Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111320\
Data File : BF122325.D
Acq On : 13 Nov 2020 12:54
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 13 14:01:26 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 10 12:25:42 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	116	0.00
2		1,4-Dioxane	0.597	0.586	1.8	116	0.00
3		Pyridine	1.643	1.624	1.2	115	0.01
4		n-Nitrosodimethylamine	0.774	0.763	1.4	115	0.01
5	S	2-Fluorophenol	1.303	1.294	0.7	116	0.00
6		Aniline	2.236	2.188	2.1	114	0.00
7	S	Phenol-d6	1.704	1.679	1.5	115	0.00
8		2-Chlorophenol	1.351	1.355	-0.3	115	0.00
9		Benzaldehyde	0.961	0.852	11.3	110	0.00
10	C	Phenol	1.678	1.655	1.4	115	0.00
11		bis(2-Chloroethyl)ether	1.334	1.282	3.9	113	0.00
12		1,3-Dichlorobenzene	1.533	1.479	3.5	112	0.00
13	C	1,4-Dichlorobenzene	1.540	1.502	2.5	114	0.00
14		1,2-Dichlorobenzene	1.438	1.415	1.6	115	0.00
15		Benzyl Alcohol	1.211	1.208	0.2	114	0.00
16		2,2'-oxybis(1-Chloropropane	1.944	1.823	6.2	110	0.00
17		2-Methylphenol	1.127	1.103	2.1	115	0.00
18		Hexachloroethane	0.536	0.532	0.7	114	0.00
19	P	n-Nitroso-di-n-propylamine	1.018	0.956	6.1	110	0.00
20		3+4-Methylphenols	1.387	1.372	1.1	115	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	114	0.00
22		Acetophenone	0.521	0.504	3.3	113	0.00
23	S	Nitrobenzene-d5	0.327	0.373	-14.1	123	0.00
24		Nitrobenzene	0.374	0.397	-6.1	117	0.00
25		Isophorone	0.728	0.688	5.5	111	0.00
26	C	2-Nitrophenol	0.126	0.177	-40.5#	142	0.00
27		2,4-Dimethylphenol	0.344	0.345	-0.3	116	0.00
28		bis(2-Chloroethoxy)methane	0.445	0.428	3.8	111	0.00
29	C	2,4-Dichlorophenol	0.304	0.314	-3.3	117	0.00
30		1,2,4-Trichlorobenzene	0.336	0.332	1.2	113	0.00
31		Naphthalene	1.063	1.038	2.4	113	0.00
32		Benzoic acid	0.155	0.212	-36.8#	142	0.00
33		4-Chloroaniline	0.463	0.455	1.7	113	0.00
34	C	Hexachlorobutadiene	0.193	0.189	2.1	113	0.00
35		Caprolactam	0.096	0.098	-2.1	115	0.00
36	C	4-Chloro-3-methylphenol	0.317	0.317	0.0	114	0.00
37		2-Methylnaphthalene	0.702	0.682	2.8	113	0.00
38		1-Methylnaphthalene	0.658	0.635	3.5	111	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	113	0.00
40		1,2,4,5-Tetrachlorobenzene	0.616	0.616	0.0	114	0.00
41	P	Hexachlorocyclopentadiene	0.360	0.382	-6.1	114	0.00
42	S	2,4,6-Tribromophenol	0.215	0.234	-8.8	118	0.00
43	C	2,4,6-Trichlorophenol	0.400	0.430	-7.5	118	0.00
44		2,4,5-Trichlorophenol	0.420	0.441	-5.0	116	0.00
45	S	2-Fluorobiphenyl	1.280	1.210	5.5	111	0.00
46		1,1'-Biphenyl	1.592	1.552	2.5	112	0.00
47		2-Chloronaphthalene	1.264	1.262	0.2	114	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111320\
Data File : BF122325.D
Acq On : 13 Nov 2020 12:54
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 13 14:01:26 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 10 12:25:42 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	0.310	0.389	-25.5#	127	0.00
49		Acenaphthylene	1.942	1.896	2.4	111	0.00
50		Dimethylphthalate	1.422	1.378	3.1	113	0.00
51		2,6-Dinitrotoluene	0.275	0.325	-18.2	124	0.00
52	C	Acenaphthene	1.202	1.208	-0.5	115	0.00
53		3-Nitroaniline	0.317	0.363	-14.5	121	0.00
54	P	2,4-Dinitrophenol	0.084	0.123	-46.4#	174#	0.00
55		Dibenzofuran	1.761	1.728	1.9	113	0.00
56	P	4-Nitrophenol	0.271	0.298	-10.0	123	0.00
57		2,4-Dinitrotoluene	0.332	0.425	-28.0#	129	0.00
58		Fluorene	1.349	1.294	4.1	111	0.00
59		2,3,4,6-Tetrachlorophenol	0.349	0.377	-8.0	117	0.00
60		Diethylphthalate	1.395	1.346	3.5	111	0.00
61		4-Chlorophenyl-phenylether	0.637	0.620	2.7	112	0.00
62		4-Nitroaniline	0.313	0.362	-15.7	123	0.00
63		Azobenzene	1.463	1.394	4.7	110	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
65		4,6-Dinitro-2-methylphenol	0.076	0.104	-36.8#	154#	0.00
66	c	n-Nitrosodiphenylamine	0.681	0.675	0.9	113	0.00
67		4-Bromophenyl-phenylether	0.235	0.236	-0.4	113	0.00
68		Hexachlorobenzene	0.254	0.255	-0.4	114	0.00
69		Atrazine	0.170	0.172	-1.2	113	0.00
70	C	Pentachlorophenol	0.146	0.161	-10.3	121	0.00
71		Phenanthrene	1.135	1.111	2.1	112	0.00
72		Anthracene	1.170	1.151	1.6	111	0.00
73		Carbazole	1.064	1.036	2.6	111	0.00
74		Di-n-butylphthalate	1.134	1.099	3.1	109	0.00
75	C	Fluoranthene	1.185	1.155	2.5	110	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	106	0.00
77		Benzidine	0.555	0.530	4.5	101	0.00
78		Pyrene	1.405	1.418	-0.9	110	0.00
79	S	Terphenyl-d14	0.944	0.936	0.8	111	0.00
80		Butylbenzylphthalate	0.505	0.552	-9.3	111	0.00
81		Benzo(a)anthracene	1.244	1.230	1.1	107	0.00
82		3,3'-Dichlorobenzidine	0.464	0.479	-3.2	108	0.00
83		Chrysene	1.253	1.228	2.0	107	0.00
84		Bis(2-ethylhexyl)phthalate	0.677	0.693	-2.4	110	0.00
85	c	Di-n-octyl phthalate	1.148	1.185	-3.2	104	0.00
86	I	Perylene-d12	1.000	1.000	0.0	95	0.00
87		Indeno(1,2,3-cd)pyrene	1.200	1.152	4.0	96	0.00
88		Benzo(b)fluoranthene	1.295	1.345	-3.9	98	0.00
89		Benzo(k)fluoranthene	1.263	1.262	0.1	98	0.00
90	C	Benzo(a)pyrene	1.129	1.139	-0.9	96	0.00
91		Dibenzo(a,h)anthracene	1.005	0.969	3.6	97	0.00
92		Benzo(g,h,i)perylene	0.978	0.938	4.1	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111320\
Data File : BF122325.D
Acq On : 13 Nov 2020 12:54
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Nov 13 14:01:26 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110920.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 10 12:25:42 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 = 1