

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100620\
 Data File : BF121841.D
 Acq On : 6 Oct 2020 9:01
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 06 11:50:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 14:29:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	130755	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	490199	20.00	ng	0.00
39) Acenaphthene-d10	9.80	164	250616	20.00	ng	0.00
64) Phenanthrene-d10	11.29	188	463536	20.00	ng	0.00
76) Chrysene-d12	13.93	240	358522	20.00	ng	0.00
86) Perylene-d12	15.36	264	317254	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	717131	81.51	ng	0.00
7) Phenol-d6	6.42	99	916734	79.42	ng	0.00
23) Nitrobenzene-d5	7.33	82	775424	84.93	ng	0.00
42) 2,4,6-Tribromophenol	10.60	330	260343	83.58	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	1304243	78.82	ng	0.00
79) Terphenyl-d14	12.87	244	1509975	85.32	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.47	88	159570	39.477	ng	99
3) Pyridine	3.20	79	449909	40.262	ng	99
4) n-Nitrosodimethylamine	3.16	42	203685	39.297	ng	100
6) Aniline	6.43	93	566318	37.670	ng	# 81
8) 2-Chlorophenol	6.56	128	362979	40.520	ng	95
9) Benzaldehyde	6.32	77	262899	34.840	ng	99
10) Phenol	6.43	94	462309	37.611	ng	90
11) bis(2-Chloroethyl)ether	6.50	93	374580	40.433	ng	99
12) 1,3-Dichlorobenzene	6.70	146	408196	40.826	ng	99
13) 1,4-Dichlorobenzene	6.78	146	408000	40.639	ng	99
14) 1,2-Dichlorobenzene	6.94	146	372590	40.286	ng	99
15) Benzyl Alcohol	6.92	79	311562	39.711	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.05	45	577816	38.759	ng	91
17) 2-Methylphenol	7.03	107	306887	40.263	ng	96
18) Hexachloroethane	7.27	117	150382	41.834	ng	99
19) n-Nitroso-di-n-propylamine	7.19	70	259954	39.076	ng	100
20) 3+4-Methylphenols	7.19	107	364813	39.837	ng	91
22) Acetophenone	7.18	105	494939	39.795	ng	96
24) Nitrobenzene	7.36	77	410915	41.614	ng	96
25) Isophorone	7.59	82	720741	39.216	ng	99
26) 2-Nitrophenol	7.67	139	197255	42.730	ng	96
27) 2,4-Dimethylphenol	7.71	122	327587	39.931	ng	99
28) bis(2-Chloroethoxy)methane	7.80	93	459290	40.368	ng	100
29) 2,4-Dichlorophenol	7.92	162	303890	40.670	ng	98
30) 1,2,4-Trichlorobenzene	7.99	180	319880	40.648	ng	98
31) Naphthalene	8.07	128	1027013	39.688	ng	99
32) Benzoic acid	7.86	122	156586	28.822	ng	99
33) 4-Chloroaniline	8.12	127	455041	40.567	ng	99
34) Hexachlorobutadiene	8.19	225	194738	42.159	ng	100
35) Caprolactam	8.51	113	98833	39.719	ng	97
36) 4-Chloro-3-methylphenol	8.62	107	325287	40.412	ng	99
37) 2-Methylnaphthalene	8.76	142	671876	39.620	ng	98
38) 1-Methylnaphthalene	8.86	142	623792	39.524	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.93	216	322023	41.135	ng	100

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41) Hexachlorocyclopentadiene	8.92	237	159814	35.748	nq	100
43) 2,4,6-Trichlorophenol	9.05	196	206982	38.790	nq	99
44) 2,4,5-Trichlorophenol	9.09	196	221375	39.244	nq	99
46) 1,1'-Biphenyl	9.23	154	802517	39.999	nq	99
47) 2-Chloronaphthalene	9.25	162	629314	39.804	nq	100
48) 2-Nitroaniline	9.35	65	214512	43.661	nq	97
49) Acenaphthylene	9.67	152	963954	39.681	nq	99
50) Dimethylphthalate	9.53	163	747996	41.180	nq	99
51) 2,6-Dinitrotoluene	9.60	165	166167	42.955	nq	86
52) Acenaphthene	9.84	154	568921	37.079	nq	99
53) 3-Nitroaniline	9.77	138	188648	42.097	nq	94
54) 2,4-Dinitrophenol	9.88	184	66951	35.661	nq	90
55) Dibenzofuran	10.01	168	828921	38.363	nq	98
56) 4-Nitrophenol	9.94	139	149617	41.865	nq	86
57) 2,4-Dinitrotoluene	10.00	165	208136	42.768	nq	92
58) Fluorene	10.35	166	669676	40.528	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.13	232	180035	39.258	nq	97
60) Diethylphthalate	10.23	149	721664	40.749	nq	99
61) 4-Chlorophenyl-phenylether	10.35	204	328356	41.483	nq	98
62) 4-Nitroaniline	10.38	138	186250	41.859	nq	96
63) Azobenzene	10.50	77	748909	39.710	nq	99
65) 4,6-Dinitro-2-methylphenol	10.42	198	116806	42.055	nq	99
66) n-Nitrosodiphenylamine	10.46	169	633860	39.858	nq	100
67) 4-Bromophenyl-phenylether	10.83	248	218861	41.470	nq	99
68) Hexachlorobenzene	10.90	284	242164	40.660	nq	# 91
69) Atrazine	10.99	200	164167	41.551	nq	99
70) Pentachlorophenol	11.10	266	113394	34.668	nq	99
71) Phenanthrene	11.32	178	992534	39.300	nq	100
72) Anthracene	11.36	178	1040358	39.918	nq	100
73) Carbazole	11.52	167	927943	39.455	nq	100
74) Di-n-butylphthalate	11.85	149	1118831	41.217	nq	99
75) Fluoranthene	12.50	202	1068209	39.622	nq	98
77) Benzidine	12.62	184	432340	36.375	nq	100
78) Pyrene	12.73	202	1096285	41.851	nq	100
80) Butylbenzylphthalate	13.34	149	468562	44.229	nq	99
81) Benzo(a)anthracene	13.92	228	937959	41.353	nq	99
82) 3,3'-Dichlorobenzidine	13.88	252	361661	40.842	nq	99
83) Chrysene	13.96	228	917243	39.935	nq	100
84) Bis(2-ethylhexyl)phthalate	13.90	149	641230	45.308	nq	98
85) Di-n-octyl phthalate	14.52	149	1066596	42.704	nq	100
87) Indeno(1,2,3-cd)pyrene	16.79	276	819591	39.669	nq	99
88) Benzo(b)fluoranthene	14.95	252	906467	42.740	nq	98
89) Benzo(k)fluoranthene	14.99	252	766398	39.461	nq	99
90) Benzo(a)pyrene	15.31	252	734680	40.752	nq	99
91) Dibenzo(a,h)anthracene	16.80	278	676644	39.343	nq	99
92) Benzo(g,h,i)perylene	17.22	276	677127	40.053	nq	99

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

