

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF050119\
 Data File : BF114059.D
 Acq On : 1 May 2019 10:50
 Operator : JU/SJ
 Sample : K2194-01MS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-042919-AMS

Manual Integrations
 APPROVED

Sohil

Quant Time: May 01 14:36:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

5/2/2019 4:53:28 PM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	89183	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	369892	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	188587	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	416379	20.00	ng	0.00
76) Chrysene-d12	14.06	240	298877	20.00	ng	-0.01
87) Perylene-d12	15.56	264	283630	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	255736	49.05	ng	0.00
7) Phenol-d6	6.52	99	352528	53.89	ng	0.00
23) Nitrobenzene-d5	7.45	82	197512	32.97	ng	-0.01
42) 2,4,6-Tribromophenol	10.72	330	123542	53.78	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	460612	38.20	ng	-0.01
79) Terphenyl-d14	13.00	244	559717	38.39	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.61	88	35263	14.346	ng	95
3) Pyridine	3.40	79	86751	12.930	ng	97
4) n-Nitrosodimethylamine	3.31	42	31811m	13.851	ng	
6) Aniline	6.55	93	95394	10.619	ng	96
8) 2-Chlorophenol	6.67	128	98137	16.485	ng	99
9) Benzaldehyde	6.44	77	20520	1.236	ng	93
10) Phenol	6.53	94	129727	16.583	ng	95
11) bis(2-Chloroethyl)ether	6.62	93	94163	15.996	ng	98
12) 1,3-Dichlorobenzene	6.83	146	110941	16.454	ng	98
13) 1,4-Dichlorobenzene	6.91	146	117547	17.333	ng	100
14) 1,2-Dichlorobenzene	7.06	146	108864	17.467	ng	97
15) Benzyl Alcohol	7.03	79	81155	18.419	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.16	45	111485	16.224	ng	92
17) 2-Methylphenol	7.14	107	83033	15.912	ng	99
18) Hexachloroethane	7.40	117	37989	15.980	ng	93
19) n-Nitroso-di-n-propylamine	7.29	70	70931	16.743	ng	96
20) 3+4-Methylphenols	7.29	107	102909	17.455	ng	# 54
22) Acetophenone	7.29	105	141257	17.161	ng	# 95
24) Nitrobenzene	7.46	77	104626	15.806	ng	97
25) Isophorone	7.70	82	207880	16.959	ng	98
26) 2-Nitrophenol	7.79	139	54488	18.043	ng	99
27) 2,4-Dimethylphenol	7.82	122	100452	20.416	ng	99
28) bis(2-Chloroethoxy)methane	7.92	93	134036	17.160	ng	100
29) 2,4-Dichlorophenol	8.03	162	99059	17.606	ng	95
30) 1,2,4-Trichlorobenzene	8.12	180	109449	17.446	ng	99
31) Naphthalene	8.19	128	320841	18.259	ng	99
32) Benzoic acid	7.82	122	100452	30.232	ng	# 14
33) 4-Chloroaniline	8.24	127	94434	12.701	ng	96
34) Hexachlorobutadiene	8.32	225	60014	15.346	ng	98
35) Caprolactam	8.58	113	29654	16.568	ng	93
36) 4-Chloro-3-methylphenol	8.72	107	91314	16.382	ng	98
37) 2-Methylnaphthalene	8.89	142	210174	17.737	ng	99
38) 1-Methylnaphthalene	8.99	142	204291	17.863	ng	97
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	104941	17.131	ng	99

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41) Hexachlorocyclopentadiene	9.05	237	37606	11.009	nq	99
43) 2,4,6-Trichlorophenol	9.16	196	74074	19.043	nq	99
44) 2,4,5-Trichlorophenol	9.21	196	80259	18.395	nq	98
46) 1,1'-Biphenyl	9.35	154	276433	19.188	nq	97
47) 2-Chloronaphthalene	9.37	162	216679	18.495	nq	99
48) 2-Nitroaniline	9.46	65	58650	19.096	nq	94
49) Acenaphthylene	9.79	152	343668	18.737	nq	97
50) Dimethylphthalate	9.64	163	291804	21.395	nq	99
51) 2,6-Dinitrotoluene	9.70	165	55485	18.962	nq	98
52) Acenaphthene	9.96	154	195334	18.021	nq	98
53) 3-Nitroaniline	9.87	138	44902	14.603	nq	96
54) 2,4-Dinitrophenol	9.99	184	5246	12.032	nq	# 5
55) Dibenzofuran	10.13	168	296190	18.242	nq	100
56) 4-Nitrophenol	10.03	139	87996	36.145	nq	100
57) 2,4-Dinitrotoluene	10.11	165	68940	18.666	nq	# 96
58) Fluorene	10.47	166	249369	20.400	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.25	232	70991	18.515	nq	# 97
60) Diethylphthalate	10.34	149	249945	19.298	nq	99
61) 4-Chlorophenyl-phenylether	10.47	204	128541	19.855	nq	95
62) 4-Nitroaniline	10.48	138	52126	17.372	nq	90
63) Azobenzene	10.63	77	228068	18.169	nq	94
65) 4,6-Dinitro-2-methylphenol	10.52	198	15713	14.288	nq	86
66) n-Nitrosodiphenylamine	10.58	169	221636	17.740	nq	97
67) 4-Bromophenyl-phenylether	10.96	248	83039	16.509	nq	97
68) Hexachlorobenzene	11.03	284	84639	15.312	nq	94
69) Atrazine	11.10	200	75611	18.604	nq	99
70) Pentachlorophenol	11.22	266	79101	25.275	nq	97
71) Phenanthrene	11.44	178	451745	21.588	nq	98
72) Anthracene	11.49	178	412371	19.515	nq	98
73) Carbazole	11.64	167	329265	17.466	nq	99
74) Di-n-butylphthalate	11.97	149	405655	18.292	nq	98
75) Fluoranthene	12.63	202	544086	23.832	nq	96
77) Benzidine	12.76	184	52755	5.924	nq	# 82
78) Pyrene	12.86	202	565503	27.056	nq	98
80) Butylbenzylphthalate	13.47	149	152269	18.517	nq	99
81) Benzo(a)anthracene	14.05	228	394875	22.424	nq	99
82) 3,3'-Dichlorobenzidine	14.00	252	65444	10.140	nq	99
83) Chrysene	14.09	228	359976	20.990	nq	98
84) Bis(2-ethylhexyl)phthalate	14.03	149	222916	21.306	nq	98
85) Di-n-octyl phthalate	14.66	149	332515	20.304	nq	99
86) Indeno(1,2,3-cd)pyrene	17.06	276	349624	21.522	nq	98
88) Benzo(b)fluoranthene	15.12	252	385992m	21.510	nq	
89) Benzo(k)fluoranthene	15.15	252	296607	19.215	nq	97
90) Benzo(a)pyrene	15.50	252	334239	21.534	nq	99
91) Dibenzo(a,h)anthracene	17.08	278	274268	18.824	nq	99
92) Benzo(g,h,i)perylene	17.52	276	311736	21.201	nq	97

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

