

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031220\
 Data File : BG044728.D
 Acq On : 12 Mar 2020 10:11
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 3/16/2020 1:48:05 PM

Quant Time: Mar 13 06:44:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 16:27:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	105661	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	420070	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	288483	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	653149	20.00	ng	0.00
76) Chrysene-d12	21.71	240	580162	20.00	ng	0.00
87) Perylene-d12	24.93	264	683921	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	504157	74.28	ng	0.00
7) Phenol-d6	7.21	99	729620	73.98	ng	0.00
23) Nitrobenzene-d5	9.22	82	711527	74.33	ng	0.00
42) 2,4,6-Tribromophenol	16.17	330	284237	75.25	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	1466901	72.82	ng	0.00
79) Terphenyl-d14	20.05	244	2215588	71.43	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.53	88	116605	35.675	ng	96
3) Pyridine	3.92	79	352898	36.471	ng	97
4) n-Nitrosodimethylamine	3.83	42	146079	39.789	ng	94
6) Aniline	7.38	93	451752	36.814	ng	97
8) 2-Chlorophenol	7.62	128	250956	37.038	ng	97
9) Benzaldehyde	7.19	77	218678	35.187	ng	98
10) Phenol	7.24	94	362729	37.152	ng	99
11) bis(2-Chloroethyl)ether	7.47	93	275360	35.339	ng	99
12) 1,3-Dichlorobenzene	7.95	146	304899	36.529	ng	98
13) 1,4-Dichlorobenzene	8.09	146	303227	36.747	ng	100
14) 1,2-Dichlorobenzene	8.42	146	286473	36.605	ng	99
15) Benzyl Alcohol	8.29	79	287574	36.344	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.59	45	499969	36.359	ng	99
17) 2-Methylphenol	8.49	107	240943	36.432	ng	95
18) Hexachloroethane	9.15	117	119568	36.805	ng	98
19) n-Nitroso-di-n-propylamine	8.86	70	247383	35.526	ng	99
20) 3+4-Methylphenols	8.82	107	332163	36.434	ng	98
22) Acetophenone	8.87	105	423170	35.559	ng	97
24) Nitrobenzene	9.26	77	358154	36.057	ng	99
25) Isophorone	9.79	82	653944	35.001	ng	99
26) 2-Nitrophenol	9.97	139	129912	38.860	ng	96
27) 2,4-Dimethylphenol	10.03	122	216018	35.504	ng	93
28) bis(2-Chloroethoxy)methane	10.27	93	402341	36.085	ng	98
29) 2,4-Dichlorophenol	10.50	162	263592	37.091	ng	98
30) 1,2,4-Trichlorobenzene	10.73	180	306065	36.019	ng	95
31) Naphthalene	10.91	128	841595	36.167	ng	99
32) Benzoic acid	10.16	122	171099	38.496	ng	94
33) 4-Chloroaniline	11.01	127	374553	35.727	ng	99
34) Hexachlorobutadiene	11.21	225	207954	37.214	ng	95
35) Caprolactam	11.78	113	99167	36.856	ng	89
36) 4-Chloro-3-methylphenol	12.13	107	304218	35.893	ng	98
37) 2-Methylnaphthalene	12.52	142	632490	36.336	ng	98
38) 1-Methylnaphthalene	12.73	142	596861	36.473	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	337988	35.575	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.88	237	186537	35.927	nq	96
43) 2,4,6-Trichlorophenol	13.12	196	230251	36.518	nq	95
44) 2,4,5-Trichlorophenol	13.19	196	236189	36.121	nq	98
46) 1,1'-Biphenyl	13.52	154	826134	35.836	nq	99
47) 2-Chloronaphthalene	13.56	162	625812	35.536	nq	98
48) 2-Nitroaniline	13.76	65	236823	38.411	nq	96
49) Acenaphthylene	14.41	152	988769	35.838	nq	99
50) Dimethylphthalate	14.14	163	837163	36.436	nq	98
51) 2,6-Dinitrotoluene	14.25	165	171714	37.077	nq	97
52) Acenaphthene	14.76	154	655478	36.277	nq	95
53) 3-Nitroaniline	14.58	138	199557	37.478	nq	99
54) 2,4-Dinitrophenol	14.78	184	77156	36.887	nq	# 80
55) Dibenzofuran	15.09	168	950200	36.264	nq	100
56) 4-Nitrophenol	14.88	139	152280	37.474	nq	99
57) 2,4-Dinitrotoluene	15.04	165	244624	38.513	nq	93
58) Fluorene	15.74	166	779121	36.147	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.31	232	222469m	36.955	nq	
60) Diethylphthalate	15.51	149	875025	36.514	nq	98
61) 4-Chlorophenyl-phenylether	15.73	204	419947	36.184	nq	99
62) 4-Nitroaniline	15.74	138	212507	37.428	nq	96
63) Azobenzene	16.03	77	896164	36.011	nq	98
65) 4,6-Dinitro-2-methylphenol	15.80	198	107381	37.040	nq	94
66) n-Nitrosodiphenylamine	15.94	169	695020	35.344	nq	99
67) 4-Bromophenyl-phenylether	16.62	248	291216	35.665	nq	96
68) Hexachlorobenzene	16.75	284	314135	35.462	nq	99
69) Atrazine	16.89	200	274496	38.101	nq	99
70) Pentachlorophenol	17.08	266	186144	38.191	nq	96
71) Phenanthrene	17.48	178	1293665	37.064	nq	99
72) Anthracene	17.57	178	1249079	36.293	nq	99
73) Carbazole	17.83	167	1207751	36.533	nq	99
74) Di-n-butylphthalate	18.40	149	1490609	37.091	nq	99
75) Fluoranthene	19.49	202	1533942	36.911	nq	99
77) Benzidine	19.66	184	811560	43.089	nq	99
78) Pyrene	19.84	202	1543919	36.272	nq	99
80) Butylbenzylphthalate	20.74	149	695698	36.739	nq	95
81) Benzo(a)anthracene	21.69	228	1502448	35.964	nq	100
82) 3,3'-Dichlorobenzidine	21.60	252	599323	37.083	nq	96
83) Chrysene	21.75	228	1459185	36.566	nq	99
84) Bis(2-ethylhexyl)phthalate	21.62	149	945770	36.189	nq	99
85) Di-n-octyl phthalate	22.85	149	1613511	36.895	nq	99
86) Indeno(1,2,3-cd)pyrene	28.59	276	1902961	36.374	nq	# 95
88) Benzo(b)fluoranthene	23.91	252	1657752	36.716	nq	99
89) Benzo(k)fluoranthene	23.97	252	1509022	35.317	nq	99
90) Benzo(a)pyrene	24.77	252	1544376	36.555	nq	99
91) Dibenzo(a,h)anthracene	28.66	278	1501976	36.036	nq	99
92) Benzo(g,h,i)perylene	29.73	276	1520168	36.100	nq	98

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

