

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG051420\
 Data File : BG045375.D
 Acq On : 14 May 2020 13:11
 Operator : CG/JU
 Sample : TEST
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TEST

Quant Time: May 14 14:38:28 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 14 13:01:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.97	152	70447	20.00	ng	0.00
21) Naphthalene-d8	10.78	136	296499	20.00	ng	0.00
39) Acenaphthene-d10	14.61	164	222044	20.00	ng	0.00
64) Phenanthrene-d10	17.36	188	528916	20.00	ng	0.00
76) Chrysene-d12	21.63	240	479256	20.00	ng	0.00
87) Perylene-d12	24.79	264	534184	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	381953	89.51	ng	0.00
7) Phenol-d6	7.14	99	545584	86.06	ng	0.00
23) Nitrobenzene-d5	9.13	82	342869	52.76	ng	0.00
42) 2,4,6-Tribromophenol	16.11	330	270261	78.79	ng	0.00
45) 2-Fluorobiphenyl	13.23	172	825333	54.50	ng	0.00
79) Terphenyl-d14	19.98	244	1438045	65.40	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.42	88	47825	25.220	ng	95
3) Pyridine	3.82	79	111469	19.091	ng	98
4) n-Nitrosodimethylamine	3.73	42	49852	27.871	ng	# 98
6) Aniline	7.29	93	112339	13.628	ng	99
8) 2-Chlorophenol	7.54	128	141040	30.755	ng	97
9) Benzaldehyde	7.10	77	101324	24.824	ng	95
10) Phenol	7.17	94	191770	30.228	ng	97
11) bis(2-Chloroethyl)ether	7.38	93	127970	25.335	ng	93
12) 1,3-Dichlorobenzene	7.86	146	147384	27.273	ng	96
13) 1,4-Dichlorobenzene	8.00	146	148063	27.497	ng	95
14) 1,2-Dichlorobenzene	8.33	146	139619	27.103	ng	95
15) Benzyl Alcohol	8.20	79	144078	27.106	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.50	45	120672	24.338	ng	97
17) 2-Methylphenol	8.42	107	131444	26.854	ng	96
18) Hexachloroethane	9.06	117	59337	29.313	ng	97
19) n-Nitroso-di-n-propylamine	8.77	70	118498	26.430	ng	97
20) 3+4-Methylphenols	8.75	107	181170	26.443	ng	98
22) Acetophenone	8.79	105	218922	27.304	ng	# 97
24) Nitrobenzene	9.17	77	180603	27.114	ng	94
25) Isophorone	9.70	82	337841	25.388	ng	99
26) 2-Nitrophenol	9.88	139	78435	27.338	ng	95
27) 2,4-Dimethylphenol	9.96	122	138465	30.416	ng	94
28) bis(2-Chloroethoxy)methane	10.18	93	187437	26.808	ng	96
29) 2,4-Dichlorophenol	10.43	162	151318	28.786	ng	96
30) 1,2,4-Trichlorobenzene	10.64	180	159975	26.879	ng	99
31) Naphthalene	10.83	128	431528	26.605	ng	99
32) Benzoic acid	10.10	122	76471	24.406	ng	97
33) 4-Chloroaniline	10.94	127	52742	6.972	ng	99
34) Hexachlorobutadiene	11.12	225	108346	26.552	ng	99
35) Caprolactam	11.69	113	56287m	26.904	ng	
36) 4-Chloro-3-methylphenol	12.07	107	173890	28.159	ng	99
37) 2-Methylnaphthalene	12.44	142	335511	26.650	ng	97
38) 1-Methylnaphthalene	12.66	142	322769	27.157	ng	96
40) 1,2,4,5-Tetrachlorobenzene	12.81	216	193975	27.132	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.79	237	254661	56.992	ng	100
43) 2,4,6-Trichlorophenol	13.05	196	135916	26.937	ng	91
44) 2,4,5-Trichlorophenol	13.13	196	148232	25.810	ng	98
46) 1,1'-Biphenyl	13.44	154	443238	26.263	ng	98
47) 2-Chloronaphthalene	13.49	162	336377	26.085	ng	98
48) 2-Nitroaniline	13.69	65	109635	25.840	ng	98
49) Acenaphthylene	14.34	152	564342	26.867	ng	99
50) Dimethylphthalate	14.07	163	474269	26.232	ng	99
51) 2,6-Dinitrotoluene	14.18	165	106557	26.350	ng	98
52) Acenaphthene	14.68	154	417233m	29.358	ng	
53) 3-Nitroaniline	14.51	138	35609	8.029	ng	# 93
54) 2,4-Dinitrophenol	14.73	184	129839	53.995	ng	92
55) Dibenzofuran	15.01	168	551582	26.621	ng	99
56) 4-Nitrophenol	14.84	139	174529	50.751	ng	93
57) 2,4-Dinitrotoluene	14.97	165	154305	26.110	ng	98
58) Fluorene	15.67	166	459302	27.138	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.25	232	141873	27.762	ng	98
60) Diethylphthalate	15.44	149	491599	26.079	ng	99
61) 4-Chlorophenyl-phenylether	15.65	204	249395	25.601	ng	97
62) 4-Nitroaniline	15.68	138	101518	22.068	ng	96
63) Azobenzene	15.95	77	465121	26.763	ng	96
65) 4,6-Dinitro-2-methylphenol	15.74	198	99142	28.517	ng	93
66) n-Nitrosodiphenylamine	15.87	169	410949	27.042	ng	98
67) 4-Bromophenyl-phenylether	16.55	248	173191	25.382	ng	98
68) Hexachlorobenzene	16.68	284	185701	26.241	ng	99
69) Atrazine	16.82	200	157710	26.487	ng	94
70) Pentachlorophenol	17.02	266	210349	48.453	ng	98
71) Phenanthrene	17.40	178	741998	27.300	ng	99
72) Anthracene	17.50	178	764890	28.055	ng	99
73) Carbazole	17.77	167	711282	25.708	ng	97
74) Di-n-butylphthalate	18.33	149	870475	26.393	ng	99
75) Fluoranthene	19.42	202	920173	26.230	ng	99
77) Benzidine	19.60	184	298101	17.887	ng	99
78) Pyrene	19.78	202	928743	28.110	ng	97
80) Butylbenzylphthalate	20.67	149	371139	27.099	ng	99
81) Benzo(a)anthracene	21.61	228	858242	27.629	ng	98
82) 3,3'-Dichlorobenzidine	21.52	252	141316	11.430	ng	# 95
83) Chrysene	21.67	228	816371	27.353	ng	98
84) Bis(2-ethylhexyl)phthalate	21.52	149	515212	27.352	ng	97
85) Di-n-octyl phthalate	22.71	149	877913	26.953	ng	99
86) Indeno(1,2,3-cd)pyrene	28.38	276	1055291	26.632	ng	# 95
88) Benzo(b)fluoranthene	23.79	252	893108	26.691	ng	99
89) Benzo(k)fluoranthene	23.86	252	862330	27.099	ng	95
90) Benzo(a)pyrene	24.64	252	825231	26.121	ng	96
91) Dibenzo(a,h)anthracene	28.45	278	860143	27.020	ng	98
92) Benzo(g,h,i)perylene	29.50	276	868588	27.215	ng	98

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

