Operator : UM/SJ
 Sample : PB95080BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB95080BS

Quant Time: Dec 07 00:34:16 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 07 00:12:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	126499	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	534162	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	309108	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	547774	20.00	ng	0.00
75) Chrysene-d12	13.99	240	333973	20.00	ng	0.00
86) Perylene-d12	15.41	264	311724	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	854001	110.29	ng	0.01
7) Phenol-d6	6.48	99	1131431	118.70	ng	0.01
23) Nitrobenzene-d5	7.41	82	775496	81.36	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	346555	118.43	ng	0.01
44) 2-Fluorobiphenyl	9.20	172	1499914	87.86	ng	0.00
78) Terphenyl-d14	12.95	244	1543490	100.45	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	122651	35.01	ng	# 82
3) Pyridine	3.24	79	346361	34.94	ng	# 84
4) n-Nitrosodimethylamine	3.18	42	238723	45.89	ng	98
6) Aniline	6.50	93	317363	25.06	ng	# 75
8) 2-Chlorophenol	6.63	128	386348	45.51	ng	93
9) Benzaldehyde	6.38	77	114637	18.70	ng	96
10) Phenol	6.49	94	438853	39.13	ng	92
11) bis(2-Chloroethyl)ether	6.58	93	361263	45.08	ng	94
12) 1,3-Dichlorobenzene	6.78	146	370768m	39.26	ng	
13) 1,4-Dichlorobenzene	6.86	146	410091	43.65	ng	98
14) 1,2-Dichlorobenzene	7.02	146	361752m	41.34	ng	
15) Benzyl Alcohol	6.98	79	320821	42.06	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.12	45	682518	39.61	ng	71
17) 2-Methylphenol	7.10	107	285617	42.56	ng	# 74
18) Hexachloroethane	7.35	117	138159	41.42	ng	# 85
19) n-Nitroso-di-n-propylamine	7.26	70	259657	43.24	ng	# 95
20) 3+4-Methylphenols	7.26	107	389235	47.51	ng	# 67
22) Acetophenone	7.26	105	519660	41.05	ng	# 84
24) Nitrobenzene	7.43	77	413057	42.33	ng	# 79
25) Isophorone	7.67	82	712607	42.20	ng	98
26) 2-Nitrophenol	7.74	139	170064	38.24	ng	98
27) 2,4-Dimethylphenol	7.78	122	334788	40.24	ng	97
28) bis(2-Chloroethoxy)methane	7.88	93	406342	39.98	ng	100
29) 2,4-Dichlorophenol	7.99	162	322005	44.00	ng	92
30) 1,2,4-Trichlorobenzene	8.07	180	359975	43.87	ng	97
31) Naphthalene	8.15	128	1071634	42.21	ng	99
32) Benzoic acid	7.90	122	145333	23.70	ng	89
33) 4-Chloroaniline	8.20	127	148529	13.69	ng	98
34) Hexachlorobutadiene	8.28	225	208214	41.27	ng	98
35) Caprolactam	8.57	113	86880	41.30	ng	97
36) 4-Chloro-3-methylphenol	8.69	107	344476	43.60	ng	92
37) 2-Methylnaphthalene	8.84	142	711517	41.25	ng	93
39) 1,2,4,5-Tetrachlorobenzene	9.01	216	363647	41.73	ng	98
40) Hexachlorocyclopentadiene	9.00	237	273429	67.69	ng	98

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120616\
 Data File : BF091457.D
 Acq On : 6 Dec 2016 14:35
 Operator : UM/SJ
 Sample : PB95080BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB95080BS

Quant Time: Dec 07 00:34:16 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 07 00:12:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.12	196	243966m	43.08	ng	
43) 2,4,5-Trichlorophenol	9.17	196	223791m	39.43	ng	
45) 1,1'-Biphenyl	9.30	154	890090	40.05	ng	97
46) 2-Chloronaphthalene	9.33	162	640547	38.70	ng	99
47) 2-Nitroaniline	9.42	65	226688	38.13	ng	# 75
48) Acenaphthylene	9.74	152	995400	38.20	ng	99
49) Dimethylphthalate	9.61	163	838188	44.78	ng	99
50) 2,6-Dinitrotoluene	9.67	165	200196	47.86	ng	# 72
51) Acenaphthene	9.91	154	646307	36.77	ng	97
52) 3-Nitroaniline	9.83	138	107273	22.15	ng	91
53) 2,4-Dinitrophenol	9.93	184	37238	20.36	ng	# 88
54) Dibenzofuran	10.08	168	942740	40.06	ng	98
55) 4-Nitrophenol	10.00	139	189285	54.12	ng	# 67
56) 2,4-Dinitrotoluene	10.07	165	258061	47.56	ng	# 80
57) Fluorene	10.42	166	705163	39.03	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.21	232	185177	38.21	ng	93
59) Diethylphthalate	10.31	149	825672	44.69	ng	# 94
60) 4-Chlorophenyl-phenylether	10.42	204	398934	44.03	ng	99
61) 4-Nitroaniline	10.45	138	166052	36.52	ng	87
62) Azobenzene	10.58	77	843061	44.72	ng	99
64) 4,6-Dinitro-2-methylphenol	10.47	198	40103	13.29	ng	# 75
65) n-Nitrosodiphenylamine	10.54	169	627078	37.76	ng	98
66) 4-Bromophenyl-phenylether	10.92	248	242933	41.26	ng	# 72
67) Hexachlorobenzene	10.97	284	245976	38.66	ng	# 84
68) Atrazine	11.06	200	218256	40.58	ng	98
69) Pentachlorophenol	11.17	266	166966	44.58	ng	97
70) Phenanthrene	11.38	178	1029568	36.17	ng	100
71) Anthracene	11.44	178	1250508	44.27	ng	98
72) Carbazole	11.59	167	935246	36.48	ng	98
73) Di-n-butylphthalate	11.93	149	1231437	42.39	ng	99
74) Fluoranthene	12.57	202	1179794	43.76	ng	96
76) Benzidine	12.69	184	95971	9.79	ng	97
77) Pyrene	12.80	202	1122099	44.27	ng	99
79) Butylbenzylphthalate	13.42	149	413662	43.59	ng	97
80) Benzo(a)anthracene	13.98	228	792303	40.57	ng	98
81) 3,3'-Dichlorobenzidine	13.95	252	151630	22.04	ng	99
82) Chrysene	14.01	228	731509	37.85	ng	98
83) Bis(2-ethylhexyl)phthalate	13.98	149	549484	45.68	ng	97
84) Di-n-octyl phthalate	14.60	149	888192	48.79	ng	95
85) Indeno(1,2,3-cd)pyrene	16.80	276	748314	42.28	ng	# 91
87) Benzo(b)fluoranthene	15.01	252	792711m	40.91	ng	
88) Benzo(k)fluoranthene	15.03	252	720744m	44.24	ng	
89) Benzo(a)pyrene	15.35	252	744977	42.81	ng	# 95
90) Dibenzo(a,h)anthracene	16.83	278	621074	38.59	ng	# 91
91) Benzo(g,h,i)perylene	17.23	276	608199	36.65	ng	# 94

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

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120616\
 Data File : BF091457.D
 Acq On : 6 Dec 2016 14:35
 Operator : UM/SJ
 Sample : PB95080BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB95080BS

Quant Time: Dec 07 00:34:16 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 07 00:12:14 2016
 Response via : Initial Calibration

