

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040620\
 Data File : BF119761.D
 Acq On : 6 Apr 2020 10:25
 Operator : MA/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 06 11:35:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 06 11:15:16 2020
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	69	0.00
2	1,4-Dioxane	0.621	0.558	10.1	60	0.00
3	Pyridine	1.709	1.497	12.4	58	0.00
4	n-Nitrosodimethylamine	0.834	0.734	12.0	59	0.00
5 S	2-Fluorophenol	1.256	1.268	-1.0	67	0.00
6	Aniline	2.215	2.059	7.0	61	0.00
7 S	Phenol-d6	1.605	1.544	3.8	63	0.00
8	2-Chlorophenol	1.320	1.356	-2.7	67	0.00
9	Benzaldehyde	1.018	0.852	16.3	56	0.00
10 C	Phenol	1.712	1.764	-3.0	68	0.00
11	bis(2-Chloroethyl)ether	1.370	1.326	3.2	64	0.00
12	1,3-Dichlorobenzene	1.428	1.456	-2.0	68	0.00
13 C	1,4-Dichlorobenzene	1.428	1.459	-2.2	68	0.00
14	1,2-Dichlorobenzene	1.335	1.365	-2.2	67	0.00
15	Benzyl Alcohol	1.207	1.168	3.2	62	0.00
16	2,2'-oxybis(1-Chloropropane	2.763	2.472	10.5	59	0.00
17	2-Methylphenol	1.209	1.206	0.2	66	0.00
18	Hexachloroethane	0.523	0.568	-8.6	72	0.00
19 P	n-Nitroso-di-n-propylamine	0.967	0.954	1.3	66	0.00
20	3+4-Methylphenols	1.482	1.520	-2.6	67	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	68	0.00
22	Acetophenone	0.478	0.490	-2.5	67	0.00
23 S	Nitrobenzene-d5	0.368	0.402	-9.2	71	0.00
24	Nitrobenzene	0.393	0.412	-4.8	69	0.00
25	Isophorone	0.746	0.738	1.1	66	0.00
26 C	2-Nitrophenol	0.155	0.191	-23.2#	80	0.00
27	2,4-Dimethylphenol	0.320	0.298	6.9	61	0.00
28	bis(2-Chloroethoxy)methane	0.441	0.404	8.4	61	0.00
29 C	2,4-Dichlorophenol	0.294	0.293	0.3	66	0.00
30	1,2,4-Trichlorobenzene	0.325	0.317	2.5	65	0.00
31	Naphthalene	1.029	1.046	-1.7	66	0.00
32	Benzoic acid	0.206	0.241	-17.0	77	0.00
33	4-Chloroaniline	0.472	0.441	6.6	61	0.00
34 C	Hexachlorobutadiene	0.182	0.188	-3.3	69	0.00
35	Caprolactam	0.096	0.093	3.1	62	0.00
36 C	4-Chloro-3-methylphenol	0.322	0.323	-0.3	66	0.00
37	2-Methylnaphthalene	0.675	0.681	-0.9	66	0.00
38	1-Methylnaphthalene	0.639	0.642	-0.5	65	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	65	0.00
40	1,2,4,5-Tetrachlorobenzene	0.586	0.631	-7.7	68	0.00
41 P	Hexachlorocyclopentadiene	0.333	0.365	-9.6	69	0.00
42 S	2,4,6-Tribromophenol	0.206	0.259	-25.7#	79	0.00
43 C	2,4,6-Trichlorophenol	0.420	0.479	-14.0	72	0.00
44	2,4,5-Trichlorophenol	0.470	0.533	-13.4	71	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040620\
 Data File : BF119761.D
 Acq On : 6 Apr 2020 10:25
 Operator : MA/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 06 11:35:19 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 06 11:15:16 2020
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
45 S	2-Fluorobiphenyl	1.350	1.442	-6.8	70	0.00
46	1,1'-Biphenyl	1.629	1.743	-7.0	67	0.00
47	2-Chloronaphthalene	1.276	1.341	-5.1	66	0.00
48	2-Nitroaniline	0.440	0.508	-15.5	70	0.00
49	Acenaphthylene	2.004	2.064	-3.0	64	0.00
50	Dimethylphthalate	1.421	1.489	-4.8	66	0.00
51	2,6-Dinitrotoluene	0.304	0.332	-9.2	67	0.00
52 C	Acenaphthene	1.249	1.306	-4.6	66	0.00
53	3-Nitroaniline	0.384	0.400	-4.2	64	0.00
54 P	2,4-Dinitrophenol	0.112	0.180	-60.7#	106	0.00
55	Dibenzofuran	1.791	1.850	-3.3	65	0.00
56 P	4-Nitrophenol	0.303	0.313	-3.3	62	0.00
57	2,4-Dinitrotoluene	0.384	0.457	-19.0	73	0.00
58	Fluorene	1.324	1.404	-6.0	66	0.00
59	2,3,4,6-Tetrachlorophenol	0.351	0.378	-7.7	66	0.00
60	Diethylphthalate	1.442	1.505	-4.4	66	0.00
61	4-Chlorophenyl-phenylether	0.648	0.705	-8.8	69	0.00
62	4-Nitroaniline	0.369	0.401	-8.7	67	0.00
63	Azobenzene	1.452	1.526	-5.1	66	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	65	0.00
65	4,6-Dinitro-2-methylphenol	0.084	0.135	-60.7#	101	0.00
66 c	n-Nitrosodiphenylamine	0.657	0.711	-8.2	69	0.00
67	4-Bromophenyl-phenylether	0.228	0.252	-10.5	70	0.00
68	Hexachlorobenzene	0.233	0.253	-8.6	69	0.00
69	Atrazine	0.180	0.175	2.8	62	0.00
70 C	Pentachlorophenol	0.137	0.146	-6.6	68	0.00
71	Phenanthrene	1.037	1.079	-4.1	66	0.00
72	Anthracene	1.054	1.075	-2.0	64	0.00
73	Carbazole	1.043	1.103	-5.8	66	0.00
74	Di-n-butylphthalate	1.116	1.298	-16.3	74	0.00
75 C	Fluoranthene	1.122	1.137	-1.3	64	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	73	0.00
77	Benzidine	0.846	0.630	25.5#	54	0.00
78	Pyrene	1.641	1.444	12.0	64	0.00
79 S	Terphenyl-d14	1.021	0.973	4.7	71	0.00
80	Butylbenzylphthalate	0.664	0.682	-2.7	76	0.00
81	Benzo(a)anthracene	1.301	1.282	1.5	70	0.00
82	3,3'-Dichlorobenzidine	0.513	0.526	-2.5	72	0.00
83	Chrysene	1.342	1.311	2.3	70	0.00
84	Bis(2-ethylhexyl)phthalate	0.791	0.857	-8.3	80	0.00
85 c	Di-n-octyl phthalate	1.392	1.570	-12.8	79	0.00
86	Indeno(1,2,3-cd)pyrene	1.264	1.349	-6.7	81	0.00
87 I	Perylene-d12	1.000	1.000	0.0	80	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040620\
Data File : BF119761.D
Acq On : 6 Apr 2020 10:25
Operator : MA/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Apr 06 11:35:19 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Apr 06 11:15:16 2020
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.336	1.289	3.5	76	0.00
89	Benzo(k)fluoranthene	1.272	1.201	5.6	72	0.00
90 C	Benzo(a)pyrene	1.214	1.182	2.6	76	0.00
91	Dibenzo(a,h)anthracene	1.062	1.037	2.4	81	0.00
92	Benzo(g,h,i)perylene	1.067	1.061	0.6	84	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 1