

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081320\
 Data File : BF121277.D
 Acq On : 13 Aug 2020 18:35
 Operator : JU/CG
 Sample : L3653-02MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 2

Instrument :
 BNA_F
 ClientSampleId :
 BW-STOCKMS

Manual Integrations
 APPROVED

Sohil
 8/14/2020 2:53:25 PM

Quant Time: Aug 13 19:02:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 13 15:23:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	105470	40.00	ng	0.00
21) Naphthalene-d8	8.00	136	414688	40.00	ng	0.00
39) Acenaphthene-d10	9.76	164	215159	40.00	ng	0.00
64) Phenanthrene-d10	11.24	188	413652	40.00	ng	0.00
76) Chrysene-d12	13.87	240	332829	40.00	ng	0.00
86) Perylene-d12	15.29	264	334801	40.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	622260	178.89	ng	0.01
7) Phenol-d6	6.36	99	755323	168.51	ng	0.00
23) Nitrobenzene-d5	7.29	82	498268	131.34	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	232146	175.27	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	925232	118.97	ng	0.00
79) Terphenyl-d14	12.82	244	998203	116.93	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.34	88	107027	72.302	ng	99
3) Pyridine	3.05	79	263456	68.371	ng	99
4) n-Nitrosodimethylamine	3.00	42	140233	88.859	ng	96
6) Aniline	6.38	93	373279	70.249	ng	# 79
8) 2-Chlorophenol	6.50	128	319154	83.441	ng	99
9) Benzaldehyde	6.27	77	197176	85.359	ng	97
10) Phenol	6.37	94	395612	81.360	ng	99
11) bis(2-Chloroethyl)ether	6.46	93	300278	85.553	ng	97
12) 1,3-Dichlorobenzene	6.66	146	329050	81.418	ng	99
13) 1,4-Dichlorobenzene	6.74	146	330389	80.881	ng	100
14) 1,2-Dichlorobenzene	6.89	146	316845	82.185	ng	99
15) Benzyl Alcohol	6.87	79	249387	83.056	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.00	45	423375	83.769	ng	99
17) 2-Methylphenol	6.99	107	256798	84.996	ng	98
18) Hexachloroethane	7.23	117	120927	83.713	ng	96
19) n-Nitroso-di-n-propylamine	7.14	70	192984	77.488	ng	98
20) 3+4-Methylphenols	7.14	107	294200	76.749	ng	92
22) Acetophenone	7.13	105	405337	75.900	ng	99
24) Nitrobenzene	7.31	77	321210	78.028	ng	100
25) Isophorone	7.55	82	586416	76.987	ng	98
26) 2-Nitrophenol	7.63	139	169294	84.877	ng	98
27) 2,4-Dimethylphenol	7.67	122	271693	85.016	ng	98
28) bis(2-Chloroethoxy)methane	7.76	93	367964	86.368	ng	99
29) 2,4-Dichlorophenol	7.87	162	259219	79.320	ng	99
30) 1,2,4-Trichlorobenzene	7.95	180	282250	78.252	ng	100
31) Naphthalene	8.03	128	887023	76.040	ng	100
32) Benzoic acid	7.80	122	161790	91.955	ng	99
33) 4-Chloroaniline	8.08	127	268284	56.356	ng	98
34) Hexachlorobutadiene	8.15	225	163281	74.819	ng	98
35) Caprolactam	8.45	113	85288m	86.645	ng	
36) 4-Chloro-3-methylphenol	8.57	107	268121	80.415	ng	96
37) 2-Methylnaphthalene	8.72	142	590495	76.925	ng	99
38) 1-Methylnaphthalene	8.82	142	557464	76.700	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.89	216	275411	76.405	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.87	237	208191	143.126	nq	97
43) 2,4,6-Trichlorophenol	9.00	196	184468	80.466	nq	99
44) 2,4,5-Trichlorophenol	9.04	196	199749	80.220	nq	98
46) 1,1'-Biphenyl	9.18	154	719181	77.369	nq	100
47) 2-Chloronaphthalene	9.20	162	552760	77.250	nq	98
48) 2-Nitroaniline	9.30	65	169968	84.507	nq	99
49) Acenaphthylene	9.62	152	886774	78.026	nq	100
50) Dimethylphthalate	9.49	163	740193	91.436	nq	100
51) 2,6-Dinitrotoluene	9.55	165	147466	81.720	nq	99
52) Acenaphthene	9.79	154	506632	75.293	nq	100
53) 3-Nitroaniline	9.72	138	117007	56.788	nq	97
54) 2,4-Dinitrophenol	9.83	184	134395	140.905	nq	91
55) Dibenzofuran	9.96	168	789474	73.676	nq	100
56) 4-Nitrophenol	9.89	139	235325	176.981	nq	100
57) 2,4-Dinitrotoluene	9.95	165	191470	80.851	nq	94
58) Fluorene	10.30	166	587400	72.417	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.09	232	166719	84.682	nq	97
60) Diethylphthalate	10.18	149	649943	78.953	nq	99
61) 4-Chlorophenyl-phenylether	10.30	204	288048	73.015	nq	100
62) 4-Nitroaniline	10.33	138	161775	78.312	nq	98
63) Azobenzene	10.46	77	620003	75.378	nq	98
65) 4,6-Dinitro-2-methylphenol	10.36	198	101911	81.568	nq	83
66) n-Nitrosodiphenylamine	10.42	169	556298	78.442	nq	100
67) 4-Bromophenyl-phenylether	10.79	248	194677	80.693	nq	100
68) Hexachlorobenzene	10.85	284	217755	79.152	nq	99
69) Atrazine	10.94	200	153631	72.425	nq	99
70) Pentachlorophenol	11.05	266	222261	179.753	nq	99
71) Phenanthrene	11.26	178	902783	76.298	nq	100
72) Anthracene	11.32	178	934963	78.742	nq	99
73) Carbazole	11.47	167	870222	74.939	nq	99
74) Di-n-butylphthalate	11.80	149	1055152	80.500	nq	100
75) Fluoranthene	12.45	202	1025459	77.832	nq	99
77) Benzidine	12.57	184	428604	97.970	nq	99
78) Pyrene	12.68	202	1054115	85.061	nq	99
80) Butylbenzylphthalate	13.30	149	468229	92.498	nq	98
81) Benzo(a)anthracene	13.86	228	857936	76.010	nq	100
82) 3,3'-Dichlorobenzidine	13.83	252	291856	65.173	nq	98
83) Chrysene	13.90	228	907236	78.726	nq	99
84) Bis(2-ethylhexyl)phthalate	13.86	149	563517	80.766	nq	99
85) Di-n-octyl phthalate	14.47	149	1117442	87.352	nq	99
87) Indeno(1,2,3-cd)pyrene	16.67	276	924911	74.313	nq	100
88) Benzo(b)fluoranthene	14.89	252	946164	84.355	nq	100
89) Benzo(k)fluoranthene	14.92	252	831787	79.858	nq	98
90) Benzo(a)pyrene	15.23	252	805039	79.520	nq	100
91) Dibenzo(a,h)anthracene	16.69	278	764379	71.675	nq	99
92) Benzo(g,h,i)perylene	17.09	276	773234	73.472	nq	99

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

