

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082819\
 Data File : BF116609.D
 Acq On : 28 Aug 2019 11:16
 Operator : HP/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Jagrut
 8/29/2019 7:46:44 PM

Quant Time: Aug 28 11:49:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 28 11:25:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	149828	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	623692	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	317433	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	529173	20.00	ng	0.00
76) Chrysene-d12	14.02	240	262966	20.00	ng	0.00
87) Perylene-d12	15.47	264	256094	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	1069874	104.34	ng	0.00
7) Phenol-d6	6.50	99	1370964	109.04	ng	0.01
23) Nitrobenzene-d5	7.44	82	1319011	115.70	ng	0.01
42) 2,4,6-Tribromophenol	10.69	330	341548	125.31	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	2140622	108.60	ng	0.00
79) Terphenyl-d14	12.97	244	1881154	133.62	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.63	88	243412	54.826	ng	99
3) Pyridine	3.39	79	706837	52.280	ng	100
4) n-Nitrosodimethylamine	3.35	42	273082	55.329	ng	99
6) Aniline	6.54	93	948242	55.707	ng	99
8) 2-Chlorophenol	6.66	128	588047	54.648	ng	97
9) Benzaldehyde	6.42	77	370893	44.150	ng	99
10) Phenol	6.52	94	770701	56.448	ng	91
11) bis(2-Chloroethyl)ether	6.61	93	579679	55.762	ng	97
12) 1,3-Dichlorobenzene	6.82	146	665000	57.096	ng	99
13) 1,4-Dichlorobenzene	6.89	146	662893	59.136	ng	98
14) 1,2-Dichlorobenzene	7.05	146	617952	58.437	ng	98
15) Benzyl Alcohol	7.01	79	549135	53.969	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.15	45	763625	53.127	ng	100
17) 2-Methylphenol	7.12	107	520539	58.385	ng	99
18) Hexachloroethane	7.39	117	254280	58.164	ng	94
19) n-Nitroso-di-n-propylamine	7.29	70	438042	51.058	ng	100
20) 3+4-Methylphenols	7.27	107	607136	55.590	ng	94
22) Acetophenone	7.28	105	852167	54.641	ng	# 92
24) Nitrobenzene	7.45	77	672675	56.978	ng	96
25) Isophorone	7.70	82	1187368	52.853	ng	99
26) 2-Nitrophenol	7.77	139	324798	75.318	ng	99
27) 2,4-Dimethylphenol	7.80	122	480483	53.400	ng	97
28) bis(2-Chloroethoxy)methane	7.90	93	739862	53.328	ng	99
29) 2,4-Dichlorophenol	8.01	162	500125	58.447	ng	98
30) 1,2,4-Trichlorobenzene	8.09	180	574965	59.686	ng	97
31) Naphthalene	8.17	128	1689824	56.007	ng	99
32) Benzoic acid	7.95	122	452805	78.718	ng	99
33) 4-Chloroaniline	8.22	127	745577	56.290	ng	99
34) Hexachlorobutadiene	8.30	225	319597	60.833	ng	98
35) Caprolactam	8.60	113	159389m	53.475	ng	
36) 4-Chloro-3-methylphenol	8.70	107	530684	54.965	ng	97
37) 2-Methylnaphthalene	8.86	142	1120708	54.099	ng	97
38) 1-Methylnaphthalene	8.96	142	1047389m	54.124	ng	
40) 1,2,4,5-Tetrachlorobenzene	9.03	216	506081	59.257	ng	99

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41) Hexachlorocyclopentadiene	9.02	237	304700	66.312	nq	98
43) 2,4,6-Trichlorophenol	9.14	196	365558	61.159	nq	97
44) 2,4,5-Trichlorophenol	9.18	196	362332	59.576	nq	97
46) 1,1'-Biphenyl	9.33	154	1399936	55.161	nq	98
47) 2-Chloronaphthalene	9.35	162	1106012	56.213	nq	96
48) 2-Nitroaniline	9.45	65	365051	60.596	nq	95
49) Acenaphthylene	9.77	152	1582384	54.970	nq	100
50) Dimethylphthalate	9.63	163	1239442	55.963	nq	99
51) 2,6-Dinitrotoluene	9.69	165	279264	61.604	nq	88
52) Acenaphthene	9.94	154	1058256	57.292	nq	95
53) 3-Nitroaniline	9.86	138	327568	60.620	nq	97
54) 2,4-Dinitrophenol	9.96	184	147214	103.285	nq	# 32
55) Dibenzofuran	10.11	168	1433114	55.659	nq	99
56) 4-Nitrophenol	10.01	139	255464	60.825	nq	98
57) 2,4-Dinitrotoluene	10.09	165	367674	63.740	nq	91
58) Fluorene	10.45	166	1068582	53.416	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.23	232	299402	59.955	nq	99
60) Diethylphthalate	10.33	149	1187052	54.532	nq	100
61) 4-Chlorophenyl-phenylether	10.45	204	537386	55.064	nq	99
62) 4-Nitroaniline	10.47	138	327993	62.062	nq	96
63) Azobenzene	10.60	77	1188797	53.115	nq	97
65) 4,6-Dinitro-2-methylphenol	10.50	198	185375	95.887	nq	97
66) n-Nitrosodiphenylamine	10.56	169	968021	56.495	nq	100
67) 4-Bromophenyl-phenylether	10.93	248	333354	60.088	nq	97
68) Hexachlorobenzene	11.00	284	362886	59.140	nq	98
69) Atrazine	11.09	200	282255	54.980	nq	97
70) Pentachlorophenol	11.19	266	230677	64.435	nq	98
71) Phenanthrene	11.41	178	1592249	55.234	nq	99
72) Anthracene	11.46	178	1592240	54.967	nq	99
73) Carbazole	11.61	167	1397659	53.753	nq	99
74) Di-n-butylphthalate	11.95	149	1677056	52.050	nq	100
75) Fluoranthene	12.59	202	1462800	49.082	nq	97
77) Benzidine	12.70	184	476715	45.058	nq	98
78) Pyrene	12.82	202	1446412	68.780	nq	99
80) Butylbenzylphthalate	13.43	149	553797	60.977	nq	95
81) Benzo(a)anthracene	14.00	228	993971	58.361	nq	99
82) 3,3'-Dichlorobenzidine	13.96	252	300596	49.527	nq	# 97
83) Chrysene	14.04	228	980764	56.815	nq	100
84) Bis(2-ethylhexyl)phthalate	14.00	149	648299	53.532	nq	100
85) Di-n-octyl phthalate	14.62	149	1035694	50.374	nq	100
86) Indeno(1,2,3-cd)pyrene	16.94	276	1037847	70.185	nq	100
88) Benzo(b)fluoranthene	15.06	252	913920	57.769	nq	98
89) Benzo(k)fluoranthene	15.09	252	827618	57.745	nq	99
90) Benzo(a)pyrene	15.42	252	815303	58.311	nq	98
91) Dibenzo(a,h)anthracene	16.96	278	797057	69.374	nq	99
92) Benzo(g,h,i)perylene	17.39	276	896911	81.886	nq	98

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

