

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091719\
 Data File : BG042876.D
 Acq On : 17 Sep 2019 20:40
 Operator : HP/JU
 Sample : K4918-12MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-12MSD

Manual Integrations
 APPROVED

mohammad
 9/18/2019 2:18:42 PM

Quant Time: Sep 18 02:18:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG091219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 12 14:06:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	58065	20.00	ng	-0.01
21) Naphthalene-d8	10.91	136	245408	20.00	ng	-0.02
39) Acenaphthene-d10	14.73	164	194204	20.00	ng	-0.01
64) Phenanthrene-d10	17.47	188	473036	20.00	ng	-0.01
76) Chrysene-d12	21.75	240	368966	20.00	ng	-0.01
87) Perylene-d12	25.01	264	388752	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.71	112	467297	139.92	ng	0.02
7) Phenol-d6	7.31	99	728504	151.97	ng	0.04
23) Nitrobenzene-d5	9.27	82	489713	103.89	ng	0.00
42) 2,4,6-Tribromophenol	16.22	330	427770	165.56	ng	0.00
45) 2-Fluorobiphenyl	13.35	172	1179101	96.11	ng	-0.02
79) Terphenyl-d14	20.07	244	1816418	109.11	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.60	88	46576	30.557	ng	# 90
3) Pyridine	4.05	79	129449	28.207	ng	92
4) n-Nitrosodimethylamine	3.92	42	107306	48.052	ng	# 69
6) Aniline	7.45	93	187183	28.941	ng	99
8) 2-Chlorophenol	7.69	128	194628	53.911	ng	97
9) Benzaldehyde	7.26	77	117881	37.788	ng	93
10) Phenol	7.33	94	279925	56.939	ng	89
11) bis(2-Chloroethyl)ether	7.54	93	171731	45.516	ng	95
12) 1,3-Dichlorobenzene	8.00	146	201010	45.649	ng	98
13) 1,4-Dichlorobenzene	8.15	146	208264	47.242	ng	99
14) 1,2-Dichlorobenzene	8.46	146	192848	45.955	ng	98
15) Benzyl Alcohol	8.37	79	211515	51.414	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.65	45	410940	49.597	ng	98
17) 2-Methylphenol	8.57	107	178429	54.250	ng	96
18) Hexachloroethane	9.19	117	76285	46.870	ng	92
19) n-Nitroso-di-n-propylamine	8.94	70	169277	47.922	ng	93
20) 3+4-Methylphenols	8.90	107	250465	55.244	ng	96
22) Acetophenone	8.95	105	299654	48.348	ng	# 93
24) Nitrobenzene	9.32	77	261464	52.845	ng	99
25) Isophorone	9.88	82	471026	47.940	ng	97
26) 2-Nitrophenol	10.03	139	116156	59.702	ng	# 93
27) 2,4-Dimethylphenol	10.10	122	202399	58.810	ng	98
28) bis(2-Chloroethoxy)methane	10.33	93	264949	48.354	ng	99
29) 2,4-Dichlorophenol	10.57	162	234945	58.746	ng	97
30) 1,2,4-Trichlorobenzene	10.78	180	254868	52.197	ng	99
31) Naphthalene	10.97	128	638106	52.664	ng	98
32) Benzoic acid	10.30	122	152726	54.013	ng	98
33) 4-Chloroaniline	11.08	127	60021	10.598	ng	95
34) Hexachlorobutadiene	11.25	225	182416	53.321	ng	98
35) Caprolactam	11.98	113	75859m	51.007	ng	
36) 4-Chloro-3-methylphenol	12.21	107	253125	58.292	ng	98
37) 2-Methylnaphthalene	12.56	142	501854	55.292	ng	99
38) 1-Methylnaphthalene	12.78	142	476430	55.966	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.93	216	342635	53.392	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.91	237	476202	129.532	nq	98
43) 2,4,6-Trichlorophenol	13.18	196	233876	54.741	nq	95
44) 2,4,5-Trichlorophenol	13.25	196	256164	58.536	nq	98
46) 1,1'-Biphenyl	13.56	154	715701	52.643	nq	98
47) 2-Chloronaphthalene	13.60	162	568377	50.538	nq	98
48) 2-Nitroaniline	13.82	65	200824	51.475	nq	95
49) Acenaphthylene	14.45	152	917304	53.373	nq	98
50) Dimethylphthalate	14.20	163	948510	65.197	nq	98
51) 2,6-Dinitrotoluene	14.30	165	174982	58.617	nq	98
52) Acenaphthene	14.79	154	546577	48.680	nq	97
53) 3-Nitroaniline	14.64	138	69401	20.715	nq #	94
54) 2,4-Dinitrophenol	14.84	184	224122	152.509	nq #	72
55) Dibenzofuran	15.13	168	907624	54.384	nq	98
56) 4-Nitrophenol	14.97	139	284860	109.966	nq	94
57) 2,4-Dinitrotoluene	15.09	165	243041	61.081	nq #	98
58) Fluorene	15.77	166	763403	55.194	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.36	232	266558	62.835	nq	97
60) Diethylphthalate	15.56	149	761608	51.573	nq	98
61) 4-Chlorophenyl-phenylether	15.77	204	450060	55.033	nq	99
62) 4-Nitroaniline	15.80	138	177195	49.067	nq #	82
63) Azobenzene	16.06	77	736160	51.767	nq	96
65) 4,6-Dinitro-2-methylphenol	15.85	198	161583	79.832	nq	96
66) n-Nitrosodiphenylamine	15.98	169	686645	53.943	nq	99
67) 4-Bromophenyl-phenylether	16.66	248	308206	53.831	nq	97
68) Hexachlorobenzene	16.78	284	321566	52.562	nq	97
69) Atrazine	16.94	200	263283	59.396	nq	100
70) Pentachlorophenol	17.12	266	431615	118.926	nq	93
71) Phenanthrene	17.51	178	1194793	52.815	nq	99
72) Anthracene	17.61	178	1211582	53.607	nq	99
73) Carbazole	17.88	167	1088808	50.686	nq	96
74) Di-n-butylphthalate	18.44	149	1193579	46.384	nq	100
75) Fluoranthene	19.52	202	1440041	49.116	nq	98
77) Benzidine	19.70	184	437606	42.451	nq	99
78) Pyrene	19.88	202	1419345	61.444	nq	96
80) Butylbenzylphthalate	20.77	149	480370	51.384	nq	98
81) Benzo(a)anthracene	21.73	228	1227197	52.311	nq	99
82) 3,3'-Dichlorobenzidine	21.64	252	220817	24.221	nq	96
83) Chrysene	21.80	228	1190411	53.571	nq	99
84) Bis(2-ethylhexyl)phthalate	21.65	149	609894	47.746	nq #	98
85) Di-n-octyl phthalate	22.88	149	959874	44.079	nq	97
86) Indeno(1,2,3-cd)pyrene	28.75	276	1284038	46.661	nq #	86
88) Benzo(b)fluoranthene	23.98	252	1197221	52.448	nq #	97
89) Benzo(k)fluoranthene	24.05	252	1145886	53.142	nq #	98
90) Benzo(a)pyrene	24.86	252	1155244	54.049	nq #	99
91) Dibenzo(a,h)anthracene	28.82	278	1088181	51.763	nq #	97
92) Benzo(g,h,i)perylene	29.92	276	1102101	53.747	nq #	93

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

