

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020819\
 Data File : BG039417.D
 Acq On : 8 Feb 2019 16:00
 Operator : JU/SJ
 Sample : PB116935BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116935BS

Manual Integrations
 APPROVED

Sohil
 2/11/2019 10:13:43 AM

Quant Time: Feb 09 00:02:55 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 29 15:41:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	45432	20.00	ng	0.00
21) Naphthalene-d8	11.00	136	202352	20.00	ng	0.00
39) Acenaphthene-d10	14.80	164	132677	20.00	ng	0.00
64) Phenanthrene-d10	17.54	188	326420	20.00	ng	0.00
76) Chrysene-d12	21.83	240	346902	20.00	ng	-0.02
87) Perylene-d12	25.22	264	364796	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.71	112	438803	165.31	ng	0.00
7) Phenol-d6	7.32	99	615062	157.41	ng	0.00
23) Nitrobenzene-d5	9.35	82	294910	91.05	ng	-0.01
42) 2,4,6-Tribromophenol	16.29	330	235732	165.00	ng	0.00
45) 2-Fluorobiphenyl	13.43	172	726232	78.52	ng	0.00
79) Terphenyl-d14	20.13	244	1305762	79.67	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.60	88	44104	37.983	ng	91
3) Pyridine	4.00	79	121689	39.583	ng	95
4) n-Nitrosodimethylamine	3.93	42	64425	41.903	ng	90
6) Aniline	7.50	93	143826	29.387	ng	98
8) 2-Chlorophenol	7.73	128	139067	47.927	ng	94
9) Benzaldehyde	7.32	77	70031	34.496	ng	96
10) Phenol	7.34	94	185465	45.534	ng	99
11) bis(2-Chloroethyl)ether	7.59	93	129767	40.319	ng	93
12) 1,3-Dichlorobenzene	8.06	146	140978	40.107	ng	96
13) 1,4-Dichlorobenzene	8.21	146	145893	41.641	ng	99
14) 1,2-Dichlorobenzene	8.53	146	137624	40.577	ng	99
15) Benzyl Alcohol	8.41	79	128485	41.885	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.70	45	254219	34.977	ng	100
17) 2-Methylphenol	8.60	107	119475	45.327	ng	98
18) Hexachloroethane	9.26	117	50120	41.581	ng	91
19) n-Nitroso-di-n-propylamine	8.98	70	101219	36.498	ng	93
20) 3+4-Methylphenols	8.93	107	164040	45.194	ng	96
22) Acetophenone	9.00	105	201141	37.005	ng	# 94
24) Nitrobenzene	9.40	77	155314	44.384	ng	97
25) Isophorone	9.91	82	296626	41.325	ng	100
26) 2-Nitrophenol	10.11	139	73272	52.046	ng	96
27) 2,4-Dimethylphenol	10.14	122	142854	49.199	ng	99
28) bis(2-Chloroethoxy)methane	10.39	93	183267	41.452	ng	96
29) 2,4-Dichlorophenol	10.64	162	147592	46.509	ng	95
30) 1,2,4-Trichlorobenzene	10.85	180	148885	41.788	ng	98
31) Naphthalene	11.05	128	424672	42.304	ng	99
32) Benzoic acid	10.25	122	61806	35.418	ng	97
33) 4-Chloroaniline	11.16	127	98698	21.205	ng	100
34) Hexachlorobutadiene	11.31	225	95127	40.148	ng	97
35) Caprolactam	11.93	113	48274m	36.761	ng	
36) 4-Chloro-3-methylphenol	12.25	107	155357	42.331	ng	95
37) 2-Methylnaphthalene	12.64	142	324561	43.652	ng	96
38) 1-Methylnaphthalene	12.86	142	303535	42.351	ng	100
40) 1,2,4,5-Tetrachlorobenzene	13.00	216	174048	41.230	ng	98

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41) Hexachlorocyclopentadiene	12.97	237	230225	100.085	nq	100
43) 2,4,6-Trichlorophenol	13.23	196	118476	45.646	nq	97
44) 2,4,5-Trichlorophenol	13.30	196	128494	47.235	nq	98
46) 1,1'-Biphenyl	13.64	154	431533	39.091	nq	98
47) 2-Chloronaphthalene	13.68	162	332872	40.084	nq	98
48) 2-Nitroaniline	13.89	65	108564	45.939	nq	94
49) Acenaphthylene	14.52	152	534849	42.683	nq	99
50) Dimethylphthalate	14.25	163	427412	39.887	nq	99
51) 2,6-Dinitrotoluene	14.37	165	93689	46.948	nq	99
52) Acenaphthene	14.87	154	325782	40.052	nq	98
53) 3-Nitroaniline	14.71	138	69838	34.211	nq	# 94
54) 2,4-Dinitrophenol	14.91	184	53683	66.231	nq	88
55) Dibenzofuran	15.19	168	531528	40.757	nq	100
56) 4-Nitrophenol	14.99	139	163648	84.142	nq	89
57) 2,4-Dinitrotoluene	15.16	165	134023	51.552	nq	# 86
58) Fluorene	15.84	166	411813	42.359	nq	95
59) 2,3,4,6-Tetrachlorophenol	15.41	232	126004	49.470	nq	# 97
60) Diethylphthalate	15.60	149	436673	39.414	nq	98
61) 4-Chlorophenyl-phenylether	15.83	204	217462	39.504	nq	96
62) 4-Nitroaniline	15.87	138	105419	47.010	nq	90
63) Azobenzene	16.12	77	395989	36.568	nq	97
65) 4,6-Dinitro-2-methylphenol	15.91	198	51279	42.058	nq	98
66) n-Nitrosodiphenylamine	16.05	169	382587	38.547	nq	98
67) 4-Bromophenyl-phenylether	16.73	248	143396	39.598	nq	95
68) Hexachlorobenzene	16.83	284	147089	40.485	nq	95
69) Atrazine	16.99	200	154969	42.725	nq	96
70) Pentachlorophenol	17.18	266	171845	68.053	nq	98
71) Phenanthrene	17.58	178	712398	42.167	nq	99
72) Anthracene	17.68	178	724879	43.559	nq	99
73) Carbazole	17.94	167	641142	38.605	nq	98
74) Di-n-butylphthalate	18.48	149	782530	40.389	nq	99
75) Fluoranthene	19.58	202	882912	45.025	nq	100
77) Benzidine	19.76	184	454875	39.995	nq	99
78) Pyrene	19.95	202	890119	41.218	nq	99
80) Butylbenzylphthalate	20.82	149	368462	40.447	nq	96
81) Benzo(a)anthracene	21.81	228	883775	43.115	nq	99
82) 3,3'-Dichlorobenzidine	21.73	252	190350	25.044	nq	99
83) Chrysene	21.89	228	821617	42.106	nq	99
84) Bis(2-ethylhexyl)phthalate	21.69	149	514510	39.589	nq	100
85) Di-n-octyl phthalate	22.95	149	881741	41.066	nq	96
86) Indeno(1,2,3-cd)pyrene	29.12	276	968816	46.276	nq	# 95
88) Benzo(b)fluoranthene	24.13	252	871993	43.831	nq	99
89) Benzo(k)fluoranthene	24.21	252	841412	42.126	nq	99
90) Benzo(a)pyrene	25.06	252	845162	44.111	nq	97
91) Dibenzo(a,h)anthracene	29.19	278	779242	43.429	nq	99
92) Benzo(g,h,i)perylene	30.34	276	780568	43.645	nq	97

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

