

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090420\
 Data File : BG046541.D
 Acq On : 4 Sep 2020 15:16
 Operator : CG/JU
 Sample : L3894-09MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 AL-FALAH-5MS

Manual Integrations
 APPROVED

mohammad
 9/8/2020 1:45:13 PM

Quant Time: Sep 04 16:59:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 15:06:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	67197	20.00	ng	0.00
21) Naphthalene-d8	10.73	136	278267	20.00	ng	0.00
39) Acenaphthene-d10	14.56	164	201690	20.00	ng	0.00
64) Phenanthrene-d10	17.31	188	503312	20.00	ng	0.00
76) Chrysene-d12	21.56	240	483080	20.00	ng	0.00
86) Perylene-d12	24.65	264	491373	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	320893	84.58	ng	0.00
7) Phenol-d6	7.10	99	425253	79.07	ng	0.00
23) Nitrobenzene-d5	9.08	82	273667	56.21	ng	0.00
42) 2,4,6-Tribromophenol	16.05	330	216896	79.37	ng	0.00
45) 2-Fluorobiphenyl	13.18	172	735424	55.66	ng	0.00
79) Terphenyl-d14	19.92	244	1255741	55.32	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.42	88	61781	39.142	ng	98
3) Pyridine	3.81	79	187399	40.579	ng	97
4) n-Nitrosodimethylamine	3.72	42	89432	52.506	ng	100
6) Aniline	7.25	93	307416	50.795	ng	99
8) 2-Chlorophenol	7.50	128	226399	56.461	ng	99
9) Benzaldehyde	7.07	77	142864	47.412	ng	95
10) Phenol	7.13	94	285335	57.601	ng	97
11) bis(2-Chloroethyl)ether	7.35	93	208815	52.443	ng	99
12) 1,3-Dichlorobenzene	7.82	146	250038	49.115	ng	99
13) 1,4-Dichlorobenzene	7.97	146	251476	49.341	ng	99
14) 1,2-Dichlorobenzene	8.28	146	244695	50.071	ng	99
15) Benzyl Alcohol	8.17	79	199426	57.755	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.46	45	314646	50.945	ng	99
17) 2-Methylphenol	8.38	107	199973	56.922	ng	99
18) Hexachloroethane	9.01	117	85174	49.288	ng	98
19) n-Nitroso-di-n-propylamine	8.74	70	172964	53.765	ng	98
20) 3+4-Methylphenols	8.71	107	278518	57.438	ng	99
22) Acetophenone	8.75	105	323718	47.791	ng	99
24) Nitrobenzene	9.13	77	245580	51.057	ng	99
25) Isophorone	9.65	82	479838	53.757	ng	99
26) 2-Nitrophenol	9.84	139	138069	54.406	ng	96
27) 2,4-Dimethylphenol	9.90	122	223496	64.032	ng	98
28) bis(2-Chloroethoxy)methane	10.13	93	287324	50.011	ng	100
29) 2,4-Dichlorophenol	10.38	162	255907	54.998	ng	97
30) 1,2,4-Trichlorobenzene	10.59	180	262880	47.616	ng	100
31) Naphthalene	10.77	128	678123	49.911	ng	99
32) Benzoic acid	10.07	122	135355	50.889	ng	96
33) 4-Chloroaniline	10.88	127	230762	38.517	ng	99
34) Hexachlorobutadiene	11.07	225	165260	46.112	ng	98
35) Caprolactam	11.67	113	98557m	56.665	ng	
36) 4-Chloro-3-methylphenol	12.02	107	265452	58.329	ng	96
37) 2-Methylnaphthalene	12.38	142	548224	51.162	ng	99
38) 1-Methylnaphthalene	12.60	142	515255	50.764	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.76	216	323000	48.564	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.74	237	382049	132.105	ng	98
43) 2,4,6-Trichlorophenol	12.99	196	239190	53.292	ng	99
44) 2,4,5-Trichlorophenol	13.07	196	262382	55.640	ng	98
46) 1,1'-Biphenyl	13.39	154	719147	51.249	ng	99
47) 2-Chloronaphthalene	13.44	162	568575	47.789	ng	99
48) 2-Nitroaniline	13.64	65	173094	52.394	ng	99
49) Acenaphthylene	14.28	152	926451	51.334	ng	100
50) Dimethylphthalate	14.02	163	880655	56.439	ng	99
51) 2,6-Dinitrotoluene	14.13	165	187566	54.531	ng	97
52) Acenaphthene	14.63	154	559738	50.049	ng	99
53) 3-Nitroaniline	14.46	138	155147	41.641	ng	99
54) 2,4-Dinitrophenol	14.68	184	255054	127.036	ng	97
55) Dibenzofuran	14.96	168	914449	50.949	ng	98
56) 4-Nitrophenol	14.79	139	317199	115.680	ng	95
57) 2,4-Dinitrotoluene	14.92	165	269428	54.935	ng	97
58) Fluorene	15.61	166	781316	51.867	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.19	232	253844	55.433	ng	98
60) Diethylphthalate	15.38	149	802572	52.754	ng	99
61) 4-Chlorophenyl-phenylether	15.60	204	420781	50.719	ng	96
62) 4-Nitroaniline	15.63	138	196107	49.884	ng	97
63) Azobenzene	15.90	77	665508	53.777	ng	97
65) 4,6-Dinitro-2-methylphenol	15.69	198	184393	64.727	ng	95
66) n-Nitrosodiphenylamine	15.82	169	722339	53.198	ng	100
67) 4-Bromophenyl-phenylether	16.50	248	302977	51.597	ng	98
68) Hexachlorobenzene	16.62	284	318108	48.915	ng	98
69) Atrazine	16.77	200	235481	42.511	ng	98
70) Pentachlorophenol	16.96	266	392409	115.490	ng	98
71) Phenanthrene	17.35	178	1282043	50.254	ng	98
72) Anthracene	17.44	178	1312727	52.154	ng	98
73) Carbazole	17.71	167	1211277	50.969	ng	99
74) Di-n-butylphthalate	18.27	149	1357616	49.680	ng	98
75) Fluoranthene	19.36	202	1583357	48.228	ng	99
77) Benzidine	19.53	184	1060808	94.502	ng	97
78) Pyrene	19.72	202	1569878	51.033	ng	98
80) Butylbenzylphthalate	20.61	149	620913	51.747	ng	97
81) Benzo(a)anthracene	21.53	228	1525754	50.672	ng	99
82) 3,3'-Dichlorobenzidine	21.45	252	440861	39.454	ng	98
83) Chrysene	21.60	228	1472764	50.849	ng	99
84) Bis(2-ethylhexyl)phthalate	21.46	149	868983	53.069	ng	99
85) Di-n-octyl phthalate	22.62	149	1493351	53.262	ng	99
87) Indeno(1,2,3-cd)pyrene	28.17	276	1902882	53.919	ng	100
88) Benzo(b)fluoranthene	23.68	252	1651267m	53.819	ng	
89) Benzo(k)fluoranthene	23.75	252	1521624	51.315	ng	99
90) Benzo(a)pyrene	24.52	252	1481617	51.745	ng	99
91) Dibenzo(a,h)anthracene	28.23	278	1544715	54.240	ng	98
92) Benzo(g,h,i)perylene	29.27	276	1592390	55.993	ng	98

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

