

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061820\
 Data File : BG045800.D
 Acq On : 18 Jun 2020 19:41
 Operator : CG/JU
 Sample : PB129622BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB129622BS

Manual Integrations
 APPROVED

mohammad
 6/19/2020 12:57:35 PM

Quant Time: Jun 19 03:38:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 15 18:34:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	52066	20.00	ng	0.00
21) Naphthalene-d8	10.73	136	205185	20.00	ng	0.00
39) Acenaphthene-d10	14.57	164	145278	20.00	ng	0.00
64) Phenanthrene-d10	17.32	188	343512	20.00	ng	0.00
76) Chrysene-d12	21.58	240	310226	20.00	ng	0.00
86) Perylene-d12	24.71	264	328840	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.53	112	445788	144.63	ng	0.00
7) Phenol-d6	7.16	99	630658	134.52	ng	0.00
23) Nitrobenzene-d5	9.08	82	409820	91.24	ng	0.00
42) 2,4,6-Tribromophenol	16.07	330	302577	137.95	ng	0.00
45) 2-Fluorobiphenyl	13.19	172	915553	92.61	ng	0.00
79) Terphenyl-d14	19.94	244	1496194	97.74	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.37	88	56319	39.895	ng	99
3) Pyridine	3.77	79	164064	39.219	ng	98
4) n-Nitrosodimethylamine	3.68	42	70547	50.241	ng	# 99
6) Aniline	7.25	93	200640	36.346	ng	98
8) 2-Chlorophenol	7.51	128	173533	52.599	ng	95
9) Benzaldehyde	7.06	77	102509	36.403	ng	97
10) Phenol	7.19	94	233760	51.459	ng	98
11) bis(2-Chloroethyl)ether	7.34	93	157981	45.786	ng	95
12) 1,3-Dichlorobenzene	7.81	146	185387	47.153	ng	99
13) 1,4-Dichlorobenzene	7.96	146	190649	48.547	ng	95
14) 1,2-Dichlorobenzene	8.27	146	176507	46.948	ng	98
15) Benzyl Alcohol	8.18	79	178451	49.186	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.45	45	147906	45.483	ng	82
17) 2-Methylphenol	8.42	107	158025	46.743	ng	98
18) Hexachloroethane	9.00	117	73298	49.143	ng	98
19) n-Nitroso-di-n-propylamine	8.73	70	147068	47.714	ng	99
20) 3+4-Methylphenols	8.75	107	206617	46.471	ng	# 68
22) Acetophenone	8.75	105	262616	46.632	ng	# 97
24) Nitrobenzene	9.12	77	222326	46.425	ng	99
25) Isophorone	9.65	82	397935	45.766	ng	98
26) 2-Nitrophenol	9.84	139	96627	48.431	ng	91
27) 2,4-Dimethylphenol	9.93	122	158782	52.126	ng	94
28) bis(2-Chloroethoxy)methane	10.13	93	220761	48.606	ng	98
29) 2,4-Dichlorophenol	10.41	162	177656	49.989	ng	94
30) 1,2,4-Trichlorobenzene	10.59	180	195763	46.530	ng	96
31) Naphthalene	10.77	128	529448	47.197	ng	98
32) Benzoic acid	10.15	122	101900	50.912	ng	# 89
33) 4-Chloroaniline	10.90	127	99587	21.199	ng	97
34) Hexachlorobutadiene	11.07	225	139574	46.647	ng	99
35) Caprolactam	11.67	113	59852m	52.043	ng	
36) 4-Chloro-3-methylphenol	12.07	107	209110	50.961	ng	99
37) 2-Methylnaphthalene	12.39	142	400349	47.927	ng	98
38) 1-Methylnaphthalene	12.61	142	386475	48.891	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.76	216	229208	47.915	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.74	237	268112	99.208	nq	98
43) 2,4,6-Trichlorophenol	13.02	196	159021	49.030	nq	97
44) 2,4,5-Trichlorophenol	13.11	196	164798	44.582	nq	94
46) 1,1'-Biphenyl	13.40	154	508406	48.880	nq	99
47) 2-Chloronaphthalene	13.44	162	396021	47.135	nq	97
48) 2-Nitroaniline	13.66	65	132779	47.510	nq	97
49) Acenaphthylene	14.29	152	645111	47.937	nq	100
50) Dimethylphthalate	14.03	163	542470	47.664	nq	98
51) 2,6-Dinitrotoluene	14.15	165	125350	49.037	nq	97
52) Acenaphthene	14.63	154	387250	47.473	nq	97
53) 3-Nitroaniline	14.49	138	70538	27.627	nq	97
54) 2,4-Dinitrophenol	14.70	184	141865	97.351	nq	97
55) Dibenzofuran	14.97	168	632459	47.727	nq	98
56) 4-Nitrophenol	14.87	139	166764	90.608	nq	98
57) 2,4-Dinitrotoluene	14.94	165	175013	47.785	nq	# 99
58) Fluorene	15.62	166	533537	48.419	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.22	232	159831	49.818	nq	97
60) Diethylphthalate	15.39	149	558095	47.969	nq	97
61) 4-Chlorophenyl-phenylether	15.61	204	293770	46.256	nq	99
62) 4-Nitroaniline	15.66	138	116314	44.388	nq	94
63) Azobenzene	15.90	77	542824	48.591	nq	98
65) 4,6-Dinitro-2-methylphenol	15.72	198	110422	52.015	nq	98
66) n-Nitrosodiphenylamine	15.83	169	466629	47.959	nq	99
67) 4-Bromophenyl-phenylether	16.51	248	204684	45.857	nq	99
68) Hexachlorobenzene	16.63	284	220140	47.563	nq	96
69) Atrazine	16.78	200	175731	46.545	nq	98
70) Pentachlorophenol	16.99	266	214030	92.452	nq	95
71) Phenanthrene	17.36	178	832172	46.972	nq	99
72) Anthracene	17.45	178	855601	48.498	nq	100
73) Carbazole	17.73	167	778730	46.358	nq	98
74) Di-n-butylphthalate	18.28	149	925968	46.555	nq	99
75) Fluoranthene	19.37	202	1025648	47.080	nq	98
77) Benzidine	19.56	184	350944	43.515	nq	99
78) Pyrene	19.73	202	1004134	48.336	nq	99
80) Butylbenzylphthalate	20.63	149	417591	48.024	nq	98
81) Benzo(a)anthracene	21.56	228	958412	47.651	nq	99
82) 3,3'-Dichlorobenzidine	21.48	252	244553	32.144	nq	99
83) Chrysene	21.62	228	924567	47.526	nq	99
84) Bis(2-ethylhexyl)phthalate	21.47	149	576888	47.697	nq	# 99
85) Di-n-octyl phthalate	22.65	149	954138	45.734	nq	100
87) Indeno(1,2,3-cd)pyrene	28.26	276	1182185	49.589	nq	99
88) Benzo(b)fluoranthene	23.72	252	1022109	49.574	nq	97
89) Benzo(k)fluoranthene	23.79	252	965954	48.951	nq	98
90) Benzo(a)pyrene	24.56	252	922034	47.873	nq	99
91) Dibenzo(a,h)anthracene	28.32	278	967700	50.620	nq	97
92) Benzo(g,h,i)perylene	29.37	276	967699	50.574	nq	98

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

