

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040120\
 Data File : BF119685.D
 Acq On : 1 Apr 2020 23:55
 Operator : MA/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 02 01:57:26 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 11:59:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	225683	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	853930	20.00	ng	0.00
39) Acenaphthene-d10	9.86	164	435207	20.00	ng	0.00
64) Phenanthrene-d10	11.35	188	851262	20.00	ng	0.00
76) Chrysene-d12	13.99	240	625125	20.00	ng	-0.01
87) Perylene-d12	15.44	264	603832	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	1092194	77.04	ng	0.00
7) Phenol-d6	6.45	99	1407959	77.74	ng	0.00
23) Nitrobenzene-d5	7.38	82	1248449	79.46	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	399082	88.88	ng	0.00
45) 2-Fluorobiphenyl	9.18	172	2269960	77.27	ng	0.00
79) Terphenyl-d14	12.93	244	2543884	79.73	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	253085	36.145	ng	97
3) Pyridine	3.23	79	694876	36.023	ng	94
4) n-Nitrosodimethylamine	3.19	42	306886	32.624	ng	89
6) Aniline	6.47	93	956369	38.263	ng	96
8) 2-Chlorophenol	6.60	128	601340	40.386	ng	99
9) Benzaldehyde	6.36	77	415358	36.154	ng	99
10) Phenol	6.46	94	803944	41.604	ng	95
11) bis(2-Chloroethyl)ether	6.55	93	602146	38.961	ng	99
12) 1,3-Dichlorobenzene	6.76	146	647783	40.197	ng	100
13) 1,4-Dichlorobenzene	6.83	146	655254	40.650	ng	98
14) 1,2-Dichlorobenzene	6.99	146	614089	40.758	ng	100
15) Benzyl Alcohol	6.96	79	530278	38.926	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.09	45	1122117	35.996	ng	100
17) 2-Methylphenol	7.06	107	539264	39.528	ng	99
18) Hexachloroethane	7.33	117	239707	40.579	ng	99
19) n-Nitroso-di-n-propylamine	7.23	70	414887	38.008	ng	97
20) 3+4-Methylphenols	7.22	107	649707	38.860	ng	# 82
22) Acetophenone	7.23	105	766568	37.569	ng	99
24) Nitrobenzene	7.40	77	644587	38.388	ng	98
25) Isophorone	7.64	82	1198257	37.595	ng	99
26) 2-Nitrophenol	7.72	139	318499	48.174	ng	# 86
27) 2,4-Dimethylphenol	7.75	122	492804	36.048	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	722653	38.342	ng	99
29) 2,4-Dichlorophenol	7.95	162	492472	39.205	ng	99
30) 1,2,4-Trichlorobenzene	8.04	180	549037	39.523	ng	98
31) Naphthalene	8.12	128	1737741	39.544	ng	99
32) Benzoic acid	7.86	122	312729	35.562	ng	97
33) 4-Chloroaniline	8.17	127	763449	37.915	ng	99
34) Hexachlorobutadiene	8.24	225	321384	41.349	ng	99
35) Caprolactam	8.54	113	159888	38.892	ng	97
36) 4-Chloro-3-methylphenol	8.65	107	530170	38.617	ng	97
37) 2-Methylnaphthalene	8.82	142	1128086	39.115	ng	97
38) 1-Methylnaphthalene	8.91	142	1074799	39.373	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.98	216	515826	40.437	ng	99

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41) Hexachlorocyclopentadiene	8.97	237	277343	38.243	nq	99
43) 2,4,6-Trichlorophenol	9.09	196	360123	39.450	nq	99
44) 2,4,5-Trichlorophenol	9.13	196	413908	40.475	nq	97
46) 1,1'-Biphenyl	9.28	154	1423650	40.164	nq	98
47) 2-Chloronaphthalene	9.30	162	1095673	39.463	nq	98
48) 2-Nitroaniline	9.40	65	398574	41.590	nq	91
49) Acenaphthylene	9.72	152	1710839	39.239	nq	98
50) Dimethylphthalate	9.58	163	1254984	40.597	nq	99
51) 2,6-Dinitrotoluene	9.64	165	282055	42.568	nq	91
52) Acenaphthene	9.89	154	1063739	39.135	nq	98
53) 3-Nitroaniline	9.81	138	347352	41.523	nq	92
54) 2,4-Dinitrophenol	9.91	184	91469	36.276	nq	98
55) Dibenzofuran	10.06	168	1543264	39.600	nq	98
56) 4-Nitrophenol	9.96	139	252697	38.293	nq	93
57) 2,4-Dinitrotoluene	10.05	165	387028	46.367	nq	# 84
58) Fluorene	10.40	166	1166142	40.465	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.18	232	309286	40.462	nq	97
60) Diethylphthalate	10.28	149	1285222	40.963	nq	99
61) 4-Chlorophenyl-phenylether	10.40	204	587282	41.672	nq	99
62) 4-Nitroaniline	10.42	138	345643	43.089	nq	95
63) Azobenzene	10.56	77	1221798	38.671	nq	99
65) 4,6-Dinitro-2-methylphenol	10.45	198	149245	41.810	nq	94
66) n-Nitrosodiphenylamine	10.52	169	1093213	39.119	nq	99
67) 4-Bromophenyl-phenylether	10.89	248	392054	40.440	nq	99
68) Hexachlorobenzene	10.96	284	402688	40.584	nq	# 91
69) Atrazine	11.05	200	312312	40.765	nq	98
70) Pentachlorophenol	11.15	266	217011	37.300	nq	98
71) Phenanthrene	11.37	178	1716566	38.889	nq	98
72) Anthracene	11.42	178	1756284	39.149	nq	99
73) Carbazole	11.57	167	1713831	38.596	nq	99
74) Di-n-butylphthalate	11.91	149	2017660	42.477	nq	99
75) Fluoranthene	12.56	202	1899841	39.770	nq	99
77) Benzidine	12.67	184	983399	37.172	nq	99
78) Pyrene	12.79	202	1948529	37.981	nq	99
80) Butylbenzylphthalate	13.41	149	885264	42.663	nq	97
81) Benzo(a)anthracene	13.97	228	1565155	38.502	nq	99
82) 3,3'-Dichlorobenzidine	13.94	252	624849	38.984	nq	99
83) Chrysene	14.02	228	1614143	38.475	nq	99
84) Bis(2-ethylhexyl)phthalate	13.97	149	1121145	45.325	nq	99
85) Di-n-octyl phthalate	14.58	149	2018595	46.401	nq	98
86) Indeno(1,2,3-cd)pyrene	16.89	276	1418095	35.890	nq	98
88) Benzo(b)fluoranthene	15.02	252	1607393	39.850	nq	98
89) Benzo(k)fluoranthene	15.05	252	1469723	38.266	nq	99
90) Benzo(a)pyrene	15.38	252	1412356	38.529	nq	98
91) Dibenzo(a,h)anthracene	16.91	278	1181313	36.844	nq	99
92) Benzo(g,h,i)perylene	17.33	276	1146540	35.595	nq	96

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

