

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031919\
 Data File : BG040006.D
 Acq On : 19 Mar 2019 17:19
 Operator : JU/SJ
 Sample : PB118005BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB118005BS

Manual Integrations
 APPROVED

Sohil
 3/20/2019 4:59:13 PM

Quant Time: Mar 20 04:31:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 04 14:38:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	47121	20.00	ng	-0.03
21) Naphthalene-d8	10.96	136	203092	20.00	ng	-0.03
39) Acenaphthene-d10	14.77	164	133036	20.00	ng	-0.02
64) Phenanthrene-d10	17.52	188	311621	20.00	ng	-0.01
76) Chrysene-d12	21.81	240	326200	20.00	ng	-0.02
87) Perylene-d12	25.17	264	320829	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.69	112	381721	138.20	ng	-0.01
7) Phenol-d6	7.30	99	520132	126.29	ng	-0.01
23) Nitrobenzene-d5	9.32	82	305351	81.32	ng	-0.02
42) 2,4,6-Tribromophenol	16.26	330	193666	124.89	ng	-0.01
45) 2-Fluorobiphenyl	13.40	172	704354	75.38	ng	-0.02
79) Terphenyl-d14	20.11	244	1110896	74.98	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.57	88	40073	31.426	ng	97
3) Pyridine	3.98	79	89728	26.284	ng	95
4) n-Nitrosodimethylamine	3.90	42	65682	40.557	ng	# 98
6) Aniline	7.47	93	170066	31.519	ng	98
8) 2-Chlorophenol	7.71	128	130798	39.034	ng	99
9) Benzaldehyde	7.29	77	74856	28.177	ng	99
10) Phenol	7.33	94	177975	39.891	ng	98
11) bis(2-Chloroethyl)ether	7.56	93	128359	36.900	ng	97
12) 1,3-Dichlorobenzene	8.03	146	136910	36.126	ng	97
13) 1,4-Dichlorobenzene	8.18	146	140432	36.379	ng	97
14) 1,2-Dichlorobenzene	8.50	146	135607	36.359	ng	96
15) Benzyl Alcohol	8.38	79	126018	38.574	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.67	45	268202	38.551	ng	99
17) 2-Methylphenol	8.58	107	113863	40.024	ng	99
18) Hexachloroethane	9.22	117	50962	36.793	ng	96
19) n-Nitroso-di-n-propylamine	8.95	70	103801	35.017	ng	95
20) 3+4-Methylphenols	8.91	107	151967	38.452	ng	98
22) Acetophenone	8.97	105	198226	36.366	ng	# 98
24) Nitrobenzene	9.37	77	157813	40.060	ng	99
25) Isophorone	9.88	82	295758	37.589	ng	99
26) 2-Nitrophenol	10.07	139	75619	39.757	ng	# 93
27) 2,4-Dimethylphenol	10.12	122	131063	44.306	ng	92
28) bis(2-Chloroethoxy)methane	10.36	93	175874	36.782	ng	96
29) 2,4-Dichlorophenol	10.60	162	140391	42.153	ng	96
30) 1,2,4-Trichlorobenzene	10.82	180	145427	38.804	ng	98
31) Naphthalene	11.02	128	412182	38.332	ng	99
32) Benzoic acid	10.26	122	68471	35.676	ng	92
33) 4-Chloroaniline	11.13	127	102287	22.433	ng	97
34) Hexachlorobutadiene	11.27	225	97432	38.045	ng	94
35) Caprolactam	11.92	113	46472m	36.527	ng	
36) 4-Chloro-3-methylphenol	12.23	107	149853	39.250	ng	99
37) 2-Methylnaphthalene	12.61	142	305267	37.973	ng	99
38) 1-Methylnaphthalene	12.83	142	293484	37.566	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.97	216	168773	37.448	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031919\
 Data File : BG040006.D
 Acq On : 19 Mar 2019 17:19
 Operator : JU/SJ
 Sample : PB118005BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleID :
 PB118005BS

