

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG062218\
 Data File : BG035334.D
 Acq On : 22 Jun 2018 15:31
 Operator : JU/SJ
 Sample : PB110230BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB110230BS

Manual Integrations
 APPROVED

Sohil
 6/25/2018 5:19:02 PM

Quant Time: Jun 22 16:56:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG060718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 08 17:16:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	53610	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	195173	20.00	ng	-0.02
38) Acenaphthene-d10	14.68	164	133438	20.00	ng	-0.02
63) Phenanthrene-d10	17.42	188	356041	20.00	ng	-0.02
75) Chrysene-d12	21.69	240	371747	20.00	ng	-0.03
86) Perylene-d12	24.91	264	361589	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	430993	135.67	ng	-0.01
7) Phenol-d6	7.22	99	618168	131.84	ng	0.00
23) Nitrobenzene-d5	9.21	82	390355	95.07	ng	-0.02
41) 2,4,6-Tribromophenol	16.16	330	279155	163.42	ng	-0.02
44) 2-Fluorobiphenyl	13.30	172	904209	88.09	ng	-0.02
78) Terphenyl-d14	20.02	244	1571429	88.54	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.55	88	48950	32.296	ng	# 75
3) Pyridine	3.95	79	140947	34.466	ng	# 67
4) n-Nitrosodimethylamine	3.86	42	56696	32.271	ng	# 70
6) Aniline	7.38	93	169632	27.989	ng	97
8) 2-Chlorophenol	7.62	128	137657	41.341	ng	96
9) Benzaldehyde	7.19	77	95197	26.359	ng	87
10) Phenol	7.24	94	208896	43.052	ng	94
11) bis(2-Chloroethyl)ether	7.48	93	139190	36.959	ng	87
12) 1,3-Dichlorobenzene	7.95	146	157663	39.684	ng	96
13) 1,4-Dichlorobenzene	8.09	146	160045	39.023	ng	99
14) 1,2-Dichlorobenzene	8.41	146	151919	39.435	ng	97
15) Benzyl Alcohol	8.29	79	162319	36.536	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.58	45	153044	40.829	ng	73
17) 2-Methylphenol	8.49	107	138828	43.397	ng	93
18) Hexachloroethane	9.14	117	57253	38.993	ng	89
19) n-Nitroso-di-n-propylamine	8.86	70	126996	31.882	ng	# 86
20) 3+4-Methylphenols	8.82	107	194650	42.451	ng	93
22) Acetophenone	8.87	105	232666	38.484	ng	# 90
24) Nitrobenzene	9.26	77	187997	44.713	ng	94
25) Isophorone	9.78	82	364746	42.075	ng	100
26) 2-Nitrophenol	9.97	139	79630	54.945	ng	# 92
27) 2,4-Dimethylphenol	10.03	122	147110	48.292	ng	91
28) bis(2-Chloroethoxy)methane	10.26	93	192772	41.380	ng	93
29) 2,4-Dichlorophenol	10.50	162	164702	49.689	ng	91
30) 1,2,4-Trichlorobenzene	10.72	180	176701	44.458	ng	98
31) Naphthalene	10.91	128	423510	41.770	ng	98
32) Benzoic acid	10.17	122	93084	45.620	ng	93
33) 4-Chloroaniline	11.01	127	84448	19.182	ng	96
34) Hexachlorobutadiene	11.19	225	133365	43.145	ng	98
35) Caprolactam	11.79	113	52192m	42.606	ng	
36) 4-Chloro-3-methylphenol	12.13	107	193723	46.885	ng	94
37) 2-Methylnaphthalene	12.51	142	332400	43.242	ng	94
39) 1,2,4,5-Tetrachlorobenzene	12.87	216	233034	46.886	ng	99
40) Hexachlorocyclopentadiene	12.85	237	323514	103.796	ng	91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	151907	51.042	nq	94
43) 2,4,5-Trichlorophenol	13.19	196	171547	52.517	nq	96
45) 1,1'-Biphenyl	13.51	154	456960	42.588	nq	98
46) 2-Chloronaphthalene	13.56	162	362582	44.695	nq	98
47) 2-Nitroaniline	13.76	65	120338	50.235	nq	95
48) Acenaphthylene	14.40	152	590635	43.420	nq	99
49) Dimethylphthalate	14.13	163	495448	42.578	nq #	98
50) 2,6-Dinitrotoluene	14.25	165	115114	54.227	nq #	89
51) Acenaphthene	14.74	154	352439	39.655	nq	99
52) 3-Nitroaniline	14.58	138	69008	33.128	nq #	96
53) 2,4-Dinitrophenol	14.78	184	133603	155.511	nq #	63
54) Dibenzofuran	15.07	168	578398	43.453	nq	94
55) 4-Nitrophenol	14.88	139	167273	95.148	nq #	86
56) 2,4-Dinitrotoluene	15.04	165	163077	57.830	nq #	82
57) Fluorene	15.72	166	482462	43.370	nq	96
58) 2,3,4,6-Tetrachlorophenol	15.29	232	173406	54.645	nq	91
59) Diethylphthalate	15.49	149	487962	41.102	nq	97
60) 4-Chlorophenyl-phenylether	15.71	204	281300	44.345	nq	97
61) 4-Nitroaniline	15.74	138	111203	49.778	nq #	76
62) Azobenzene	16.01	77	479766	41.239	nq	94
64) 4,6-Dinitro-2-methylphenol	15.79	198	95376	58.672	nq	82
65) n-Nitrosodiphenylamine	15.93	169	429651	40.183	nq	99
66) 4-Bromophenyl-phenylether	16.60	248	194514	45.367	nq	98
67) Hexachlorobenzene	16.73	284	198000	45.371	nq	93
68) Atrazine	16.88	200	193650	42.470	nq	95
69) Pentachlorophenol	17.07	266	239099	81.478	nq	98
70) Phenanthrene	17.46	178	780005	43.127	nq	96
71) Anthracene	17.55	178	791116	43.668	nq	98
72) Carbazole	17.82	167	723258	43.027	nq	99
73) Di-n-butylphthalate	18.37	149	798674	39.864	nq #	98
74) Fluoranthene	19.47	202	1037261	44.073	nq	97
76) Benzidine	19.64	184	510699	36.926	nq	98
77) Pyrene	19.83	202	1022970	41.742	nq	98
79) Butylbenzylphthalate	20.71	149	364441	40.953	nq #	96
80) Benzo(a)anthracene	21.67	228	1042920	43.902	nq	99
81) 3,3'-Dichlorobenzidine	21.58	252	242281	27.108	nq #	97
82) Chrysene	21.73	228	969280	44.564	nq	98
83) Bis(2-ethylhexyl)phthalate	21.57	149	497102	39.533	nq #	99
84) Di-n-octyl phthalate	22.78	149	852550	39.520	nq #	92
85) Indeno(1,2,3-cd)pyrene	28.58	276	1136105	44.599	nq #	89
87) Benzo(b)fluoranthene	23.89	252	1004910	45.129	nq #	95
88) Benzo(k)fluoranthene	23.95	252	968367	44.456	nq #	96
89) Benzo(a)pyrene	24.76	252	969239	46.257	nq #	96
90) Dibenzo(a,h)anthracene	28.65	278	929192	44.923	nq #	97
91) Benzo(g,h,i)perylene	29.73	276	946749	46.770	nq #	92

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

