

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041620\
 Data File : BF120025.D
 Acq On : 16 Apr 2020 10:00
 Operator : MA/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 16 10:46:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 11:01:23 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	112	0.00
2	1,4-Dioxane	0.515	0.469	8.9	100	0.00
3	Pyridine	1.414	1.342	5.1	109	-0.01
4	n-Nitrosodimethylamine	0.723	0.670	7.3	106	-0.01
5 S	2-Fluorophenol	1.157	1.163	-0.5	115	0.00
6	Aniline	1.957	1.955	0.1	113	0.00
7 S	Phenol-d6	1.491	1.498	-0.5	111	0.00
8	2-Chlorophenol	1.293	1.293	0.0	113	0.00
9	Benzaldehyde	0.824	0.784	4.9	109	0.00
10 C	Phenol	1.692	1.670	1.3	111	0.00
11	bis(2-Chloroethyl)ether	1.306	1.302	0.3	113	0.00
12	1,3-Dichlorobenzene	1.416	1.394	1.6	112	0.00
13 C	1,4-Dichlorobenzene	1.442	1.400	2.9	112	0.00
14	1,2-Dichlorobenzene	1.329	1.314	1.1	110	0.00
15	Benzyl Alcohol	1.158	1.131	2.3	110	0.00
16	2,2'-oxybis(1-Chloropropane	2.542	2.462	3.1	108	0.00
17	2-Methylphenol	1.154	1.150	0.3	112	0.00
18	Hexachloroethane	0.539	0.537	0.4	112	0.00
19 P	n-Nitroso-di-n-propylamine	0.959	0.973	-1.5	111	0.00
20	3+4-Methylphenols	1.411	1.418	-0.5	108	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	115	0.00
22	Acetophenone	0.468	0.451	3.6	110	0.00
23 S	Nitrobenzene-d5	0.387	0.366	5.4	112	0.00
24	Nitrobenzene	0.389	0.373	4.1	113	0.00
25	Isophorone	0.745	0.712	4.4	107	0.00
26 C	2-Nitrophenol	0.194	0.189	2.6	105	0.00
27	2,4-Dimethylphenol	0.293	0.274	6.5	105	0.00
28	bis(2-Chloroethoxy)methane	0.439	0.426	3.0	113	0.00
29 C	2,4-Dichlorophenol	0.300	0.285	5.0	112	0.00
30	1,2,4-Trichlorobenzene	0.340	0.320	5.9	112	0.00
31	Naphthalene	1.052	0.995	5.4	113	0.00
32	Benzoic acid	0.257	0.233	9.3	107	0.00
33	4-Chloroaniline	0.458	0.433	5.5	114	0.00
34 C	Hexachlorobutadiene	0.204	0.195	4.4	112	0.00
35	Caprolactam	0.092	0.084	8.7	110	0.00
36 C	4-Chloro-3-methylphenol	0.325	0.309	4.9	113	0.00
37	2-Methylnaphthalene	0.713	0.673	5.6	113	0.00
38	1-Methylnaphthalene	0.672	0.620	7.7	103	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
40	1,2,4,5-Tetrachlorobenzene	0.632	0.592	6.3	104	0.00
41 P	Hexachlorocyclopentadiene	0.359	0.423	-17.8	131	0.00
42 S	2,4,6-Tribromophenol	0.245	0.215	12.2	99	0.00
43 C	2,4,6-Trichlorophenol	0.436	0.413	5.3	110	0.00
44	2,4,5-Trichlorophenol	0.491	0.468	4.7	105	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041620\
 Data File : BF120025.D
 Acq On : 16 Apr 2020 10:00
 Operator : MA/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Apr 16 10:46:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 11:01:23 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	1.409	1.349	4.3	107	0.00
46	1,1'-Biphenyl	1.688	1.608	4.7	108	0.00
47	2-Chloronaphthalene	1.300	1.233	5.2	110	0.00
48	2-Nitroaniline	0.471	0.448	4.9	113	0.00
49	Acenaphthylene	1.978	1.853	6.3	109	0.00
50	Dimethylphthalate	1.547	1.458	5.8	111	0.00
51	2,6-Dinitrotoluene	0.339	0.321	5.3	111	0.00
52 C	Acenaphthene	1.319	1.149	12.9	104	0.00
53	3-Nitroaniline	0.387	0.371	4.1	117	0.00
54 P	2,4-Dinitrophenol	0.203	0.248	-22.2	143	-0.01
55	Dibenzofuran	1.810	1.708	5.6	111	0.00
56 P	4-Nitrophenol	0.288	0.266	7.6	110	0.00
57	2,4-Dinitrotoluene	0.446	0.422	5.4	112	0.00
58	Fluorene	1.448	1.308	9.7	111	0.00
59	2,3,4,6-Tetrachlorophenol	0.376	0.364	3.2	113	0.00
60	Diethylphthalate	1.603	1.540	3.9	115	0.00
61	4-Chlorophenyl-phenylether	0.742	0.674	9.2	110	0.00
62	4-Nitroaniline	0.369	0.368	0.3	117	0.00
63	Azobenzene	1.437	1.386	3.5	113	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	120	0.00
65	4,6-Dinitro-2-methylphenol	0.132	0.129	2.3	116	0.00
66 c	n-Nitrosodiphenylamine	0.665	0.627	5.7	114	0.00
67	4-Bromophenyl-phenylether	0.249	0.225	9.6	107	0.00
68	Hexachlorobenzene	0.252	0.226	10.3	106	0.00
69	Atrazine	0.185	0.171	7.6	105	0.00
70 C	Pentachlorophenol	0.145	0.126	13.1	100	0.00
71	Phenanthrene	1.064	0.982	7.7	112	0.00
72	Anthracene	1.087	0.999	8.1	110	0.00
73	Carbazole	1.028	0.954	7.2	111	0.00
74	Di-n-butylphthalate	1.352	1.258	7.0	109	0.00
75 C	Fluoranthene	1.148	1.061	7.6	114	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	137	0.00
77	Benzidine	0.676	0.598	11.5	126	0.00
78	Pyrene	1.686	1.513	10.3	113	0.00
79 S	Terphenyl-d14	1.189	1.010	15.1	107	0.00
80	Butylbenzylphthalate	0.818	0.733	10.4	117	0.00
81	Benzo(a)anthracene	1.316	1.213	7.8	128	0.00
82	3,3'-Dichlorobenzidine	0.491	0.456	7.1	127	0.00
83	Chrysene	1.337	1.242	7.1	130	0.00
84	Bis(2-ethylhexyl)phthalate	1.108	0.986	11.0	123	0.00
85 c	Di-n-octyl phthalate	1.886	1.740	7.7	127	0.00
86	Indeno(1,2,3-cd)pyrene	1.178	1.023	13.2	111	-0.01
87 I	Perylene-d12	1.000	1.000	0.0	139	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041620\
Data File : BF120025.D
Acq On : 16 Apr 2020 10:00
Operator : MA/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Apr 16 10:46:55 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 15 11:01:23 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	1.439	1.273	11.5	132	0.00
89	Benzo(k)fluoranthene	1.241	1.221	1.6	120	0.00
90 C	Benzo(a)pyrene	1.210	1.137	6.0	132	0.00
91	Dibenzo(a,h)anthracene	1.072	0.943	12.0	111	-0.01
92	Benzo(q,h,i)perylene	1.058	0.876	17.2	109	-0.01

(#) = Out of Range

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