

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040919\
 Data File : BG040400.D
 Acq On : 9 Apr 2019 15:44
 Operator : JU/SJ
 Sample : K2317-16MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleID :
 COMP-3MS

Manual Integrations
 APPROVED

Sohil
 4/10/2019 11:22:52 AM

Quant Time: Apr 10 04:03:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	48612	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	192849	20.00	ng	-0.02
39) Acenaphthene-d10	14.68	164	121802	20.00	ng	-0.02
64) Phenanthrene-d10	17.43	188	331580	20.00	ng	-0.02
76) Chrysene-d12	21.70	240	361939	20.00	ng	-0.03
87) Perylene-d12	24.98	264	381967	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	414222	141.65	ng	-0.02
7) Phenol-d6	7.21	99	568097	140.57	ng	-0.02
23) Nitrobenzene-d5	9.22	82	328308	90.49	ng	-0.02
42) 2,4,6-Tribromophenol	16.17	330	218632	166.09	ng	-0.02
45) 2-Fluorobiphenyl	13.30	172	704643	85.72	ng	-0.02
79) Terphenyl-d14	20.03	244	1288630	80.10	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	61154	44.476	ng	97
3) Pyridine	3.90	79	144280	38.972	ng	96
4) n-Nitrosodimethylamine	3.82	42	74040	43.235	ng	# 91
6) Aniline	7.38	93	175617	33.448	ng	96
8) 2-Chlorophenol	7.61	128	147742	43.161	ng	96
9) Benzaldehyde	7.19	77	72871	26.287	ng	97
10) Phenol	7.23	94	210032	47.882	ng	97
11) bis(2-Chloroethyl)ether	7.47	93	145273	41.350	ng	97
12) 1,3-Dichlorobenzene	7.94	146	159515	42.868	ng	97
13) 1,4-Dichlorobenzene	8.08	146	162422	43.826	ng	100
14) 1,2-Dichlorobenzene	8.40	146	154956	43.722	ng	99
15) Benzyl Alcohol	8.29	79	132985	40.860	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.56	45	276062	40.942	ng	99
17) 2-Methylphenol	8.49	107	135371	44.598	ng	97
18) Hexachloroethane	9.12	117	58977	41.836	ng	97
19) n-Nitroso-di-n-propylamine	8.85	70	115242	39.773	ng	97
20) 3+4-Methylphenols	8.82	107	181256	44.080	ng	96
22) Acetophenone	8.88	105	227330	47.256	ng	# 98
24) Nitrobenzene	9.26	77	171390	45.618	ng	99
25) Isophorone	9.78	82	330778	44.309	ng	99
26) 2-Nitrophenol	9.97	139	80758	45.363	ng	95
27) 2,4-Dimethylphenol	10.02	122	150741	53.307	ng	98
28) bis(2-Chloroethoxy)methane	10.26	93	199722	45.221	ng	97
29) 2,4-Dichlorophenol	10.50	162	150028	48.879	ng	98
30) 1,2,4-Trichlorobenzene	10.72	180	155004	45.991	ng	99
31) Naphthalene	10.92	128	467676	47.312	ng	99
33) 4-Chloroaniline	11.03	127	154618	36.670	ng	97
34) Hexachlorobutadiene	11.17	225	91994	42.679	ng	98
35) Caprolactam	11.81	113	54724m	44.927	ng	
36) 4-Chloro-3-methylphenol	12.14	107	179490	50.867	ng	99
37) 2-Methylnaphthalene	12.51	142	345226	47.222	ng	99
38) 1-Methylnaphthalene	12.73	142	331600	46.775	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	175013	45.208	ng	99
41) Hexachlorocyclopentadiene	12.84	237	205286	96.577	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,6-Trichlorophenol	13.12	196	123396	49.439	ng	96
44) 2,4,5-Trichlorophenol	13.19	196	139896	51.423	ng	98
46) 1,1'-Biphenyl	13.51	154	448760	46.630	ng	100
47) 2-Chloronaphthalene	13.56	162	350023	47.415	ng	100
48) 2-Nitroaniline	13.78	65	123961	49.782	ng	98
49) Acenaphthylene	14.41	152	586028	46.821	ng	99
50) Dimethylphthalate	14.13	163	569910	59.785	ng	99
51) 2,6-Dinitrotoluene	14.26	165	106715	49.857	ng	97
52) Acenaphthene	14.75	154	343543	47.900	ng	99
53) 3-Nitroaniline	14.60	138	97501	43.033	ng	# 92
54) 2,4-Dinitrophenol	14.80	184	78760	69.189	ng	89
55) Dibenzofuran	15.07	168	573583	49.771	ng	99
56) 4-Nitrophenol	14.90	139	192106	105.055	ng	97
57) 2,4-Dinitrotoluene	15.05	165	156276	52.545	ng	99
58) Fluorene	15.73	166	453803	50.085	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.30	232	136025	55.099	ng	99
60) Diethylphthalate	15.48	149	499668	51.223	ng	99
61) 4-Chlorophenyl-phenylether	15.71	204	240158	49.586	ng	95
62) 4-Nitroaniline	15.76	138	130764	53.779	ng	98
63) Azobenzene	16.00	77	461780	50.187	ng	99
65) 4,6-Dinitro-2-methylphenol	15.80	198	82230	44.142	ng	99
66) n-Nitrosodiphenylamine	15.93	169	438549	45.197	ng	98
67) 4-Bromophenyl-phenylether	16.61	248	165274	44.292	ng	97
68) Hexachlorobenzene	16.72	284	168819	44.997	ng	97
69) Atrazine	16.87	200	171584	47.538	ng	98
70) Pentachlorophenol	17.07	266	189705	87.460	ng	98
71) Phenanthrene	17.47	178	820301	47.067	ng	98
72) Anthracene	17.56	178	841771	48.254	ng	99
73) Carbazole	17.84	167	812716	49.228	ng	100
74) Di-n-butylphthalate	18.37	149	978161	49.985	ng	99
75) Fluoranthene	19.48	202	1034228	50.114	ng	99
77) Benzidine	19.66	184	420599	32.180	ng	98
78) Pyrene	19.84	202	1046032	43.948	ng	98
80) Butylbenzylphthalate	20.70	149	484329	47.553	ng	98
81) Benzo(a)anthracene	21.68	228	986417	44.247	ng	99
82) 3,3'-Dichlorobenzidine	21.60	252	253336	29.873	ng	98
83) Chrysene	21.75	228	971660	45.351	ng	100
84) Bis(2-ethylhexyl)phthalate	21.55	149	688798	47.310	ng	99
85) Di-n-octyl phthalate	22.77	149	1175370	47.384	ng	99
86) Indeno(1,2,3-cd)pyrene	28.75	276	1233161	47.059	ng	99
88) Benzo(b)fluoranthene	23.93	252	1055551	45.137	ng	98
89) Benzo(k)fluoranthene	24.00	252	1028707	47.834	ng	99
90) Benzo(a)pyrene	24.83	252	1054049	48.039	ng	99
91) Dibenzo(a,h)anthracene	28.81	278	989529	48.558	ng	99
92) Benzo(g,h,i)perylene	29.93	276	1030526	49.206	ng	98

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

