

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120718\
 Data File : BG038420.D
 Acq On : 8 Dec 2018 2:02
 Operator : JU/SJ
 Sample : PB115253BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115253BS

Quant Time: Dec 08 04:00:01 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG120418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 04 17:15:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	27825	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	119225	20.00	ng	0.00
39) Acenaphthene-d10	14.71	164	73285	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	184062	20.00	ng	-0.01
76) Chrysene-d12	21.73	240	179437	20.00	ng	-0.01
87) Perylene-d12	25.03	264	193715	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	190944	121.44	ng	0.00
7) Phenol-d6	7.22	99	261333	114.99	ng	0.00
23) Nitrobenzene-d5	9.25	82	134228	55.90	ng	0.00
42) 2,4,6-Tribromophenol	16.20	330	122573	136.24	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	315478	61.34	ng	0.00
79) Terphenyl-d14	20.05	244	618270	73.81	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.50	88	22877	31.745	ng	# 91
3) Pyridine	3.90	79	69369	32.176	ng	90
4) n-Nitrosodimethylamine	3.82	42	39663	41.999	ng	# 77
6) Aniline	7.40	93	94838	32.674	ng	97
8) 2-Chlorophenol	7.63	128	73296	40.138	ng	99
9) Benzaldehyde	7.21	77	30114	20.926	ng	99
10) Phenol	7.25	94	92491	38.520	ng	92
11) bis(2-Chloroethyl)ether	7.48	93	64777	32.940	ng	97
12) 1,3-Dichlorobenzene	7.95	146	72918	34.165	ng	94
13) 1,4-Dichlorobenzene	8.10	146	77847	35.253	ng	97
14) 1,2-Dichlorobenzene	8.42	146	73378	35.092	ng	98
15) Benzyl Alcohol	8.31	79	70495	37.731	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.59	45	147430	32.960	ng	96
17) 2-Methylphenol	8.51	107	66278	39.324	ng	94
18) Hexachloroethane	9.15	117	28529	35.795	ng	94
19) n-Nitroso-di-n-propylamine	8.87	70	56204	33.496	ng	91
20) 3+4-Methylphenols	8.83	107	89939	37.180	ng	98
22) Acetophenone	8.90	105	107684	37.238	ng	# 97
24) Nitrobenzene	9.29	77	87694	35.943	ng	99
25) Isophorone	9.80	82	184262	43.662	ng	98
26) 2-Nitrophenol	10.00	139	45477	40.228	ng	90
27) 2,4-Dimethylphenol	10.04	122	81807	50.382	ng	99
28) bis(2-Chloroethoxy)methane	10.28	93	99596	41.167	ng	97
29) 2,4-Dichlorophenol	10.53	162	112648	60.776	ng	99
30) 1,2,4-Trichlorobenzene	10.74	180	96991	44.577	ng	99
31) Naphthalene	10.94	128	236994	41.614	ng	99
32) Benzoic acid	10.19	122	38033	34.481	ng	91
33) 4-Chloroaniline	11.06	127	24535	9.732	ng	95
34) Hexachlorobutadiene	11.20	225	50692	37.241	ng	98
35) Caprolactam	11.84	113	29494	38.480	ng	93
36) 4-Chloro-3-methylphenol	12.16	107	85145	41.345	ng	99
37) 2-Methylnaphthalene	12.54	142	170547	42.051	ng	97
38) 1-Methylnaphthalene	12.75	142	162204	41.757	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.90	216	96139	38.124	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG120718\
 Data File : BG038420.D
 Acq On : 8 Dec 2018 2:02
 Operator : JU/SJ
 Sample : PB115253BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115253BS

Quant Time: Dec 08 04:00:01 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG120418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 04 17:15:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.87	237	134700	97.201	ng	98
43) 2,4,6-Trichlorophenol	13.14	196	68127	41.705	ng	99
44) 2,4,5-Trichlorophenol	13.21	196	77894	45.011	ng	96
46) 1,1'-Biphenyl	13.54	154	224187	36.190	ng	98
47) 2-Chloronaphthalene	13.59	162	179048	38.303	ng	98
48) 2-Nitroaniline	13.80	65	63515	40.500	ng	98
49) Acenaphthylene	14.43	152	298907	42.457	ng	99
50) Dimethylphthalate	14.15	163	246042	41.857	ng	99
51) 2,6-Dinitrotoluene	14.29	165	55782	41.544	ng	95
52) Acenaphthene	14.77	154	168599	39.283	ng	98
53) 3-Nitroaniline	14.62	138	42467	29.366	ng #	99
54) 2,4-Dinitrophenol	14.82	184	71774	97.512	ng #	84
55) Dibenzofuran	15.11	168	290277	40.911	ng	97
56) 4-Nitrophenol	14.92	139	86111	75.381	ng #	76
57) 2,4-Dinitrotoluene	15.07	165	80174	43.292	ng	97
58) Fluorene	15.75	166	231619	44.308	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.32	232	67847	45.141	ng	98
60) Diethylphthalate	15.51	149	256761	39.428	ng	98
61) 4-Chlorophenyl-phenylether	15.74	204	126559	40.307	ng	99
62) 4-Nitroaniline	15.79	138	58854	41.379	ng	85
63) Azobenzene	16.03	77	229481	40.265	ng	95
65) 4,6-Dinitro-2-methylphenol	15.83	198	50026	41.481	ng	96
66) n-Nitrosodiphenylamine	15.96	169	213409	41.231	ng	99
67) 4-Bromophenyl-phenylether	16.63	248	84395	42.459	ng	98
68) Hexachlorobenzene	16.74	284	86598	42.239	ng	98
69) Atrazine	16.90	200	83609	43.291	ng	96
70) Pentachlorophenol	17.10	266	104680	74.544	ng	96
71) Phenanthrene	17.50	178	407083	44.163	ng	98
72) Anthracene	17.58	178	412320	46.191	ng	99
73) Carbazole	17.86	167	373492	42.157	ng	100
74) Di-n-butylphthalate	18.39	149	457689	44.221	ng	98
75) Fluoranthene	19.50	202	505560	46.944	ng	97
77) Benzidine	19.69	184	181891	30.040	ng	99
78) Pyrene	19.86	202	502250	40.004	ng	98
80) Butylbenzylphthalate	20.73	149	202048	40.013	ng	98
81) Benzo(a)anthracene	21.72	228	473059	42.895	ng	99
82) 3,3'-Dichlorobenzidine	21.63	252	143220	33.383	ng	98
83) Chrysene	21.78	228	440852	42.242	ng	98
84) Bis(2-ethylhexyl)phthalate	21.58	149	278462	38.906	ng	99
85) Di-n-octyl phthalate	22.81	149	674064	54.926	ng	100
86) Indeno(1,2,3-cd)pyrene	28.82	276	514554	43.383	ng	99
88) Benzo(b)fluoranthene	23.98	252	477789	43.651	ng	98
89) Benzo(k)fluoranthene	24.05	252	451257	41.395	ng	99
90) Benzo(a)pyrene	24.88	252	450311	42.414	ng #	97
91) Dibenzo(a,h)anthracene	28.89	278	421162	41.825	ng	97
92) Benzo(g,h,i)perylene	30.02	276	420157	42.055	ng #	95

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

