

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040319\
 Data File : BG040338.D
 Acq On : 3 Apr 2019 23:09
 Operator : JU/SJ
 Sample : K1695-08MSD
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DPC-91723-SO-TP5-BMSD

Manual Integrations
 APPROVED

sohil
 4/4/2019 4:47:02 AM

Quant Time: Apr 04 03:16:40 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	62464	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	243782	20.00	ng	-0.02
39) Acenaphthene-d10	14.68	164	153727	20.00	ng	-0.02
64) Phenanthrene-d10	17.43	188	381993	20.00	ng	-0.02
76) Chrysene-d12	21.70	240	390790	20.00	ng	-0.03
87) Perylene-d12	24.98	264	411250	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	628421	167.25	ng	-0.01
7) Phenol-d6	7.21	99	884733	170.38	ng	-0.01
23) Nitrobenzene-d5	9.22	82	532168	116.03	ng	-0.02
42) 2,4,6-Tribromophenol	16.18	330	345296	207.84	ng	-0.01
45) 2-Fluorobiphenyl	13.30	172	1134471	109.35	ng	-0.02
79) Terphenyl-d14	20.02	244	1763890	101.54	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.50	88	73805	41.773	ng	98
3) Pyridine	3.90	79	198213	41.668	ng	95
4) n-Nitrosodimethylamine	3.82	42	108518	49.316	ng	# 83
6) Aniline	7.38	93	171469	25.416	ng	98
8) 2-Chlorophenol	7.61	128	229152	52.098	ng	97
9) Benzaldehyde	7.19	77	154440	43.358	ng	94
10) Phenol	7.23	94	309911	54.984	ng	99
11) bis(2-Chloroethyl)ether	7.47	93	226438	50.159	ng	95
12) 1,3-Dichlorobenzene	7.93	146	236957	49.558	ng	94
13) 1,4-Dichlorobenzene	8.08	146	241219	50.653	ng	99
14) 1,2-Dichlorobenzene	8.40	146	229407	50.374	ng	98
15) Benzyl Alcohol	8.29	79	219961	52.597	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.57	45	422370	48.750	ng	98
17) 2-Methylphenol	8.48	107	215268	55.193	ng	96
18) Hexachloroethane	9.12	117	88452	48.830	ng	98
19) n-Nitroso-di-n-propylamine	8.85	70	185073	49.709	ng	99
20) 3+4-Methylphenols	8.82	107	294665	55.769	ng	99
22) Acetophenone	8.88	105	351610	57.820	ng	99
24) Nitrobenzene	9.27	77	264423	55.676	ng	99
25) Isophorone	9.78	82	533609	56.545	ng	97
26) 2-Nitrophenol	9.98	139	130339	57.917	ng	100
27) 2,4-Dimethylphenol	10.02	122	244008	68.261	ng	97
28) bis(2-Chloroethoxy)methane	10.26	93	318199	56.993	ng	99
29) 2,4-Dichlorophenol	10.51	162	247621	63.819	ng	99
30) 1,2,4-Trichlorobenzene	10.72	180	247546	58.104	ng	99
31) Naphthalene	10.91	128	712430	57.014	ng	99
32) Benzoic acid	10.15	122	74457	34.474	ng	93
33) 4-Chloroaniline	11.03	127	76659	14.382	ng	96
34) Hexachlorobutadiene	11.18	225	148630	54.548	ng	97
35) Caprolactam	11.83	113	96782m	62.855	ng	
36) 4-Chloro-3-methylphenol	12.14	107	284359	63.749	ng	96
37) 2-Methylnaphthalene	12.51	142	539861	58.417	ng	98
38) 1-Methylnaphthalene	12.73	142	520954	58.132	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	281137	57.539	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.84	237	326418	121.673	ng	97
43) 2,4,6-Trichlorophenol	13.11	196	208032	66.039	ng	97
44) 2,4,5-Trichlorophenol	13.19	196	230048	66.999	ng	99
46) 1,1'-Biphenyl	13.51	154	697267	57.405	ng	100
47) 2-Chloronaphthalene	13.56	162	538874	57.837	ng	98
48) 2-Nitroaniline	13.77	65	196213	62.434	ng	95
49) Acenaphthylene	14.41	152	905462	57.319	ng	98
50) Dimethylphthalate	14.13	163	803552	66.789	ng	99
51) 2,6-Dinitrotoluene	14.27	165	165705	61.339	ng	99
52) Acenaphthene	14.75	154	522592	57.733	ng	99
53) 3-Nitroaniline	14.59	138	62729	21.937	ng	96
54) 2,4-Dinitrophenol	14.81	184	205301	142.898	ng	97
55) Dibenzofuran	15.08	168	863822	59.389	ng	99
56) 4-Nitrophenol	14.91	139	309268	134.003	ng	97
57) 2,4-Dinitrotoluene	15.05	165	239292	63.748	ng	100
58) Fluorene	15.73	166	673212	58.870	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.30	232	217906	69.936	ng	99
60) Diethylphthalate	15.49	149	757796	61.552	ng	98
61) 4-Chlorophenyl-phenylether	15.71	204	365041	59.719	ng	96
62) 4-Nitroaniline	15.76	138	189069	61.610	ng	99
63) Azobenzene	16.00	77	689023	59.332	ng	98
65) 4,6-Dinitro-2-methylphenol	15.81	198	142465	66.383	ng	95
66) n-Nitrosodiphenylamine	15.93	169	654069	58.513	ng	98
67) 4-Bromophenyl-phenylether	16.60	248	241710	56.228	ng	94
68) Hexachlorobenzene	16.72	284	251797	58.257	ng	97
69) Atrazine	16.87	200	247459	59.511	ng	98
70) Pentachlorophenol	17.07	266	318744	127.557	ng	96
71) Phenanthrene	17.47	178	1146878	57.120	ng	98
72) Anthracene	17.56	178	1182104	58.821	ng	99
73) Carbazole	17.83	167	1137367	59.801	ng	98
74) Di-n-butylphthalate	18.37	149	1321182	58.603	ng	99
75) Fluoranthene	19.47	202	1423936	59.892	ng	98
77) Benzidine	19.66	184	545489	38.655	ng	99
78) Pyrene	19.84	202	1427809	55.559	ng	97
80) Butylbenzylphthalate	20.71	149	659245	59.948	ng	98
81) Benzo(a)anthracene	21.68	228	1336396	55.520	ng	96
82) 3,3'-Dichlorobenzidine	21.59	252	248766	27.168	ng	96
83) Chrysene	21.75	228	1316395	56.905	ng	98
84) Bis(2-ethylhexyl)phthalate	21.55	149	930030	59.163	ng	99
85) Di-n-octyl phthalate	22.77	149	1579553	58.977	ng	100
86) Indeno(1,2,3-cd)pyrene	28.75	276	1699182	60.056	ng	100
88) Benzo(b)fluoranthene	23.93	252	1459563	57.969	ng	97
89) Benzo(k)fluoranthene	24.00	252	1385988	59.859	ng	99
90) Benzo(a)pyrene	24.83	252	1427472	60.425	ng	99
91) Dibenzo(a,h)anthracene	28.81	278	1377002	62.760	ng	97
92) Benzo(g,h,i)perylene	29.94	276	1420266	62.987	ng	98

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

