Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103020\
Data File : BF122176.D
Acq On : 30 Oct 2020 12:08
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 30 13:11:35 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 30 10:16:54 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)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	71	0.00
2		1,4-Dioxane	0.596	0.603	-1.2	73	0.00
3		Pyridine	1.567	1.461	6.8	66	0.00
4		n-Nitrosodimethylamine	0.757	0.773	-2.1	73	0.00
5	S	2-Fluorophenol	1.355	1.397	-3.1	76	0.00
6		Aniline	2.148	2.103	2.1	73	0.00
7	S	Phenol-d6	1.744	1.722	1.3	73	0.00
8		2-Chlorophenol	1.370	1.367	0.2	73	0.00
9		Benzaldehyde	1.006	0.948	5.8	71	0.00
10	C	Phenol	1.800	1.772	1.6	73	0.00
11		bis(2-Chloroethyl)ether	1.392	1.356	2.6	71	0.00
12		1,3-Dichlorobenzene	1.541	1.502	2.5	72	-0.01
13	C	1,4-Dichlorobenzene	1.547	1.540	0.5	73	0.00
14		1,2-Dichlorobenzene	1.414	1.395	1.3	72	0.00
15		Benzyl Alcohol	1.128	1.184	-5.0	76	0.00
16		2,2'-oxybis(1-Chloropropane	2.142	1.977	7.7	67	0.00
17		2-Methylphenol	1.171	1.113	5.0	71	0.00
18		Hexachloroethane	0.559	0.550	1.6	71	0.00
19	P	n-Nitroso-di-n-propylamine	0.993	0.990	0.3	75	0.00
20		3+4-Methylphenols	1.412	1.482	-5.0	79	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	71	0.00
22		Acetophenone	0.515	0.527	-2.3	76	0.00
23	S	Nitrobenzene-d5	0.390	0.396	-1.5	73	0.00
24		Nitrobenzene	0.417	0.423	-1.4	73	0.00
25		Isophorone	0.737	0.732	0.7	73	0.00
26	C	2-Nitrophenol	0.192	0.194	-1.0	71	0.00
27		2,4-Dimethylphenol	0.327	0.320	2.1	71	0.00
28		bis(2-Chloroethoxy)methane	0.461	0.449	2.6	70	0.00
29	C	2,4-Dichlorophenol	0.306	0.308	-0.7	72	0.00
30		1,2,4-Trichlorobenzene	0.326	0.320	1.8	71	0.00
31		Naphthalene	1.071	1.059	1.1	73	0.00
32		Benzoic acid	0.165	0.154	6.7	65	0.00
33		4-Chloroaniline	0.456	0.465	-2.0	74	0.00
34	C	Hexachlorobutadiene	0.193	0.191	1.0	70	0.00
35		Caprolactam	0.097	0.101	-4.1	75	0.00
36	C	4-Chloro-3-methylphenol	0.329	0.343	-4.3	76	0.00
37		2-Methylnaphthalene	0.694	0.682	1.7	72	0.00
38		1-Methylnaphthalene	0.643	0.628	2.3	72	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	72	0.00
40		1,2,4,5-Tetrachlorobenzene	0.646	0.631	2.3	72	0.00
41	P	Hexachlorocyclopentadiene	0.127	0.098	22.8	54	0.00
42	S	2,4,6-Tribromophenol	0.247	0.247	0.0	73	0.00
43	C	2,4,6-Trichlorophenol	0.398	0.382	4.0	70	0.00
44		2,4,5-Trichlorophenol	0.430	0.379	11.9	65	0.00
45	S	2-Fluorobiphenyl	1.359	1.299	4.4	73	0.00
46		1,1'-Biphenyl	1.626	1.599	1.7	74	0.00
47		2-Chloronaphthalene	1.265	1.254	0.9	73	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103020\
Data File : BF122176.D
Acq On : 30 Oct 2020 12:08
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 30 13:11:35 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 30 10:16:54 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)
48		2-Nitroaniline	0.417	0.449	-7.7	79	0.00
49		Acenaphthylene	1.942	1.880	3.2	72	0.00
50		Dimethylphthalate	1.462	1.411	3.5	71	0.00
51		2,6-Dinitrotoluene	0.321	0.322	-0.3	72	0.00
52	C	Acenaphthene	1.155	1.149	0.5	75	0.00
53		3-Nitroaniline	0.365	0.371	-1.6	75	0.00
54	P	2,4-Dinitrophenol	0.129	0.173	-34.1#	94	0.00
55		Dibenzofuran	1.690	1.694	-0.2	75	0.00
56	P	4-Nitrophenol	0.198	0.216	-9.1	78	0.00
57		2,4-Dinitrotoluene	0.395	0.424	-7.3	78	0.00
58		Fluorene	1.361	1.378	-1.2	76	0.00
59		2,3,4,6-Tetrachlorophenol	0.333	0.323	3.0	69	0.00
60		Diethylphthalate	1.410	1.412	-0.1	74	0.00
61		4-Chlorophenyl-phenylether	0.653	0.650	0.5	75	0.00
62		4-Nitroaniline	0.365	0.346	5.2	70	0.00
63		Azobenzene	1.491	1.490	0.1	73	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	74	0.00
65		4,6-Dinitro-2-methylphenol	0.115	0.115	0.0	68	0.00
66	c	n-Nitrosodiphenylamine	0.689	0.670	2.8	75	0.00
67		4-Bromophenyl-phenylether	0.231	0.224	3.0	73	0.00
68		Hexachlorobenzene	0.262	0.252	3.8	73	0.00
69		Atrazine	0.168	0.161	4.2	71	0.00
70	C	Pentachlorophenol	0.084	0.073	13.1	65	0.00
71		Phenanthrene	1.097	1.083	1.3	76	0.00
72		Anthracene	1.137	1.133	0.4	77	0.00
73		Carbazole	1.011	1.024	-1.3	77	0.00
74		Di-n-butylphthalate	1.163	1.141	1.9	74	0.00
75	C	Fluoranthene	1.176	1.186	-0.9	77	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	76	0.00
77		Benzidine	0.617	0.498	19.3	64	0.00
78		Pyrene	1.524	1.512	0.8	77	0.00
79	S	Terphenyl-d14	1.049	1.038	1.0	79	0.00
80		Butylbenzylphthalate	0.604	0.598	1.0	75	0.00
81		Benzo(a)anthracene	1.311	1.317	-0.5	79	0.00
82		3,3'-Dichlorobenzidine	0.486	0.472	2.9	74	0.00
83		Chrysene	1.287	1.266	1.6	76	0.00
84		Bis(2-ethylhexyl)phthalate	0.821	0.809	1.5	75	0.00
85	c	Di-n-octyl phthalate	1.327	1.298	2.2	74	0.00
86	I	Perylene-d12	1.000	1.000	0.0	73	0.01
87		Indeno(1,2,3-cd)pyrene	1.299	1.306	-0.5	75	0.01
88		Benzo(b)fluoranthene	1.352	1.495	-10.6	78	0.00
89		Benzo(k)fluoranthene	1.284	1.328	-3.4	74	0.00
90	C	Benzo(a)pyrene	1.168	1.233	-5.6	75	0.00
91		Dibenzo(a,h)anthracene	1.069	1.060	0.8	73	0.01
92		Benzo(g,h,i)perylene	1.070	1.074	-0.4	75	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103020\
Data File : BF122176.D
Acq On : 30 Oct 2020 12:08
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Oct 30 13:11:35 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 30 10:16:54 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)

(#) = Out of Range SPCC's out = 0 CCC's out = 0