

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092618\
 Data File : BG036993.D
 Acq On : 26 Sep 2018 22:25
 Operator : JU/SJ
 Sample : PB113238BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113238BS

Manual Integrations
 APPROVED

Sohil
 9/27/2018 1:37:45 PM

Quant Time: Sep 27 08:41:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 24 10:41:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	41971	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	171294	20.00	ng	-0.01
38) Acenaphthene-d10	14.67	164	110054	20.00	ng	0.00
63) Phenanthrene-d10	17.41	188	277400	20.00	ng	0.00
75) Chrysene-d12	21.69	240	290550	20.00	ng	0.00
86) Perylene-d12	24.89	264	299107	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	329582	150.60	ng	0.00
7) Phenol-d6	7.21	99	465865	137.59	ng	0.00
23) Nitrobenzene-d5	9.21	82	301370	94.22	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	193942	147.29	ng	0.00
44) 2-Fluorobiphenyl	13.29	172	657685	89.01	ng	-0.01
78) Terphenyl-d14	20.02	244	1139947	103.17	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	40747	40.689	ng	96
3) Pyridine	3.92	79	116586	40.320	ng	92
4) n-Nitrosodimethylamine	3.84	42	69569	54.132	ng	# 92
6) Aniline	7.37	93	161238	36.699	ng	97
8) 2-Chlorophenol	7.61	128	130405	51.318	ng	93
9) Benzaldehyde	7.18	77	82221	36.748	ng	98
10) Phenol	7.23	94	177137	50.690	ng	92
11) bis(2-Chloroethyl)ether	7.46	93	133690	41.358	ng	99
12) 1,3-Dichlorobenzene	7.93	146	134311	43.112	ng	96
13) 1,4-Dichlorobenzene	8.08	146	137326	43.074	ng	99
14) 1,2-Dichlorobenzene	8.39	146	132230	42.799	ng	97
15) Benzyl Alcohol	8.28	79	122727	44.670	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.57	45	270464	43.582	ng	97
17) 2-Methylphenol	8.48	107	116885	48.018	ng	98
18) Hexachloroethane	9.12	117	53368	45.488	ng	97
19) n-Nitroso-di-n-propylamine	8.84	70	113553	40.313	ng	96
20) 3+4-Methylphenols	8.81	107	162601	48.017	ng	98
22) Acetophenone	8.86	105	197045	48.951	ng	98
24) Nitrobenzene	9.25	77	169073	49.496	ng	97
25) Isophorone	9.76	82	318323	54.463	ng	100
26) 2-Nitrophenol	9.96	139	72814	53.478	ng	96
27) 2,4-Dimethylphenol	10.02	122	124642	58.768	ng	98
28) bis(2-Chloroethoxy)methane	10.25	93	180524	50.055	ng	97
29) 2,4-Dichlorophenol	10.49	162	135861	55.259	ng	95
30) 1,2,4-Trichlorobenzene	10.71	180	141144	45.873	ng	98
31) Naphthalene	10.90	128	412290	50.586	ng	100
32) Benzoic acid	10.15	122	44493m	29.880	ng	
33) 4-Chloroaniline	11.01	127	102743	27.726	ng	96
34) Hexachlorobutadiene	11.18	225	96923	45.551	ng	96
35) Caprolactam	11.79	113	51392m	50.786	ng	
36) 4-Chloro-3-methylphenol	12.13	107	162268	55.313	ng	98
37) 2-Methylnaphthalene	12.50	142	317456	51.141	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.87	216	181164	47.197	ng	98
40) Hexachlorocyclopentadiene	12.84	237	212988	107.889	ng	97

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092618\
 Data File : BG036993.D
 Acq On : 26 Sep 2018 22:25
 Operator : JU/SJ
 Sample : PB113238BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113238BS