Manual Integrations
 APPROVED

Sohil
 3/20/2019 4:59:13 PM

Quant Time: Mar 20 04:31:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 04 14:38:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.94	237	229259	94.920	nq	98
43) 2,4,6-Trichlorophenol	13.21	196	113848	43.147	nq	99
44) 2,4,5-Trichlorophenol	13.28	196	121908	40.989	nq	97
46) 1,1'-Biphenyl	13.61	154	411662	37.622	nq	98
47) 2-Chloronaphthalene	13.65	162	315726	38.355	nq	98
48) 2-Nitroaniline	13.87	65	109003	41.205	nq	98
49) Acenaphthylene	14.50	152	508904	37.660	nq	98
50) Dimethylphthalate	14.22	163	416858	37.473	nq	98
51) 2,6-Dinitrotoluene	14.35	165	95263	40.395	nq	98
52) Acenaphthene	14.83	154	307069	37.755	nq	99
53) 3-Nitroaniline	14.69	138	66751	28.158	nq #	97
54) 2,4-Dinitrophenol	14.89	184	119324	80.217	nq	97
55) Dibenzofuran	15.17	168	502607	37.665	nq	98
56) 4-Nitrophenol	14.99	139	158530	74.755	nq	91
57) 2,4-Dinitrotoluene	15.13	165	134550	40.604	nq	92
58) Fluorene	15.82	166	391939	37.363	nq	97
59) 2,3,4,6-Tetrachlorophenol	15.39	232	119194	41.036	nq	100
60) Diethylphthalate	15.57	149	424487	38.547	nq	99
61) 4-Chlorophenyl-phenylether	15.80	204	209604	37.444	nq	99
62) 4-Nitroaniline	15.85	138	99025	38.531	nq	97
63) Azobenzene	16.10	77	391914	39.261	nq	98
65) 4,6-Dinitro-2-methylphenol	15.89	198	82781	44.928	nq	97
66) n-Nitrosodiphenylamine	16.02	169	366990	40.679	nq	97
67) 4-Bromophenyl-phenylether	16.70	248	136192	39.045	nq	95
68) Hexachlorobenzene	16.81	284	139120	38.477	nq	98
69) Atrazine	16.96	200	138722	39.071	nq	97
70) Pentachlorophenol	17.16	266	180386	78.997	nq	98
71) Phenanthrene	17.56	178	660216	38.464	nq	99
72) Anthracene	17.65	178	669065	40.000	nq	98
73) Carbazole	17.92	167	590990	39.192	nq	98
74) Di-n-butylphthalate	18.45	149	723267	40.766	nq	99
75) Fluoranthene	19.56	202	802834	39.577	nq	98
77) Benzidine	19.74	184	389887	37.666	nq	100
78) Pyrene	19.93	202	796240	37.564	nq	98
80) Butylbenzylphthalate	20.79	149	349737	42.588	nq	97
81) Benzo(a)anthracene	21.78	228	789864	38.636	nq	99
82) 3,3'-Dichlorobenzidine	21.70	252	205729	27.563	nq	98
83) Chrysene	21.85	228	751693	37.926	nq	99
84) Bis(2-ethylhexyl)phthalate	21.65	149	489076	42.381	nq	98
85) Di-n-octyl phthalate	22.90	149	833594	43.626	nq	96
86) Indeno(1,2,3-cd)pyrene	29.03	276	905495	40.272	nq	98
88) Benzo(b)fluoranthene	24.09	252	801577m	41.898	nq	
89) Benzo(k)fluoranthene	24.16	252	750995	42.170	nq	99
90) Benzo(a)pyrene	25.01	252	772734	43.710	nq	98
91) Dibenzo(a,h)anthracene	29.10	278	730729	42.162	nq	98
92) Benzo(g,h,i)perylene	30.27	276	750375	43.109	nq	95

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

