

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090718\
 Data File : BG036606.D
 Acq On : 7 Sep 2018 20:55
 Operator : JU/SJ
 Sample : PB112680BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112680BS

Manual Integrations
 APPROVED

Sohil
 9/10/2018 11:05:50 AM

Quant Time: Sep 08 01:42:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 31 13:03:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	52677	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	239020	20.00	ng	0.00
38) Acenaphthene-d10	14.68	164	162183	20.00	ng	0.00
63) Phenanthrene-d10	17.42	188	432345	20.00	ng	0.00
75) Chrysene-d12	21.69	240	429458	20.00	ng	0.00
86) Perylene-d12	24.90	264	454979	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	376280	113.31	ng	0.00
7) Phenol-d6	7.22	99	505940	106.41	ng	0.00
23) Nitrobenzene-d5	9.22	82	295104	66.99	ng	0.00
41) 2,4,6-Tribromophenol	16.17	330	226761	117.18	ng	0.00
44) 2-Fluorobiphenyl	13.30	172	732434	64.48	ng	0.00
78) Terphenyl-d14	20.03	244	1251521	68.11	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	51348	33.133	ng	98
3) Pyridine	3.92	79	149228	34.304	ng	94
4) n-Nitrosodimethylamine	3.84	42	75839	39.142	ng	95
6) Aniline	7.38	93	197080	32.656	ng	96
8) 2-Chlorophenol	7.62	128	161985	44.981	ng	98
9) Benzaldehyde	7.19	77	89883	27.453	ng	97
10) Phenol	7.24	94	223148	44.849	ng	98
11) bis(2-Chloroethyl)ether	7.48	93	154429	39.150	ng	95
12) 1,3-Dichlorobenzene	7.95	146	160360	38.998	ng	99
13) 1,4-Dichlorobenzene	8.09	146	162561	39.156	ng	99
14) 1,2-Dichlorobenzene	8.41	146	158165	39.892	ng	98
15) Benzyl Alcohol	8.29	79	156715	40.888	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.58	45	284739	35.794	ng	99
17) 2-Methylphenol	8.49	107	147906	44.769	ng	98
18) Hexachloroethane	9.14	117	60150	40.410	ng	91
19) n-Nitroso-di-n-propylamine	8.86	70	129498	36.444	ng	97
20) 3+4-Methylphenols	8.82	107	203510	44.121	ng	98
22) Acetophenone	8.87	105	244553	38.963	ng	99
24) Nitrobenzene	9.26	77	192956	40.619	ng	95
25) Isophorone	9.78	82	360583	41.203	ng	99
26) 2-Nitrophenol	9.97	139	88048	46.774	ng	100
27) 2,4-Dimethylphenol	10.03	122	162114	46.516	ng	96
28) bis(2-Chloroethoxy)methane	10.26	93	214050	39.853	ng	98
29) 2,4-Dichlorophenol	10.50	162	174451	45.674	ng	93
30) 1,2,4-Trichlorobenzene	10.73	180	175657	39.313	ng	96
31) Naphthalene	10.91	128	501509	41.973	ng	99
32) Benzoic acid	10.16	122	108603	42.589	ng	93
33) 4-Chloroaniline	11.01	127	142200	25.971	ng	97
34) Hexachlorobutadiene	11.20	225	117402	39.107	ng	94
35) Caprolactam	11.79	113	65884m	43.301	ng	
36) 4-Chloro-3-methylphenol	12.13	107	198572	44.743	ng	99
37) 2-Methylnaphthalene	12.51	142	392524	44.898	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.88	216	226681	39.495	ng	98
40) Hexachlorocyclopentadiene	12.86	237	271332	90.860	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	154246	44.118	nq	97
43) 2,4,5-Trichlorophenol	13.19	196	169746	45.520	nq	97
45) 1,1'-Biphenyl	13.52	154	515649	38.303	nq	100
46) 2-Chloronaphthalene	13.56	162	394217	38.357	nq	98
47) 2-Nitroaniline	13.76	65	143433	44.443	nq	97
48) Acenaphthylene	14.40	152	677474	42.178	nq	99
49) Dimethylphthalate	14.13	163	539060	40.705	nq	98
50) 2,6-Dinitrotoluene	14.25	165	122168	44.831	nq	95
51) Acenaphthene	14.74	154	401689	39.049	nq	100
52) 3-Nitroaniline	14.58	138	105201	35.824	nq	99
53) 2,4-Dinitrophenol	14.79	184	103521	100.292	nq	96
54) Dibenzofuran	15.08	168	649597	40.308	nq	99
55) 4-Nitrophenol	14.89	139	204196	85.044	nq	99
56) 2,4-Dinitrotoluene	15.04	165	174312	49.080	nq	89
57) Fluorene	15.73	166	527395	43.775	nq	96
58) 2,3,4,6-Tetrachlorophenol	15.30	232	173783	49.601	nq	95
59) Diethylphthalate	15.50	149	541692	40.587	nq	100
60) 4-Chlorophenyl-phenylether	15.72	204	299744	40.291	nq	99
61) 4-Nitroaniline	15.75	138	136268	44.344	nq	92
62) Azobenzene	16.01	77	534575	39.366	nq	97
64) 4,6-Dinitro-2-methylphenol	15.80	198	90154	47.469	nq	94
65) n-Nitrosodiphenylamine	15.93	169	483994	37.675	nq	99
66) 4-Bromophenyl-phenylether	16.61	248	202911	40.086	nq	99
67) Hexachlorobenzene	16.73	284	204130	39.438	nq	98
68) Atrazine	16.88	200	204419	42.394	nq	99
69) Pentachlorophenol	17.07	266	253390	72.032	nq	97
70) Phenanthrene	17.47	178	907229	41.296	nq	98
71) Anthracene	17.56	178	918354	41.936	nq	98
72) Carbazole	17.82	167	809334	37.006	nq	99
73) Di-n-butylphthalate	18.38	149	927646	38.090	nq	99
74) Fluoranthene	19.47	202	1098248	41.003	nq	99
76) Benzidine	19.65	184	566677	41.358	nq	99
77) Pyrene	19.83	202	1054951	39.946	nq	98
79) Butylbenzylphthalate	20.72	149	427983	40.721	nq	93
80) Benzo(a)anthracene	21.67	228	1087148	41.785	nq	99
81) 3,3'-Dichlorobenzidine	21.58	252	277992	26.810	nq	99
82) Chrysene	21.74	228	1005710	40.782	nq	99
83) Bis(2-ethylhexyl)phthalate	21.58	149	584579	39.747	nq	100
84) Di-n-octyl phthalate	22.79	149	993268	40.433	nq	97
85) Indeno(1,2,3-cd)pyrene	28.55	276	1201287	41.940	nq	99
87) Benzo(b)fluoranthene	23.89	252	1078290	42.679	nq	98
88) Benzo(k)fluoranthene	23.95	252	1033274	41.862	nq	99
89) Benzo(a)pyrene	24.75	252	1047493	43.146	nq	99
90) Dibenzo(a,h)anthracene	28.62	278	962582	41.070	nq	98
91) Benzo(g,h,i)perylene	29.70	276	996548	42.311	nq #	95

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

