

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092118\
 Data File : BG036878.D
 Acq On : 22 Sep 2018 3:50
 Operator : JU/SJ
 Sample : PB113112BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113112BS

Manual Integrations
 APPROVED

Sohil
 9/24/2018 9:29:51 AM

Quant Time: Sep 22 05:19:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 20 20:07:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	43477	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	193452	20.00	ng	0.00
38) Acenaphthene-d10	14.68	164	135357	20.00	ng	0.00
63) Phenanthrene-d10	17.42	188	353094	20.00	ng	0.00
75) Chrysene-d12	21.69	240	359702	20.00	ng	0.00
86) Perylene-d12	24.90	264	376148	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	418561	184.63	ng	0.00
7) Phenol-d6	7.21	99	597904	170.46	ng	0.00
23) Nitrobenzene-d5	9.21	82	364002	100.76	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	271637	167.73	ng	0.00
44) 2-Fluorobiphenyl	13.30	172	887510	97.66	ng	0.00
78) Terphenyl-d14	20.03	244	1521191	111.20	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	41172	39.689	ng	96
3) Pyridine	3.92	79	115309	38.497	ng	99
4) n-Nitrosodimethylamine	3.84	42	61451	46.159	ng	# 98
6) Aniline	7.37	93	171461	37.674	ng	99
8) 2-Chlorophenol	7.61	128	142223	54.030	ng	99
9) Benzaldehyde	7.19	77	99648	42.994	ng	94
10) Phenol	7.24	94	197032	54.430	ng	93
11) bis(2-Chloroethyl)ether	7.47	93	138279	41.296	ng	100
12) 1,3-Dichlorobenzene	7.94	146	136601	42.328	ng	98
13) 1,4-Dichlorobenzene	8.08	146	139078	42.112	ng	98
14) 1,2-Dichlorobenzene	8.40	146	134633	42.067	ng	98
15) Benzyl Alcohol	8.28	79	134960	47.421	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.57	45	265152	41.246	ng	98
17) 2-Methylphenol	8.49	107	129792	51.474	ng	97
18) Hexachloroethane	9.13	117	51840	42.655	ng	96
19) n-Nitroso-di-n-propylamine	8.85	70	122089	41.842	ng	93
20) 3+4-Methylphenols	8.82	107	181231	51.664	ng	96
22) Acetophenone	8.87	105	214596	47.205	ng	# 99
24) Nitrobenzene	9.25	77	177995	46.139	ng	97
25) Isophorone	9.77	82	348149	52.744	ng	99
26) 2-Nitrophenol	9.96	139	77992	50.720	ng	98
27) 2,4-Dimethylphenol	10.02	122	142416	59.457	ng	98
28) bis(2-Chloroethoxy)methane	10.25	93	201101	49.374	ng	98
29) 2,4-Dichlorophenol	10.50	162	158686	57.150	ng	96
30) 1,2,4-Trichlorobenzene	10.72	180	155034	44.616	ng	98
31) Naphthalene	10.90	128	447077	48.571	ng	100
32) Benzoic acid	10.16	122	59724m	34.418	ng	
33) 4-Chloroaniline	11.02	127	94236	22.517	ng	98
34) Hexachlorobutadiene	11.19	225	102749	42.758	ng	98
35) Caprolactam	11.79	113	62517m	54.703	ng	
36) 4-Chloro-3-methylphenol	12.13	107	180841	54.583	ng	99
37) 2-Methylnaphthalene	12.50	142	359476	51.278	ng	96
39) 1,2,4,5-Tetrachlorobenzene	12.87	216	207184	43.886	ng	99
40) Hexachlorocyclopentadiene	12.85	237	269773	111.108	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	138733	52.919	ng	96
43) 2,4,5-Trichlorophenol	13.18	196	156507	55.987	ng	97
45) 1,1'-Biphenyl	13.51	154	482793	46.573	ng	97
46) 2-Chloronaphthalene	13.55	162	371038	45.513	ng	99
47) 2-Nitroaniline	13.75	65	136222	48.310	ng	97
48) Acenaphthylene	14.40	152	636448	49.821	ng	99
49) Dimethylphthalate	14.13	163	519691	47.699	ng	99
50) 2,6-Dinitrotoluene	14.25	165	116535	49.023	ng	95
51) Acenaphthene	14.74	154	363001	47.016	ng	98
52) 3-Nitroaniline	14.58	138	71507	28.546	ng	94
53) 2,4-Dinitrophenol	14.79	184	60306	56.605	ng	97
54) Dibenzofuran	15.08	168	608169	46.569	ng	98
55) 4-Nitrophenol	14.89	139	181850	113.638	ng	99
56) 2,4-Dinitrotoluene	15.04	165	166228	50.113	ng	93
57) Fluorene	15.72	166	490779	50.279	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.30	232	154770	54.577	ng	97
59) Diethylphthalate	15.49	149	516399	46.992	ng	99
60) 4-Chlorophenyl-phenylether	15.71	204	278674	45.081	ng	99
61) 4-Nitroaniline	15.75	138	124002	47.062	ng	94
62) Azobenzene	16.00	77	509503	48.253	ng	99
64) 4,6-Dinitro-2-methylphenol	15.80	198	66237	35.881	ng	97
65) n-Nitrosodiphenylamine	15.93	169	452595	46.430	ng	97
66) 4-Bromophenyl-phenylether	16.61	248	185200	46.113	ng	98
67) Hexachlorobenzene	16.73	284	186292	45.037	ng	95
68) Atrazine	16.87	200	197221	49.967	ng	99
69) Pentachlorophenol	17.07	266	210902	80.981	ng	99
70) Phenanthrene	17.46	178	834412	49.039	ng	98
71) Anthracene	17.56	178	855507	50.847	ng	100
72) Carbazole	17.83	167	764797	46.621	ng	99
73) Di-n-butylphthalate	18.38	149	879196	48.260	ng	99
74) Fluoranthene	19.47	202	1042057	50.877	ng	99
76) Benzidine	19.65	184	479569	44.104	ng	99
77) Pyrene	19.83	202	1037507	48.066	ng	99
79) Butylbenzylphthalate	20.71	149	411438	47.821	ng	96
80) Benzo(a)anthracene	21.67	228	1034934	49.289	ng	99
81) 3,3'-Dichlorobenzidine	21.59	252	237381	29.701	ng	98
82) Chrysene	21.74	228	962195	47.427	ng	98
83) Bis(2-ethylhexyl)phthalate	21.57	149	561384	46.708	ng	99
84) Di-n-octyl phthalate	22.78	149	953944	48.205	ng	97
85) Indeno(1,2,3-cd)pyrene	28.56	276	1172621	53.825	ng	# 94
87) Benzo(b)fluoranthene	23.88	252	1044213	52.091	ng	98
88) Benzo(k)fluoranthene	23.95	252	969283	48.196	ng	98
89) Benzo(a)pyrene	24.75	252	993344	51.257	ng	100
90) Dibenzo(a,h)anthracene	28.62	278	949194	52.260	ng	97
91) Benzo(g,h,i)perylene	29.71	276	972253	53.337	ng	97

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

