Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100821\
Data File : BF125775.D
Acq On : 8 Oct 2021 10:17
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 08 12:00:30 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 07 17:16:23 2021
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	111	0.00
2		1,4-Dioxane	0.521	0.502	3.6	111	0.02
3		Pyridine	1.349	1.349	0.0	111	0.02
4		n-Nitrosodimethylamine	0.485	0.474	2.3	110	0.02
5	S	2-Fluorophenol	1.224	1.172	4.2	110	0.00
6		Aniline	1.815	1.801	0.8	112	0.00
7	S	Phenol-d6	1.576	1.509	4.3	111	0.00
8		2-Chlorophenol	1.341	1.296	3.4	110	0.00
9		Benzaldehyde	0.855	0.857	-0.2	108	0.00
10	C	Phenol	1.533	1.457	5.0	109	0.00
11		bis(2-Chloroethyl)ether	1.242	1.206	2.9	111	0.00
12		1,3-Dichlorobenzene	1.463	1.414	3.3	111	0.00
13	C	1,4-Dichlorobenzene	1.492	1.413	5.3	108	0.00
14		1,2-Dichlorobenzene	1.397	1.340	4.1	110	0.00
15		Benzyl Alcohol	1.044	1.025	1.8	110	0.00
16		2,2'-oxybis(1-Chloropropane	1.588	1.536	3.3	110	0.00
17		2-Methylphenol	1.068	1.040	2.6	112	0.00
18		Hexachloroethane	0.514	0.497	3.3	109	0.00
19	P	n-Nitroso-di-n-propylamine	0.832	0.804	3.4	111	0.00
20		3+4-Methylphenols	1.317	1.262	4.2	110	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22		Acetophenone	0.440	0.414	5.9	112	0.00
23	S	Nitrobenzene-d5	0.322	0.315	2.2	111	0.00
24		Nitrobenzene	0.312	0.305	2.2	112	0.00
25		Isophorone	0.570	0.553	3.0	112	0.00
26	C	2-Nitrophenol	0.158	0.162	-2.5	110	0.00
27		2,4-Dimethylphenol	0.263	0.271	-3.0	117	0.00
28		bis(2-Chloroethoxy)methane	0.396	0.380	4.0	111	0.00
29	C	2,4-Dichlorophenol	0.283	0.278	1.8	113	0.00
30		1,2,4-Trichlorobenzene	0.310	0.299	3.5	112	0.00
31		Naphthalene	0.967	0.917	5.2	110	0.00
32		Benzoic acid	0.185	0.196	-5.9	113	0.01
33		4-Chloroaniline	0.405	0.397	2.0	113	0.00
34	C	Hexachlorobutadiene	0.173	0.167	3.5	111	0.00
35		Caprolactam	0.082	0.084	-2.4	116	0.00
36	C	4-Chloro-3-methylphenol	0.276	0.271	1.8	112	0.00
37		2-Methylnaphthalene	0.626	0.604	3.5	112	0.00
38		1-Methylnaphthalene	0.608	0.580	4.6	111	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
40		1,2,4,5-Tetrachlorobenzene	0.537	0.510	5.0	109	0.00
41	P	Hexachlorocyclopentadiene	0.263	0.267	-1.5	110	0.00
42	S	2,4,6-Tribromophenol	0.197	0.198	-0.5	115	0.00
43	C	2,4,6-Trichlorophenol	0.389	0.377	3.1	111	0.00
44		2,4,5-Trichlorophenol	0.414	0.406	1.9	111	0.00
45	S	2-Fluorobiphenyl	1.160	1.123	3.2	111	0.00
46		1,1'-Biphenyl	1.470	1.405	4.4	112	0.00
47		2-Chloronaphthalene	1.194	1.147	3.9	112	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100821\
Data File : BF125775.D
Acq On : 8 Oct 2021 10:17
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 08 12:00:30 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 07 17:16:23 2021
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	0.287	0.296	-3.1	114	0.00
49		Acenaphthylene	1.748	1.695	3.0	112	0.00
50		Dimethylphthalate	1.317	1.263	4.1	111	0.00
51		2,6-Dinitrotoluene	0.266	0.274	-3.0	115	0.00
52	C	Acenaphthene	1.092	1.044	4.4	111	0.00
53		3-Nitroaniline	0.315	0.326	-3.5	116	0.00
54	P	2,4-Dinitrophenol	0.113	0.119	-5.3	113	0.00
55		Dibenzofuran	1.636	1.589	2.9	113	0.00
56	P	4-Nitrophenol	0.235	0.249	-6.0	116	0.00
57		2,4-Dinitrotoluene	0.336	0.354	-5.4	115	0.00
58		Fluorene	1.199	1.133	5.5	113	0.00
59		2,3,4,6-Tetrachlorophenol	0.312	0.314	-0.6	117	0.00
60		Diethylphthalate	1.256	1.247	0.7	115	0.00
61		4-Chlorophenyl-phenylether	0.577	0.541	6.2	111	0.00
62		4-Nitroaniline	0.291	0.310	-6.5	120	0.00
63		Azobenzene	1.181	1.154	2.3	113	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
65		4,6-Dinitro-2-methylphenol	0.100	0.110	-10.0	120	0.00
66	c	n-Nitrosodiphenylamine	0.682	0.644	5.6	113	0.00
67		4-Bromophenyl-phenylether	0.221	0.211	4.5	114	0.00
68		Hexachlorobenzene	0.251	0.242	3.6	118	0.00
69		Atrazine	0.198	0.197	0.5	118	0.00
70	C	Pentachlorophenol	0.134	0.140	-4.5	118	0.00
71		Phenanthrene	1.070	1.013	5.3	115	0.00
72		Anthracene	1.069	1.010	5.5	114	0.00
73		Carbazole	0.995	0.976	1.9	119	0.00
74		Di-n-butylphthalate	1.098	1.102	-0.4	120	0.00
75	C	Fluoranthene	0.966	0.969	-0.3	123	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	131	0.00
77		Benzidine	0.732	0.767	-4.8	136	0.00
78		Pyrene	2.032	1.951	4.0	124	0.00
79	S	Terphenyl-d14	1.313	1.288	1.9	125	0.00
80		Butylbenzylphthalate	0.651	0.749	-15.1	150	0.00
81		Benzo(a)anthracene	1.298	1.253	3.5	129	0.00
82		3,3'-Dichlorobenzidine	0.410	0.401	2.2	123	0.00
83		Chrysene	1.283	1.246	2.9	129	0.00
84		Bis(2-ethylhexyl)phthalate	0.709	0.775	-9.3	141	0.00
85	c	Di-n-octyl phthalate	1.181	1.089	7.8	111	0.00
86	I	Perylene-d12	1.000	1.000	0.0	102	0.00
87		Indeno(1,2,3-cd)pyrene	1.526	1.529	-0.2	104	0.00
88		Benzo(b)fluoranthene	1.129	1.145	-1.4	104	0.00
89		Benzo(k)fluoranthene	1.159	1.139	1.7	103	0.00
90	C	Benzo(a)pyrene	1.000	0.999	0.1	102	0.00
91		Dibenzo(a,h)anthracene	1.259	1.266	-0.6	103	0.00
92		Benzo(g,h,i)perylene	1.286	1.289	-0.2	104	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100821\
Data File : BF125775.D
Acq On : 8 Oct 2021 10:17
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Oct 08 12:00:30 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF100721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 07 17:16:23 2021
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