

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
 Data File : BF119838.D
 Acq On : 8 Apr 2020 21:30
 Operator : MA/SJ
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 4/10/2020 9:57:06 AM

Quant Time: Apr 09 01:17:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 08 21:57:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	217262	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	806776	20.00	ng	0.00
39) Acenaphthene-d10	9.82	164	405653	20.00	ng	0.00
64) Phenanthrene-d10	11.30	188	712728	20.00	ng	0.00
76) Chrysene-d12	13.94	240	426304	20.00	ng	0.00
87) Perylene-d12	15.37	264	377334	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.38	112	1704072	124.86	ng	0.00
7) Phenol-d6	6.42	99	2302440	132.06	ng	0.02
23) Nitrobenzene-d5	7.35	82	2158138	145.38	ng	0.01
42) 2,4,6-Tribromophenol	10.62	330	650682	155.48	ng	0.01
45) 2-Fluorobiphenyl	9.15	172	4002690	146.18	ng	0.00
79) Terphenyl-d14	12.90	244	3636180	167.11	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.45	88	387479	57.484	ng	99
3) Pyridine	3.17	79	1010526	54.417	ng	97
4) n-Nitrosodimethylamine	3.15	42	532126	58.761	ng	94
6) Aniline	6.44	93	1540632	64.028	ng	# 56
8) 2-Chlorophenol	6.56	128	989403	69.024	ng	95
9) Benzaldehyde	6.32	77	438675	39.664	ng	97
10) Phenol	6.43	94	1313999	70.635	ng	91
11) bis(2-Chloroethyl)ether	6.52	93	1033120	69.438	ng	98
12) 1,3-Dichlorobenzene	6.72	146	999926	64.453	ng	99
13) 1,4-Dichlorobenzene	6.79	146	1119352	72.133	ng	98
14) 1,2-Dichlorobenzene	6.95	146	1020344	70.347	ng	98
15) Benzyl Alcohol	6.93	79	884534	67.448	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.06	45	1942123	64.715	ng	96
17) 2-Methylphenol	7.03	107	927396	70.613	ng	97
18) Hexachloroethane	7.29	117	404802	71.183	ng	96
19) n-Nitroso-di-n-propylamine	7.21	70	723108	68.812	ng	97
20) 3+4-Methylphenols	7.20	107	998256	62.021	ng	94
22) Acetophenone	7.20	105	1262775	65.505	ng	# 94
24) Nitrobenzene	7.37	77	1068563	67.356	ng	99
25) Isophorone	7.61	82	2188547	72.678	ng	99
26) 2-Nitrophenol	7.68	139	546161	87.436	ng	91
27) 2,4-Dimethylphenol	7.72	122	789979	61.163	ng	99
28) bis(2-Chloroethoxy)methane	7.82	93	1196325	67.184	ng	99
29) 2,4-Dichlorophenol	7.92	162	933436	78.654	ng	97
30) 1,2,4-Trichlorobenzene	8.00	180	1012124	77.117	ng	98
31) Naphthalene	8.09	128	3055809	73.603	ng	98
32) Benzoic acid	7.89	122	811170	97.634	ng	98
33) 4-Chloroaniline	8.13	127	1402929	73.745	ng	99
34) Hexachlorobutadiene	8.20	225	603771	82.221	ng	99
35) Caprolactam	8.55	113	309360m	79.650	ng	
36) 4-Chloro-3-methylphenol	8.63	107	987990	76.170	ng	96
37) 2-Methylnaphthalene	8.77	142	2094201	76.858	ng	99
38) 1-Methylnaphthalene	8.87	142	1954328	75.777	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.95	216	941086	79.149	ng	99

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41) Hexachlorocyclopentadiene	8.93	237	589505	87.210	nq	99
43) 2,4,6-Trichlorophenol	9.06	196	684299	80.423	nq	99
44) 2,4,5-Trichlorophenol	9.10	196	776134	81.426	nq	100
46) 1,1'-Biphenyl	9.25	154	2523805	76.390	nq	96
47) 2-Chloronaphthalene	9.27	162	1950300	75.361	nq	96
48) 2-Nitroaniline	9.36	65	743025	83.181	nq	97
49) Acenaphthylene	9.69	152	2932189	72.150	nq	99
50) Dimethylphthalate	9.56	163	2389483	82.929	nq	100
51) 2,6-Dinitrotoluene	9.62	165	526347	85.224	nq #	84
52) Acenaphthene	9.86	154	1955342m	77.178	nq	
53) 3-Nitroaniline	9.79	138	588642	75.494	nq	97
54) 2,4-Dinitrophenol	9.89	184	332556	146.881	nq	94
55) Dibenzofuran	10.03	168	2383182	65.608	nq	97
56) 4-Nitrophenol	9.95	139	430360	69.967	nq	97
57) 2,4-Dinitrotoluene	10.02	165	620808	79.794	nq #	87
58) Fluorene	10.37	166	1840747	68.527	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.15	232	515143	72.302	nq	99
60) Diethylphthalate	10.25	149	2206109	75.437	nq	98
61) 4-Chlorophenyl-phenylether	10.36	204	915459	69.692	nq	96
62) 4-Nitroaniline	10.41	138	505173	67.564	nq	99
63) Azobenzene	10.52	77	2006620	68.138	nq	99
65) 4,6-Dinitro-2-methylphenol	10.43	198	356395	119.249	nq	77
66) n-Nitrosodiphenylamine	10.49	169	1686657	72.086	nq	98
67) 4-Bromophenyl-phenylether	10.85	248	653266	80.481	nq	96
68) Hexachlorobenzene	10.92	284	649909	78.230	nq	98
69) Atrazine	11.02	200	444786	69.341	nq	96
70) Pentachlorophenol	11.11	266	380866	78.187	nq	99
71) Phenanthrene	11.33	178	2823545	76.401	nq	97
72) Anthracene	11.39	178	2827979	75.291	nq	99
73) Carbazole	11.54	167	2590637	69.682	nq	98
74) Di-n-butylphthalate	11.87	149	3273549	82.311	nq	98
75) Fluoranthene	12.52	202	2897709	72.450	nq	97
77) Benzidine	12.64	184	1334797m	73.986	nq	
78) Pyrene	12.75	202	2711389	77.498	nq	98
80) Butylbenzylphthalate	13.37	149	1347727	95.243	nq	100
81) Benzo(a)anthracene	13.94	228	2061600	74.367	nq	98
82) 3,3'-Dichlorobenzidine	13.90	252	745160	68.173	nq	97
83) Chrysene	13.97	228	2146341	75.020	nq	98
84) Bis(2-ethylhexyl)phthalate	13.93	149	1799889	106.701	nq	98
85) Di-n-octyl phthalate	14.54	149	2986756	100.677	nq	100
86) Indeno(1,2,3-cd)pyrene	16.80	276	1901921	70.585	nq	99
88) Benzo(b)fluoranthene	14.97	252	2204084	87.443	nq	99
89) Benzo(k)fluoranthene	15.00	252	1728357	72.011	nq	99
90) Benzo(a)pyrene	15.33	252	1773616	77.428	nq	98
91) Dibenzo(a,h)anthracene	16.83	278	1547742	77.249	nq	99
92) Benzo(g,h,i)perylene	17.24	276	1563552	77.679	nq	98

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

