

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\
 Data File : BG044357.D
 Acq On : 10 Feb 2020 15:24
 Operator : CG/JU
 Sample : PB126601BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB126601BS

Manual Integrations
 APPROVED

mohammad
 2/12/2020 9:49:30 AM

Quant Time: Feb 10 16:20:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 10 16:11:50 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	64763	20.00	ng	0.00
21) Naphthalene-d8	10.70	136	276992	20.00	ng	0.00
39) Acenaphthene-d10	14.55	164	196309	20.00	ng	0.00
64) Phenanthrene-d10	17.29	188	472959	20.00	ng	0.00
76) Chrysene-d12	21.53	240	447777	20.00	ng	0.00
87) Perylene-d12	24.62	264	500905	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	459766	125.29	ng	0.00
7) Phenol-d6	7.08	99	622064	126.24	ng	0.00
23) Nitrobenzene-d5	9.06	82	325894	66.42	ng	0.00
42) 2,4,6-Tribromophenol	16.03	330	284076	123.27	ng	0.00
45) 2-Fluorobiphenyl	13.16	172	788880	67.29	ng	0.00
79) Terphenyl-d14	19.90	244	1458267	66.99	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.36	88	57971	35.161	ng	99
3) Pyridine	3.76	79	139500	31.758	ng	98
4) n-Nitrosodimethylamine	3.67	42	71911	39.177	ng	99
6) Aniline	7.22	93	218551	35.731	ng	98
8) 2-Chlorophenol	7.47	128	184589	45.769	ng	97
9) Benzaldehyde	7.04	77	81975	29.209	ng	96
10) Phenol	7.10	94	227961	46.744	ng	97
11) bis(2-Chloroethyl)ether	7.32	93	170054	40.787	ng	99
12) 1,3-Dichlorobenzene	7.79	146	186881	38.810	ng	99
13) 1,4-Dichlorobenzene	7.94	146	186535	39.112	ng	99
14) 1,2-Dichlorobenzene	8.26	146	178886	39.683	ng	98
15) Benzyl Alcohol	8.14	79	151622	43.461	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.43	45	216510	38.367	ng	98
17) 2-Methylphenol	8.35	107	159916	45.959	ng	98
18) Hexachloroethane	8.99	117	67878	37.928	ng	97
19) n-Nitroso-di-n-propylamine	8.71	70	133681	40.705	ng	98
20) 3+4-Methylphenols	8.68	107	219595	45.794	ng	100
22) Acetophenone	8.72	105	255240	37.879	ng	99
24) Nitrobenzene	9.10	77	190671	39.808	ng	97
25) Isophorone	9.63	82	376363	41.156	ng	99
26) 2-Nitrophenol	9.82	139	105389	42.404	ng	98
27) 2,4-Dimethylphenol	9.89	122	184199	52.503	ng	98
28) bis(2-Chloroethoxy)methane	10.11	93	239503	39.260	ng	98
29) 2,4-Dichlorophenol	10.36	162	190731	44.961	ng	98
30) 1,2,4-Trichlorobenzene	10.57	180	186283	38.296	ng	99
31) Naphthalene	10.76	128	552839	41.137	ng	99
32) Benzoic acid	10.03	122	55426	31.075	ng	91
33) 4-Chloroaniline	10.86	127	136540	22.340	ng	98
34) Hexachlorobutadiene	11.05	225	111045	37.100	ng	99
35) Caprolactam	11.63	113	72871m	46.893	ng	
36) 4-Chloro-3-methylphenol	12.00	107	201626	45.332	ng	100
37) 2-Methylnaphthalene	12.37	142	423791	42.473	ng	99
38) 1-Methylnaphthalene	12.58	142	399271	42.443	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.74	216	208352	35.481	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.72	237	258555	91.762	ng	100
43) 2,4,6-Trichlorophenol	12.98	196	151977	40.040	ng	99
44) 2,4,5-Trichlorophenol	13.05	196	169627	43.753	ng	98
46) 1,1'-Biphenyl	13.38	154	540661	38.182	ng	98
47) 2-Chloronaphthalene	13.42	162	423657	37.599	ng	99
48) 2-Nitroaniline	13.62	65	128503	38.643	ng	99
49) Acenaphthylene	14.26	152	686393	41.532	ng	99
50) Dimethylphthalate	14.00	163	575922	40.813	ng	99
51) 2,6-Dinitrotoluene	14.11	165	130780	41.566	ng	98
52) Acenaphthene	14.60	154	420876	39.289	ng	98
53) 3-Nitroaniline	14.44	138	93038	26.728	ng	98
54) 2,4-Dinitrophenol	14.65	184	153840	82.277	ng	99
55) Dibenzofuran	14.94	168	670581	41.346	ng	98
56) 4-Nitrophenol	14.76	139	257913	101.150	ng	97
57) 2,4-Dinitrotoluene	14.90	165	189092	42.880	ng	99
58) Fluorene	15.59	166	541827	42.415	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.17	232	157239	44.902	ng	99
60) Diethylphthalate	15.37	149	597088	42.465	ng	99
61) 4-Chlorophenyl-phenylether	15.59	204	279714	40.384	ng	100
62) 4-Nitroaniline	15.61	138	141775	40.760	ng	99
63) Azobenzene	15.88	77	517221	41.971	ng	100
65) 4,6-Dinitro-2-methylphenol	15.67	198	117140	44.869	ng	96
66) n-Nitrosodiphenylamine	15.80	169	517995	39.081	ng	98
67) 4-Bromophenyl-phenylether	16.48	248	192922	37.440	ng	99
68) Hexachlorobenzene	16.60	284	210801	36.516	ng	99
69) Atrazine	16.75	200	211402	46.534	ng	100
70) Pentachlorophenol	16.95	266	240442	82.271	ng	95
71) Phenanthrene	17.33	178	934332	40.796	ng	99
72) Anthracene	17.42	178	966751	42.696	ng	98
73) Carbazole	17.69	167	890302	42.652	ng	99
74) Di-n-butylphthalate	18.26	149	1109024	42.994	ng	98
75) Fluoranthene	19.34	202	1164154	43.941	ng	100
77) Benzidine	19.52	184	582849	46.927	ng	99
78) Pyrene	19.70	202	1164695	40.502	ng	99
80) Butylbenzylphthalate	20.60	149	498072	38.007	ng	98
81) Benzo(a)anthracene	21.51	228	1113340	40.343	ng	98
82) 3,3'-Dichlorobenzidine	21.43	252	312647	29.190	ng	99
83) Chrysene	21.58	228	1060130	39.742	ng	100
84) Bis(2-ethylhexyl)phthalate	21.44	149	706890	39.190	ng	100
85) Di-n-octyl phthalate	22.60	149	1238680	39.690	ng	99
86) Indeno(1,2,3-cd)pyrene	28.11	276	1367636	39.298	ng	100
88) Benzo(b)fluoranthene	23.65	252	1170625	40.557	ng	98
89) Benzo(k)fluoranthene	23.71	252	1164382	41.390	ng	99
90) Benzo(a)pyrene	24.47	252	1090389	39.955	ng	99
91) Dibenzo(a,h)anthracene	28.18	278	1126946	41.726	ng	98
92) Benzo(g,h,i)perylene	29.20	276	1149217	42.505	ng	100

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

