

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082720\
 Data File : BF121443.D
 Acq On : 27 Aug 2020 18:36
 Operator : JU/CG
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Quant Time: Aug 28 03:10:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 28 03:08:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	288158	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1081583	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	563740	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	1060884	20.00	ng	0.00
76) Chrysene-d12	14.02	240	924172	20.00	ng	0.00
86) Perylene-d12	15.49	264	911598	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	392510	22.32	ng	-0.01
7) Phenol-d6	6.47	99	546180	22.45	ng	-0.02
23) Nitrobenzene-d5	7.42	82	229088	13.35	ng	-0.01
42) 2,4,6-Tribromophenol	10.68	330	105004	16.63	ng	-0.01
45) 2-Fluorobiphenyl	9.22	172	901610	25.19	ng	-0.01
79) Terphenyl-d14	12.97	244	1051984	24.00	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.62	88	90715	11.147	ng	98
3) Pyridine	3.37	79	241317	10.694	ng	99
4) n-Nitrosodimethylamine	3.30	42	103922	10.581	ng	# 99
6) Aniline	6.52	93	311826	11.299	ng	100
8) 2-Chlorophenol	6.64	128	191452	10.452	ng	98
9) Benzaldehyde	6.40	77	167822	13.617	ng	99
10) Phenol	6.49	94	252745	10.646	ng	98
11) bis(2-Chloroethyl)ether	6.59	93	212875	10.392	ng	98
12) 1,3-Dichlorobenzene	6.80	146	242325	11.526	ng	97
13) 1,4-Dichlorobenzene	6.87	146	242425	11.512	ng	98
14) 1,2-Dichlorobenzene	7.03	146	233333	11.913	ng	98
15) Benzyl Alcohol	6.99	79	174468	10.693	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.13	45	320863	11.115	ng	98
17) 2-Methylphenol	7.10	107	165911	10.785	ng	99
18) Hexachloroethane	7.37	117	77826	10.188	ng	97
19) n-Nitroso-di-n-propylamine	7.26	70	149585	10.773	ng	97
20) 3+4-Methylphenols	7.25	107	215493	11.491	ng	# 71
22) Acetophenone	7.26	105	308757	11.554	ng	# 97
24) Nitrobenzene	7.43	77	132632	7.236	ng	94
25) Isophorone	7.67	82	388468	10.658	ng	98
26) 2-Nitrophenol	7.76	139	27992	4.382	ng	97
27) 2,4-Dimethylphenol	7.79	122	156098	10.379	ng	97
28) bis(2-Chloroethoxy)methane	7.89	93	267919	10.862	ng	98
29) 2,4-Dichlorophenol	7.99	162	146829	9.551	ng	99
30) 1,2,4-Trichlorobenzene	8.08	180	201948	11.329	ng	99
31) Naphthalene	8.16	128	628418	11.633	ng	97
32) Benzoic acid	7.86	122	34546	4.082	ng	92
33) 4-Chloroaniline	8.21	127	263598	11.150	ng	99
34) Hexachlorobutadiene	8.29	225	121305	11.402	ng	98
35) Caprolactam	8.55	113	46026	8.720	ng	97
36) 4-Chloro-3-methylphenol	8.68	107	163289	10.238	ng	98
37) 2-Methylnaphthalene	8.86	142	412858	11.445	ng	98
38) 1-Methylnaphthalene	8.95	142	389110	11.407	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.02	216	192052	11.650	ng	100

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41) Hexachlorocyclopentadiene	9.01	237	62048	7.778	nq	100
43) 2,4,6-Trichlorophenol	9.13	196	99128	8.877	nq	99
44) 2,4,5-Trichlorophenol	9.16	196	105322	9.478	nq	97
46) 1,1'-Biphenyl	9.32	154	501316	11.882	nq	97
47) 2-Chloronaphthalene	9.34	162	401691	11.640	nq	98
48) 2-Nitroaniline	9.43	65	44531	5.025	nq	97
49) Acenaphthylene	9.76	152	606616	11.490	nq	97
50) Dimethylphthalate	9.62	163	438489	10.960	nq	98
51) 2,6-Dinitrotoluene	9.67	165	33949	4.785	nq #	85
52) Acenaphthene	9.93	154	376288	11.281	nq	99
53) 3-Nitroaniline	9.84	138	49401	5.514	nq #	78
54) 2,4-Dinitrophenol	9.94	184	8693	3.835	nq #	1
55) Dibenzofuran	10.10	168	564318	11.749	nq	96
56) 4-Nitrophenol	9.99	139	36807	5.607	nq	90
57) 2,4-Dinitrotoluene	10.07	165	35419	4.140	nq	94
58) Fluorene	10.45	166	442938	12.117	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.22	232	82213	8.608	nq	98
60) Diethylphthalate	10.32	149	448548	10.946	nq	97
61) 4-Chlorophenyl-phenylether	10.44	204	220229	12.005	nq	96
62) 4-Nitroaniline	10.45	138	61329	6.753	nq	85
63) Azobenzene	10.60	77	449684	11.347	nq	98
65) 4,6-Dinitro-2-methylphenol	10.47	198	11689	3.766	nq	75
66) n-Nitrosodiphenylamine	10.55	169	381561	11.397	nq	99
67) 4-Bromophenyl-phenylether	10.93	248	132370	11.323	nq	95
68) Hexachlorobenzene	10.99	284	153742	11.247	nq	97
69) Atrazine	11.07	200	110833	10.723	nq	99
70) Pentachlorophenol	11.18	266	51691	7.571	nq	98
71) Phenanthrene	11.40	178	654781	12.078	nq	97
72) Anthracene	11.46	178	644102	12.033	nq	96
73) Carbazole	11.61	167	608803	11.751	nq	97
74) Di-n-butylphthalate	11.95	149	700295	11.405	nq	98
75) Fluoranthene	12.59	202	715599	12.047	nq	97
77) Benzidine	12.72	184	241764	11.644	nq	99
78) Pyrene	12.82	202	738837	11.311	nq	96
80) Butylbenzylphthalate	13.44	149	245454	9.028	nq	93
81) Benzo(a)anthracene	14.01	228	661916	11.740	nq	99
82) 3,3'-Dichlorobenzidine	13.97	252	238739	11.446	nq	99
83) Chrysene	14.04	228	646307	11.555	nq	97
84) Bis(2-ethylhexyl)phthalate	14.01	149	380022	10.564	nq	98
85) Di-n-octyl phthalate	14.62	149	571743	9.133	nq	98
87) Indeno(1,2,3-cd)pyrene	16.94	276	633737	11.087	nq	98
88) Benzo(b)fluoranthene	15.06	252	631129	10.619	nq	98
89) Benzo(k)fluoranthene	15.09	252	649328	12.034	nq	97
90) Benzo(a)pyrene	15.42	252	584771	11.112	nq	99
91) Dibenzo(a,h)anthracene	16.96	278	523499	11.158	nq	98
92) Benzo(g,h,i)perylene	17.38	276	497974	11.042	nq	100

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