Manual Integrations
 APPROVED

Sohil
 9/27/2018 1:37:45 PM

Quant Time: Sep 27 08:41:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 24 10:41:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	116445	54.630	nq	95
43) 2,4,5-Trichlorophenol	13.18	196	129056	56.781	nq	97
45) 1,1'-Biphenyl	13.50	154	418947	49.705	nq	99
46) 2-Chloronaphthalene	13.55	162	318859	48.105	nq	99
47) 2-Nitroaniline	13.75	65	123070	53.680	nq	94
48) Acenaphthylene	14.39	152	543558	52.332	nq	99
49) Dimethylphthalate	14.12	163	429653	48.501	nq	100
50) 2,6-Dinitrotoluene	14.24	165	96003	49.671	nq	99
51) Acenaphthene	14.73	154	311092	49.557	nq	99
52) 3-Nitroaniline	14.58	138	76317	37.471	nq	99
53) 2,4-Dinitrophenol	14.78	184	34661	42.256	nq #	86
54) Dibenzofuran	15.07	168	513929	48.401	nq	99
55) 4-Nitrophenol	14.88	139	144803	111.291	nq	93
56) 2,4-Dinitrotoluene	15.03	165	137780	51.087	nq	97
57) Fluorene	15.72	166	416090	52.428	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.29	232	131773	57.151	nq	98
59) Diethylphthalate	15.48	149	438796	49.111	nq	99
60) 4-Chlorophenyl-phenylether	15.70	204	231252	46.011	nq	99
61) 4-Nitroaniline	15.74	138	107659	50.254	nq	96
62) Azobenzene	16.00	77	443434	51.652	nq	99
64) 4,6-Dinitro-2-methylphenol	15.80	198	51732	35.670	nq	98
65) n-Nitrosodiphenylamine	15.92	169	380262	49.654	nq	98
66) 4-Bromophenyl-phenylether	16.60	248	150510	47.701	nq	98
67) Hexachlorobenzene	16.72	284	151957	46.760	nq	99
68) Atrazine	16.87	200	162760	52.488	nq	98
69) Pentachlorophenol	17.06	266	160631	78.509	nq	96
70) Phenanthrene	17.45	178	706167	52.826	nq	99
71) Anthracene	17.55	178	717882	54.310	nq	99
72) Carbazole	17.82	167	649442	50.392	nq	98
73) Di-n-butylphthalate	18.37	149	746321	52.145	nq	99
74) Fluoranthene	19.46	202	874760	54.363	nq	99
76) Benzidine	19.65	184	418933	47.697	nq	98
77) Pyrene	19.82	202	861889	49.433	nq	99
79) Butylbenzylphthalate	20.70	149	357935	51.504	nq	100
80) Benzo(a)anthracene	21.66	228	869562	51.269	nq	99
81) 3,3'-Dichlorobenzidine	21.58	252	234345	36.299	nq	97
82) Chrysene	21.73	228	827091	50.471	nq	99
83) Bis(2-ethylhexyl)phthalate	21.56	149	490728	50.546	nq	99
84) Di-n-octyl phthalate	22.77	149	831009	51.988	nq	100
85) Indeno(1,2,3-cd)pyrene	28.54	276	976086	55.467	nq	99
87) Benzo(b)fluoranthene	23.87	252	860282	53.970	nq	97
88) Benzo(k)fluoranthene	23.94	252	824258	51.541	nq	99
89) Benzo(a)pyrene	24.74	252	831624	53.965	nq	100
90) Dibenzo(a,h)anthracene	28.60	278	774793	53.645	nq	95
91) Benzo(g,h,i)perylene	29.68	276	807127	55.683	nq	98

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092618\
 Data File : BG036993.D
 Acq On : 26 Sep 2018 22:25
 Operator : JU/SJ
 Sample : PB113238BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113238BS

Manual Integrations
 APPROVED

Sohil
 9/27/2018 1:37:45 PM

Quant Time: Sep 27 08:41:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 24 10:41:23 2018
 Response via : Initial Calibration

