

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042720\
 Data File : BF120211.D
 Acq On : 27 Apr 2020 11:45
 Operator : MA/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 27 12:20:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 27 12:17:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	148334	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	542747	20.00	ng	0.00
39) Acenaphthene-d10	9.71	164	281416	20.00	ng	0.00
64) Phenanthrene-d10	11.19	188	551962	20.00	ng	0.00
76) Chrysene-d12	13.84	240	423308	20.00	ng	0.00
87) Perylene-d12	15.24	264	394735	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.28	112	751459	87.55	ng	0.00
7) Phenol-d6	6.33	99	931211	84.20	ng	0.00
23) Nitrobenzene-d5	7.24	82	872438	83.16	ng	0.00
42) 2,4,6-Tribromophenol	10.50	330	255107	74.04	ng	0.00
45) 2-Fluorobiphenyl	9.03	172	1647023	83.09	ng	0.00
79) Terphenyl-d14	12.79	244	1768174	70.29	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.28	88	139478	36.484	ng	99
3) Pyridine	2.96	79	411474	39.229	ng	99
4) n-Nitrosodimethylamine	2.92	42	209162	39.029	ng	98
6) Aniline	6.33	93	594563	40.958	ng	# 48
8) 2-Chlorophenol	6.46	128	406885	42.427	ng	99
9) Benzaldehyde	6.22	77	240589	40.976	ng	98
10) Phenol	6.34	94	523395	41.713	ng	85
11) bis(2-Chloroethyl)ether	6.41	93	397944	41.076	ng	99
12) 1,3-Dichlorobenzene	6.61	146	435638	41.492	ng	100
13) 1,4-Dichlorobenzene	6.69	146	435082	40.680	ng	99
14) 1,2-Dichlorobenzene	6.84	146	413719	41.973	ng	99
15) Benzyl Alcohol	6.82	79	346501	40.337	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.95	45	701683	37.220	ng	63
17) 2-Methylphenol	6.95	107	352213	41.153	ng	# 91
18) Hexachloroethane	7.18	117	163666	40.946	ng	94
19) n-Nitroso-di-n-propylamine	7.09	70	287573	40.434	ng	98
20) 3+4-Methylphenols	7.10	107	468665	44.774	ng	# 83
22) Acetophenone	7.09	105	548270	43.139	ng	98
24) Nitrobenzene	7.26	77	437306	41.436	ng	100
25) Isophorone	7.50	82	804393	39.774	ng	98
26) 2-Nitrophenol	7.57	139	222230	42.148	ng	91
27) 2,4-Dimethylphenol	7.62	122	327991	41.246	ng	98
28) bis(2-Chloroethoxy)methane	7.71	93	482726	40.563	ng	99
29) 2,4-Dichlorophenol	7.82	162	336986	41.342	ng	97
30) 1,2,4-Trichlorobenzene	7.90	180	369973	40.058	ng	100
31) Naphthalene	7.97	128	1188144	41.608	ng	98
32) Benzoic acid	7.77	122	233422	33.422	ng	95
33) 4-Chloroaniline	8.03	127	507334	40.811	ng	97
34) Hexachlorobutadiene	8.09	225	222153	40.182	ng	99
35) Caprolactam	8.41	113	100612	40.351	ng	96
36) 4-Chloro-3-methylphenol	8.53	107	360595	40.861	ng	97
37) 2-Methylnaphthalene	8.67	142	779586	40.268	ng	98
38) 1-Methylnaphthalene	8.77	142	735587	40.312	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.83	216	358100	40.289	ng	99

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41) Hexachlorocyclopentadiene	8.82	237	181612	35.933	ng	99
43) 2,4,6-Trichlorophenol	8.96	196	247992	40.404	ng	99
44) 2,4,5-Trichlorophenol	9.00	196	278948	40.351	ng	95
46) 1,1'-Biphenyl	9.13	154	990984	41.720	ng	99
47) 2-Chloronaphthalene	9.16	162	753734	41.192	ng	100
48) 2-Nitroaniline	9.26	65	278959	42.069	ng	91
49) Acenaphthylene	9.57	152	1157058	41.579	ng	98
50) Dimethylphthalate	9.44	163	886513	40.727	ng	100
51) 2,6-Dinitrotoluene	9.50	165	194569	40.819	ng	91
52) Acenaphthene	9.75	154	693772	37.374	ng	98
53) 3-Nitroaniline	9.67	138	229544	42.165	ng	95
54) 2,4-Dinitrophenol	9.79	184	122765	43.018	ng	97
55) Dibenzofuran	9.92	168	1040863	40.861	ng	100
56) 4-Nitrophenol	9.86	139	147996	36.537	ng	99
57) 2,4-Dinitrotoluene	9.91	165	254789	40.571	ng	95
58) Fluorene	10.26	166	837004	41.086	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.04	232	212966	40.280	ng	97
60) Diethylphthalate	10.14	149	902443	40.002	ng	99
61) 4-Chlorophenyl-phenylether	10.25	204	409374	39.206	ng	98
62) 4-Nitroaniline	10.29	138	235770	45.444	ng	95
63) Azobenzene	10.41	77	834829	41.291	ng	98
65) 4,6-Dinitro-2-methylphenol	10.32	198	139566	38.381	ng	93
66) n-Nitrosodiphenylamine	10.37	169	746549	40.669	ng	99
67) 4-Bromophenyl-phenylether	10.74	248	254864	37.129	ng	98
68) Hexachlorobenzene	10.81	284	260546	37.416	ng	92
69) Atrazine	10.90	200	203119	39.770	ng	99
70) Pentachlorophenol	11.01	266	137564	34.281	ng	99
71) Phenanthrene	11.22	178	1193406	40.625	ng	98
72) Anthracene	11.27	178	1229411	40.968	ng	98
73) Carbazole	11.43	167	1180584	41.601	ng	98
74) Di-n-butylphthalate	11.76	149	1458353	39.091	ng	100
75) Fluoranthene	12.41	202	1323126	41.744	ng	95
77) Benzidine	12.53	184	591044	41.306	ng	98
78) Pyrene	12.64	202	1355704	37.990	ng	97
80) Butylbenzylphthalate	13.26	149	634732	36.656	ng	96
81) Benzo(a)anthracene	13.83	228	1114830	40.033	ng	100
82) 3,3'-Dichlorobenzidine	13.79	252	416994	40.132	ng	100
83) Chrysene	13.86	228	1129821	39.929	ng	99
84) Bis(2-ethylhexyl)phthalate	13.82	149	829391	35.370	ng	98
85) Di-n-octyl phthalate	14.43	149	1451353	36.351	ng	99
86) Indeno(1,2,3-cd)pyrene	16.61	276	1009573	40.476	ng	100
88) Benzo(b)fluoranthene	14.85	252	1098728	38.693	ng	100
89) Benzo(k)fluoranthene	14.87	252	958175	39.108	ng	99
90) Benzo(a)pyrene	15.19	252	948866	39.742	ng	97
91) Dibenzo(a,h)anthracene	16.63	278	834810	39.473	ng	98
92) Benzo(g,h,i)perylene	17.02	276	820985	39.327	ng	96

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

