

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022420\
 Data File : BF119052.D
 Acq On : 25 Feb 2020 7:14
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 25 07:43:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 20 17:40:17 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	100	0.00
2	1,4-Dioxane	0.533	0.516	3.2	103	-0.02
3	Pyridine	1.455	1.405	3.4	103	-0.02
4	n-Nitrosodimethylamine	0.441	0.455	-3.2	109	-0.02
5 S	2-Fluorophenol	1.197	1.220	-1.9	106	0.00
6	Aniline	1.796	1.747	2.7	100	0.00
7 S	Phenol-d6	1.455	1.454	0.1	104	0.00
8	2-Chlorophenol	1.292	1.312	-1.5	105	0.00
9	Benzaldehyde	0.972	0.927	4.6	100	0.00
10 C	Phenol	1.574	1.549	1.6	103	0.00
11	bis(2-Chloroethyl)ether	1.204	1.209	-0.4	106	0.00
12	1,3-Dichlorobenzene	1.548	1.570	-1.4	107	0.00
13 C	1,4-Dichlorobenzene	1.549	1.576	-1.7	106	0.00
14	1,2-Dichlorobenzene	1.413	1.440	-1.9	106	0.00
15	Benzyl Alcohol	1.052	1.081	-2.8	105	0.00
16	2,2'-oxybis(1-Chloropropane	1.043	1.040	0.3	106	0.00
17	2-Methylphenol	1.047	1.036	1.1	104	0.00
18	Hexachloroethane	0.554	0.565	-2.0	108	0.00
19 P	n-Nitroso-di-n-propylamine	0.848	0.829	2.2	104	0.00
20	3+4-Methylphenols	1.283	1.263	1.6	106	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
22	Acetophenone	0.463	0.473	-2.2	105	0.00
23 S	Nitrobenzene-d5	0.379	0.388	-2.4	105	0.00
24	Nitrobenzene	0.375	0.380	-1.3	105	0.00
25	Isophorone	0.658	0.649	1.4	103	0.00
26 C	2-Nitrophenol	0.198	0.204	-3.0	106	0.00
27	2,4-Dimethylphenol	0.269	0.271	-0.7	104	0.00
28	bis(2-Chloroethoxy)methane	0.428	0.432	-0.9	104	0.00
29 C	2,4-Dichlorophenol	0.317	0.319	-0.6	102	0.00
30	1,2,4-Trichlorobenzene	0.376	0.382	-1.6	104	0.00
31	Naphthalene	1.037	1.048	-1.1	103	0.00
32	Benzoic acid	0.184	0.180	2.2	102	0.00
33	4-Chloroaniline	0.446	0.438	1.8	100	0.00
34 C	Hexachlorobutadiene	0.234	0.243	-3.8	107	0.00
35	Caprolactam	0.098	0.093	5.1	96	0.00
36 C	4-Chloro-3-methylphenol	0.321	0.312	2.8	99	0.00
37	2-Methylnaphthalene	0.698	0.696	0.3	102	0.00
38	1-Methylnaphthalene	0.656	0.655	0.2	101	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
40	1,2,4,5-Tetrachlorobenzene	0.674	0.694	-3.0	102	0.00
41 P	Hexachlorocyclopentadiene	0.198	0.190	4.0	96	0.00
42 S	2,4,6-Tribromophenol	0.262	0.254	3.1	96	0.00
43 C	2,4,6-Trichlorophenol	0.426	0.442	-3.8	103	0.00
44	2,4,5-Trichlorophenol	0.447	0.448	-0.2	99	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022420\
 Data File : BF119052.D
 Acq On : 25 Feb 2020 7:14
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 25 07:43:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 20 17:40:17 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	1.358	1.357	0.1	102	0.00
46	1,1'-Biphenyl	1.616	1.659	-2.7	101	0.00
47	2-Chloronaphthalene	1.303	1.329	-2.0	101	0.00
48	2-Nitroaniline	0.363	0.361	0.6	98	0.00
49	Acenaphthylene	1.881	1.901	-1.1	99	0.00
50	Dimethylphthalate	1.529	1.526	0.2	100	0.00
51	2,6-Dinitrotoluene	0.340	0.332	2.4	97	0.00
52 C	Acenaphthene	1.134	1.146	-1.1	99	0.00
53	3-Nitroaniline	0.377	0.358	5.0	93	0.00
54 P	2,4-Dinitrophenol	0.140	0.127	9.3	88	0.00
55	Dibenzofuran	1.693	1.695	-0.1	97	0.00
56 P	4-Nitrophenol	0.226	0.198	12.4	84	0.00
57	2,4-Dinitrotoluene	0.440	0.427	3.0	94	0.00
58	Fluorene	1.316	1.311	0.4	97	0.00
59	2,3,4,6-Tetrachlorophenol	0.377	0.370	1.9	98	0.00
60	Diethylphthalate	1.441	1.421	1.4	97	0.00
61	4-Chlorophenyl-phenylether	0.696	0.704	-1.1	99	0.00
62	4-Nitroaniline	0.385	0.358	7.0	91	0.00
63	Azobenzene	1.253	1.250	0.2	99	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
65	4,6-Dinitro-2-methylphenol	0.121	0.120	0.8	91	0.00
66 c	n-Nitrosodiphenylamine	0.655	0.684	-4.4	95	0.00
67	4-Bromophenyl-phenylether	0.261	0.279	-6.9	98	0.00
68	Hexachlorobenzene	0.301	0.316	-5.0	96	0.00
69	Atrazine	0.229	0.215	6.1	86	0.00
70 C	Pentachlorophenol	0.121	0.122	-0.8	93	0.00
71	Phenanthrene	1.091	1.113	-2.0	93	0.00
72	Anthracene	1.090	1.121	-2.8	92	0.00
73	Carbazole	0.971	0.950	2.2	88	0.00
74	Di-n-butylphthalate	1.176	1.222	-3.9	94	0.00
75 C	Fluoranthene	1.158	1.098	5.2	85	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	70	0.00
77	Benzidine	0.813	0.718	11.7	64	0.00
78	Pyrene	1.606	1.720	-7.1	82	0.00
79 S	Terphenyl-d14	1.162	1.273	-9.6	85	0.00
80	Butylbenzylphthalate	0.676	0.731	-8.1	82	0.00
81	Benzo(a)anthracene	1.363	1.403	-2.9	77	0.00
82	3,3'-Dichlorobenzidine	0.529	0.553	-4.5	75	0.00
83	Chrysene	1.358	1.383	-1.8	77	0.00
84	Bis(2-ethylhexyl)phthalate	0.854	0.957	-12.1	84	0.00
85 c	Di-n-octyl phthalate	1.361	1.482	-8.9	77	0.00
86	Indeno(1,2,3-cd)pyrene	1.315	1.833	-39.4#	81	0.01
87 I	Perylene-d12	1.000	1.000	0.0	56	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022420\
Data File : BF119052.D
Acq On : 25 Feb 2020 7:14
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 25 07:43:30 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 20 17:40:17 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	1.318	1.251	5.1	80	0.00
89	Benzo(k)fluoranthene	1.222	1.231	-0.7	78	0.00
90 C	Benzo(a)pyrene	1.190	1.192	-0.2	60	0.00
91	Dibenzo(a,h)anthracene	1.047	1.307	-24.8	79	0.01
92	Benzo(g,h,i)perylene	1.034	1.340	-29.6#	84	0.01

(#) = Out of Range

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