

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG073020\
 Data File : BG046115.D
 Acq On : 30 Jul 2020 17:55
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Sohil
 7/31/2020 12:04:18 PM

Quant Time: Jul 30 18:48:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 16:37:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	115254	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	404505	20.00	ng	0.00
39) Acenaphthene-d10	14.59	164	258587	20.00	ng	0.00
64) Phenanthrene-d10	17.33	188	594421	20.00	ng	0.00
76) Chrysene-d12	21.58	240	585564	20.00	ng	0.00
86) Perylene-d12	24.68	264	640013	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.54	112	1003602	141.21	ng	0.00
7) Phenol-d6	7.13	99	1341099	131.07	ng	0.01
23) Nitrobenzene-d5	9.12	82	1192157	169.63	ng	0.01
42) 2,4,6-Tribromophenol	16.07	330	662650	276.11	ng	0.00
45) 2-Fluorobiphenyl	13.21	172	2459909	148.38	ng	0.00
79) Terphenyl-d14	0.00	244	0d	0.00	ng	

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.45	88	231670	70.654	ng	99
3) Pyridine	3.85	79	647127m	65.270	ng	
4) n-Nitrosodimethylamine	3.76	42	282032	84.125	ng	98
6) Aniline	7.28	93	815048	62.267	ng	97
8) 2-Chlorophenol	7.53	128	541924	70.833	ng	96
9) Benzaldehyde	7.10	77	359056	55.385	ng	95
10) Phenol	7.15	94	674581	65.722	ng	97
11) bis(2-Chloroethyl)ether	7.38	93	544214	64.119	ng	99
12) 1,3-Dichlorobenzene	7.85	146	660517	74.901	ng	97
13) 1,4-Dichlorobenzene	8.00	146	655292	75.538	ng	100
14) 1,2-Dichlorobenzene	8.31	146	614414	73.590	ng	97
15) Benzyl Alcohol	8.19	79	509634	63.522	ng	100
16) 2,2'-oxybis(1-Chloropropan	8.49	45	882543	79.695	ng	98
17) 2-Methylphenol	8.40	107	472727	61.380	ng	99
18) Hexachloroethane	9.04	117	230333	69.006	ng	95
19) n-Nitroso-di-n-propylamine	8.77	70	447811	62.298	ng	96
20) 3+4-Methylphenols	8.74	107	656965	62.589	ng	98
22) Acetophenone	8.78	105	809855	74.943	ng	# 97
24) Nitrobenzene	9.16	77	641462	82.697	ng	98
25) Isophorone	9.69	82	1211801	70.633	ng	98
26) 2-Nitrophenol	9.86	139	328543	129.947	ng	97
27) 2,4-Dimethylphenol	9.93	122	478868	78.416	ng	99
28) bis(2-Chloroethoxy)methane	10.16	93	634584	65.869	ng	97
29) 2,4-Dichlorophenol	10.40	162	578996	91.589	ng	98
30) 1,2,4-Trichlorobenzene	10.62	180	640224	91.402	ng	98
31) Naphthalene	10.81	128	1620422	74.726	ng	97
32) Benzoic acid	10.12	122	391203	118.305	ng	98
33) 4-Chloroaniline	10.91	127	723337	74.647	ng	97
34) Hexachlorobutadiene	11.10	225	428748	104.834	ng	100
35) Caprolactam	11.71	113	200912	83.708	ng	# 86
36) 4-Chloro-3-methylphenol	12.04	107	572351	79.492	ng	98
37) 2-Methylnaphthalene	12.41	142	1260423	81.767	ng	96
38) 1-Methylnaphthalene	12.63	142	1194421	82.104	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.78	216	725784	99.567	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.77	237	444074	102.670	ng	96
43) 2,4,6-Trichlorophenol	13.02	196	503305	102.094	ng	96
44) 2,4,5-Trichlorophenol	13.10	196	545429	97.853	ng	99
46) 1,1'-Biphenyl	13.42	154	1522324	78.259	ng	96
47) 2-Chloronaphthalene	13.46	162	1229878	80.132	ng	94
48) 2-Nitroaniline	13.66	65	378244	105.301	ng	98
49) Acenaphthylene	14.31	152	1878377	75.854	ng	95
50) Dimethylphthalate	14.05	163	1506738	77.642	ng	98
51) 2,6-Dinitrotoluene	14.16	165	386054	128.206	ng	96
52) Acenaphthene	14.65	154	1314525	84.159	ng	98
53) 3-Nitroaniline	14.49	138	397484	106.520	ng	97
54) 2,4-Dinitrophenol	14.69	184	239146	225.143	ng	92
55) Dibenzofuran	14.99	168	1834128	80.107	ng	94
56) 4-Nitrophenol	14.80	139	346343	109.544	ng	99
57) 2,4-Dinitrotoluene	14.95	165	541012	145.407	ng	97
58) Fluorene	15.63	166	1493319	79.505	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.21	232	506428	117.677	ng	99
60) Diethylphthalate	15.41	149	1500683	73.737	ng	96
61) 4-Chlorophenyl-phenylether	15.63	204	804843	86.131	ng	98
62) 4-Nitroaniline	15.65	138	398213	99.282	ng	97
63) Azobenzene	15.92	77	1404589	64.302	ng	96
65) 4,6-Dinitro-2-methylphenol	15.71	198	353340	202.954	ng	98
66) n-Nitrosodiphenylamine	15.84	169	1366810	74.202	ng	94
67) 4-Bromophenyl-phenylether	16.52	248	585971	88.264	ng	96
68) Hexachlorobenzene	16.64	284	642147	96.390	ng	98
69) Atrazine	16.79	200	549389	100.878	ng	100
70) Pentachlorophenol	16.98	266	459490	124.054	ng	98
71) Phenanthrene	17.37	178	2298912	73.155	ng	93
72) Anthracene	17.46	178	2301429	72.164	ng	# 91
73) Carbazole	17.72	167	2260813	70.724	ng	93
74) Di-n-butylphthalate	18.29	149	2317003	59.981	ng	# 94
75) Fluoranthene	19.38	202	2720058	74.874	ng	89
77) Benzidine	19.55	184	1329370	79.400	ng	97
78) Pyrene	19.74	202	2723896	68.147	ng	# 90
80) Butylbenzylphthalate	20.63	149	1183124	62.637	ng	96
81) Benzo(a)anthracene	21.56	228	2811292	73.015	ng	90
82) 3,3'-Dichlorobenzidine	21.47	252	1300730	91.424	ng	94
83) Chrysene	21.63	228	2717898	74.096	ng	89
84) Bis(2-ethylhexyl)phthalate	21.49	149	1557440	56.803	ng	95
85) Di-n-octyl phthalate	22.67	149	2757539	57.955	ng	96
87) Indeno(1,2,3-cd)pyrene	28.22	276	3951865	84.403	ng	99
88) Benzo(b)fluoranthene	23.71	252	3242265	79.803	ng	95
89) Benzo(k)fluoranthene	23.78	252	3093289	80.519	ng	96
90) Benzo(a)pyrene	24.55	252	3086666	80.793	ng	95
91) Dibenzo(a,h)anthracene	28.29	278	3203987	84.842	ng	99
92) Benzo(g,h,i)perylene	29.33	276	3276486	86.215	ng	97

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

