

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070220\
 Data File : BF120761.D
 Acq On : 2 Jul 2020 16:32
 Operator : JU/CG
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

mohammad
 7/7/2020 8:25:21 AM

Quant Time: Jul 02 17:42:42 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF070220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 02 17:30:37 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	136264	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	504869	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	249171	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	450144	20.00	ng	0.00
76) Chrysene-d12	14.01	240	335174	20.00	ng	0.00
86) Perylene-d12	15.47	264	285792	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	707616	86.92	ng	0.00
7) Phenol-d6	6.47	99	894500	83.55	ng	0.00
23) Nitrobenzene-d5	7.42	82	725135	82.42	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	215997	69.84	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	1340811	75.85	ng	0.00
79) Terphenyl-d14	12.96	244	1362051	76.59	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.59	88	158698	45.572	ng	100
3) Pyridine	3.33	79	442983	44.854	ng	100
4) n-Nitrosodimethylamine	3.29	42	218034	54.966	ng	100
6) Aniline	6.51	93	577438	44.512	ng	100
8) 2-Chlorophenol	6.63	128	382473	43.205	ng	100
9) Benzaldehyde	6.40	77	241729	44.696	ng	100
10) Phenol	6.49	94	458641m	44.420	ng	
11) bis(2-Chloroethyl)ether	6.59	93	373399	42.473	ng	100
12) 1,3-Dichlorobenzene	6.79	146	415118	42.468	ng	100
13) 1,4-Dichlorobenzene	6.87	146	410838	41.879	ng	100
14) 1,2-Dichlorobenzene	7.02	146	394902	40.818	ng	100
15) Benzyl Alcohol	6.99	79	339377	50.396	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.13	45	585936	49.473	ng	100
17) 2-Methylphenol	7.10	107	333113	43.640	ng	100
18) Hexachloroethane	7.36	117	156030	44.142	ng	100
19) n-Nitroso-di-n-propylamine	7.26	70	265149	43.182	ng	100
20) 3+4-Methylphenols	7.25	107	419668	40.658	ng	100
22) Acetophenone	7.26	105	502558	39.226	ng	100
24) Nitrobenzene	7.43	77	404304	44.027	ng	100
25) Isophorone	7.67	82	744504	43.679	ng	100
26) 2-Nitrophenol	7.75	139	154097	31.370	ng	100
27) 2,4-Dimethylphenol	7.78	122	302055	41.017	ng	100
28) bis(2-Chloroethoxy)methane	7.89	93	443888	42.283	ng	100
29) 2,4-Dichlorophenol	7.99	162	303077	40.943	ng	100
30) 1,2,4-Trichlorobenzene	8.07	180	341499	39.389	ng	100
31) Naphthalene	8.16	128	1075862	42.709	ng	100
32) Benzoic acid	7.89	122	201817	39.002	ng	100
33) 4-Chloroaniline	8.20	127	456552	41.506	ng	100
34) Hexachlorobutadiene	8.27	225	207809	39.699	ng	100
35) Caprolactam	8.57	113	87805m	36.546	ng	
36) 4-Chloro-3-methylphenol	8.67	107	317450	42.802	ng	100
37) 2-Methylnaphthalene	8.85	142	710744	41.754	ng	100
38) 1-Methylnaphthalene	8.95	142	665956	41.711	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.01	216	313770	39.446	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.00	237	192923	53.072	nq	100
43) 2,4,6-Trichlorophenol	9.12	196	213942	39.875	nq	100
44) 2,4,5-Trichlorophenol	9.16	196	238165	39.504	nq	100
46) 1,1'-Biphenyl	9.31	154	830749	41.140	nq	100
47) 2-Chloronaphthalene	9.33	162	661249	40.693	nq	100
48) 2-Nitroaniline	9.43	65	187698	39.607	nq	100
49) Acenaphthylene	9.75	152	1013309	40.157	nq	100
50) Dimethylphthalate	9.61	163	736994	38.592	nq	100
51) 2,6-Dinitrotoluene	9.67	165	142009	33.635	nq	100
52) Acenaphthene	9.92	154	639675	44.292	nq	100
53) 3-Nitroaniline	9.84	138	171885	34.680	nq	100
54) 2,4-Dinitrophenol	9.94	184	42612	25.486	nq	100
55) Dibenzofuran	10.09	168	922535	41.904	nq	100
56) 4-Nitrophenol	9.99	139	129659	42.341	nq	100
57) 2,4-Dinitrotoluene	10.07	165	174868	33.872	nq	100
58) Fluorene	10.44	166	686235	40.044	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.21	232	175945m	38.633	nq	
60) Diethylphthalate	10.31	149	766980	40.860	nq	100
61) 4-Chlorophenyl-phenylether	10.43	204	345890	39.765	nq	100
62) 4-Nitroaniline	10.45	138	164120	33.112	nq	100
63) Azobenzene	10.59	77	773703	45.154	nq	100
65) 4,6-Dinitro-2-methylphenol	10.47	198	64844	26.100	nq	100
66) n-Nitrosodiphenylamine	10.55	169	618748	41.794	nq	100
67) 4-Bromophenyl-phenylether	10.92	248	216664	39.545	nq	100
68) Hexachlorobenzene	10.98	284	228192	39.250	nq	100
69) Atrazine	11.07	200	162493	40.774	nq	100
70) Pentachlorophenol	11.17	266	133408	68.548	nq	100
71) Phenanthrene	11.40	178	988922	42.241	nq	100
72) Anthracene	11.45	178	1009626	43.001	nq	100
73) Carbazole	11.60	167	958849	42.952	nq	100
74) Di-n-butylphthalate	11.94	149	1205434	44.638	nq	100
75) Fluoranthene	12.59	202	1052513	40.017	nq	100
77) Benzidine	12.70	184	409393	41.146	nq	100
78) Pyrene	12.82	202	1085285	39.824	nq	100
80) Butylbenzylphthalate	13.44	149	481643	41.911	nq	100
81) Benzo(a)anthracene	14.00	228	882502	40.921	nq	100
82) 3,3'-Dichlorobenzidine	13.97	252	307069	38.147	nq	100
83) Chrysene	14.04	228	896841	39.639	nq	100
84) Bis(2-ethylhexyl)phthalate	14.00	149	621506	43.373	nq	100
85) Di-n-octyl phthalate	14.62	149	1116108	41.111	nq	100
87) Indeno(1,2,3-cd)pyrene	16.93	276	776913	43.028	nq	100
88) Benzo(b)fluoranthene	15.05	252	805624	44.031	nq	100
89) Benzo(k)fluoranthene	15.08	252	782258	48.804	nq	100
90) Benzo(a)pyrene	15.41	252	717816	45.658	nq	100
91) Dibenzo(a,h)anthracene	16.94	278	652866	44.188	nq	100
92) Benzo(g,h,i)perylene	17.36	276	588511	39.810	nq	100

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

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