

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\
 Data File : BF118682.D
 Acq On : 24 Jan 2020 13:20
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 1/28/2020 8:12:07 AM

Quant Time: Jan 27 01:15:42 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 27 01:13:13 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	272966	20.00	ng	-0.02
21) Naphthalene-d8	8.15	136	1000852	20.00	ng	-0.02
39) Acenaphthene-d10	9.90	164	567733	20.00	ng	-0.03
64) Phenanthrene-d10	11.38	188	1071400	20.00	ng	-0.03
76) Chrysene-d12	14.02	240	828245	20.00	ng	-0.03
87) Perylene-d12	15.47	264	801996	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	1123446	76.44	ng	-0.02
7) Phenol-d6	6.49	99	1412888	73.96	ng	-0.02
23) Nitrobenzene-d5	7.43	82	1295311	72.39	ng	-0.02
42) 2,4,6-Tribromophenol	10.69	330	503956	76.75	ng	-0.02
45) 2-Fluorobiphenyl	9.22	172	2523113	73.05	ng	-0.02
79) Terphenyl-d14	12.96	244	3117538	82.10	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.60	88	270870	38.051	ng	95
3) Pyridine	3.36	79	759589	37.858	ng	96
4) n-Nitrosodimethylamine	3.30	42	279356m	32.476	ng	
6) Aniline	6.53	93	899398	35.756	ng	96
8) 2-Chlorophenol	6.65	128	628220	38.082	ng	96
9) Benzaldehyde	6.41	77	423974	36.494	ng	95
10) Phenol	6.50	94	793143	36.433	ng	96
11) bis(2-Chloroethyl)ether	6.60	93	592017	36.653	ng	96
12) 1,3-Dichlorobenzene	6.81	146	737401	38.190	ng	98
13) 1,4-Dichlorobenzene	6.88	146	751562	38.775	ng	98
14) 1,2-Dichlorobenzene	7.04	146	697518	38.282	ng	97
15) Benzyl Alcohol	7.00	79	547999	36.174	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.14	45	756766	33.638	ng	99
17) 2-Methylphenol	7.11	107	510122	37.123	ng	97
18) Hexachloroethane	7.38	117	260026	36.997	ng	97
19) n-Nitroso-di-n-propylamine	7.27	70	427639	33.154	ng	98
20) 3+4-Methylphenols	7.26	107	632008	37.671	ng	# 83
22) Acetophenone	7.27	105	844517	35.620	ng	# 95
24) Nitrobenzene	7.45	77	650278	35.492	ng	93
25) Isophorone	7.69	82	1153092	35.331	ng	97
26) 2-Nitrophenol	7.76	139	322355	41.258	ng	91
27) 2,4-Dimethylphenol	7.80	122	466867	34.604	ng	99
28) bis(2-Chloroethoxy)methane	7.89	93	751981	35.759	ng	99
29) 2,4-Dichlorophenol	8.00	162	557715	37.394	ng	99
30) 1,2,4-Trichlorobenzene	8.09	180	655207	37.888	ng	98
31) Naphthalene	8.17	128	1765369	39.047	ng	98
32) Benzoic acid	7.90	122	371720	36.711	ng	96
33) 4-Chloroaniline	8.21	127	786326	37.629	ng	98
34) Hexachlorobutadiene	8.29	225	393397	37.355	ng	99
35) Caprolactam	8.57	113	176229	36.106	ng	87
36) 4-Chloro-3-methylphenol	8.69	107	551787	37.465	ng	98
37) 2-Methylnaphthalene	8.86	142	1231317	38.583	ng	99
38) 1-Methylnaphthalene	8.96	142	1173340	38.700	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.03	216	669755	37.703	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.02	237	286876	32.008	nq	98
43) 2,4,6-Trichlorophenol	9.13	196	446164	37.849	nq	99
44) 2,4,5-Trichlorophenol	9.17	196	456577	37.925	nq	97
46) 1,1'-Biphenyl	9.32	154	1564363	39.981	nq	96
47) 2-Chloronaphthalene	9.35	162	1275798	38.684	nq	97
48) 2-Nitroaniline	9.43	65	354211	34.939	nq	93
49) Acenaphthylene	9.76	152	1820094	39.229	nq	96
50) Dimethylphthalate	9.62	163	1456089	39.366	nq	98
51) 2,6-Dinitrotoluene	9.68	165	326909	39.469	nq #	74
52) Acenaphthene	9.93	154	1194304	38.496	nq	97
53) 3-Nitroaniline	9.85	138	363736	38.459	nq	95
54) 2,4-Dinitrophenol	9.95	184	136605	41.312	nq #	32
55) Dibenzofuran	10.10	168	1707593	39.717	nq	95
56) 4-Nitrophenol	10.00	139	267760	37.503	nq	88
57) 2,4-Dinitrotoluene	10.08	165	423650	40.573	nq	93
58) Fluorene	10.45	166	1295555	38.384	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.22	232	402461	38.008	nq	98
60) Diethylphthalate	10.32	149	1423822	39.702	nq	99
61) 4-Chlorophenyl-phenylether	10.44	204	710487	38.704	nq	99
62) 4-Nitroaniline	10.46	138	352881	36.374	nq	94
63) Azobenzene	10.60	77	1266248	36.458	nq	94
65) 4,6-Dinitro-2-methylphenol	10.49	198	200503	42.559	nq	94
66) n-Nitrosodiphenylamine	10.56	169	1196895	37.043	nq	98
67) 4-Bromophenyl-phenylether	10.93	248	487199	36.807	nq	94
68) Hexachlorobenzene	11.00	284	542766	36.167	nq	97
69) Atrazine	11.08	200	401573	35.858	nq	99
70) Pentachlorophenol	11.19	266	306279	36.708	nq	97
71) Phenanthrene	11.41	178	2010344	38.893	nq	94
72) Anthracene	11.46	178	2010054	39.301	nq	95
73) Carbazole	11.61	167	1793165	38.223	nq	96
74) Di-n-butylphthalate	11.95	149	2192866	42.072	nq	97
75) Fluoranthene	12.59	202	2166093	39.732	nq	97
77) Benzidine	12.71	184	821435	37.119	nq	98
78) Pyrene	12.82	202	2200322	39.130	nq	95
80) Butylbenzylphthalate	13.44	149	946967	41.149	nq	92
81) Benzo(a)anthracene	14.01	228	1927280	38.682	nq	96
82) 3,3'-Dichlorobenzidine	13.97	252	771747	43.470	nq	98
83) Chrysene	14.04	228	1860340	38.959	nq	96
84) Bis(2-ethylhexyl)phthalate	14.00	149	1257862	44.739	nq	97
85) Di-n-octyl phthalate	14.61	149	2058816	49.472	nq	96
86) Indeno(1,2,3-cd)pyrene	16.94	276	1674495	40.358	nq	98
88) Benzo(b)fluoranthene	15.05	252	1821468	36.132	nq	98
89) Benzo(k)fluoranthene	15.08	252	1829826	39.179	nq	98
90) Benzo(a)pyrene	15.42	252	1644278	37.811	nq	98
91) Dibenzo(a,h)anthracene	16.96	278	1387662	36.075	nq	98
92) Benzo(g,h,i)perylene	17.39	276	1293998	34.646	nq	98

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

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