Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125935.D
Acq On : 26 Oct 2021 15:33
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 26 17:09:53 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 25 18:49:16 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	121	0.00
2		1,4-Dioxane	0.656	0.630	4.0	122	0.00
3		Pyridine	1.737	1.731	0.3	119	0.00
4		n-Nitrosodimethylamine	1.114	1.022	8.3	111	0.00
5	S	2-Fluorophenol	1.343	1.285	4.3	121	0.00
6		Aniline	2.169	2.154	0.7	120	0.00
7	S	Phenol-d6	1.842	1.695	8.0	121	0.00
8		2-Chlorophenol	1.355	1.318	2.7	121	0.00
9		Benzaldehyde	1.102	1.036	6.0	120	0.00
10	C	Phenol	1.776	1.715	3.4	120	0.00
11		bis(2-Chloroethyl)ether	1.430	1.397	2.3	121	0.00
12		1,3-Dichlorobenzene	1.497	1.416	5.4	117	0.00
13	C	1,4-Dichlorobenzene	1.498	1.429	4.6	119	0.00
14		1,2-Dichlorobenzene	1.452	1.339	7.8	121	0.00
15		Benzyl Alcohol	1.394	1.384	0.7	120	0.00
16		2,2'-oxybis(1-Chloropropane	3.502	3.306	5.6	117	0.00
17		2-Methylphenol	1.171	1.173	-0.2	123	0.00
18		Hexachloroethane	0.568	0.562	1.1	123	0.00
19	P	n-Nitroso-di-n-propylamine	1.052	1.046	0.6	125	0.00
20		3+4-Methylphenols	1.496	1.403	6.2	121	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
22		Acetophenone	0.516	0.482	6.6	123	0.00
23	S	Nitrobenzene-d5	0.387	0.398	-2.8	127	0.00
24		Nitrobenzene	0.400	0.402	-0.5	124	0.00
25		Isophorone	0.737	0.710	3.7	124	0.00
26	C	2-Nitrophenol	0.130	0.167	-28.5#	147	0.00
27		2,4-Dimethylphenol	0.288	0.273	5.2	122	0.00
28		bis(2-Chloroethoxy)methane	0.463	0.443	4.3	127	0.00
29	C	2,4-Dichlorophenol	0.282	0.277	1.8	125	0.00
30		1,2,4-Trichlorobenzene	0.331	0.318	3.9	124	0.00
31		Naphthalene	1.017	0.959	5.7	124	0.00
32		Benzoic acid	0.174	0.201	-15.5	126	0.00
33		4-Chloroaniline	0.419	0.405	3.3	126	0.00
34	C	Hexachlorobutadiene	0.208	0.196	5.8	122	0.00
35		Caprolactam	0.092	0.096	-4.3	129	0.00
36	C	4-Chloro-3-methylphenol	0.330	0.320	3.0	125	0.00
37		2-Methylnaphthalene	0.655	0.619	5.5	125	0.00
38		1-Methylnaphthalene	0.632	0.605	4.3	125	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
40		1,2,4,5-Tetrachlorobenzene	0.551	0.533	3.3	127	0.00
41	P	Hexachlorocyclopentadiene	0.298	0.317	-6.4	131	0.00
42	S	2,4,6-Tribromophenol	0.174	0.183	-5.2	131	0.00
43	C	2,4,6-Trichlorophenol	0.376	0.385	-2.4	130	0.00
44		2,4,5-Trichlorophenol	0.397	0.396	0.3	126	0.00
45	S	2-Fluorobiphenyl	1.131	1.103	2.5	125	0.00
46		1,1'-Biphenyl	1.465	1.411	3.7	127	0.00
47		2-Chloronaphthalene	1.104	1.053	4.6	124	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125935.D
Acq On : 26 Oct 2021 15:33
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 26 17:09:53 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 25 18:49:16 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	0.394	0.457	-16.0	136	0.00
49		Acenaphthylene	1.668	1.618	3.0	127	0.00
50		Dimethylphthalate	1.239	1.213	2.1	126	0.00
51		2,6-Dinitrotoluene	0.244	0.282	-15.6	138	0.00
52	C	Acenaphthene	1.084	1.081	0.3	132	0.00
53		3-Nitroaniline	0.280	0.313	-11.8	136	0.00
54	P	2,4-Dinitrophenol	0.085	0.132	-55.3#	187#	0.00
55		Dibenzofuran	1.619	1.550	4.3	127	0.00
56	P	4-Nitrophenol	0.218	0.258	-18.3	139	0.00
57		2,4-Dinitrotoluene	0.300	0.372	-24.0	142	0.00
58		Fluorene	1.207	1.091	9.6	129	0.00
59		2,3,4,6-Tetrachlorophenol	0.324	0.338	-4.3	134	0.00
60		Diethylphthalate	1.264	1.280	-1.3	130	0.00
61		4-Chlorophenyl-phenylether	0.596	0.541	9.2	129	0.00
62		4-Nitroaniline	0.257	0.292	-13.6	139	0.00
63		Azobenzene	1.458	1.450	0.5	129	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	129	0.00
65		4,6-Dinitro-2-methylphenol	0.079	0.117	-48.1#	178#	0.00
66	c	n-Nitrosodiphenylamine	0.599	0.582	2.8	130	0.00
67		4-Bromophenyl-phenylether	0.214	0.213	0.5	132	0.00
68		Hexachlorobenzene	0.225	0.221	1.8	132	0.00
69		Atrazine	0.190	0.199	-4.7	134	0.00
70	C	Pentachlorophenol	0.127	0.146	-15.0	142	0.00
71		Phenanthrene	1.038	0.990	4.6	127	0.00
72		Anthracene	1.036	0.999	3.6	129	0.00
73		Carbazole	0.985	0.942	4.4	127	0.00
74		Di-n-butylphthalate	1.049	1.111	-5.9	135	0.00
75	C	Fluoranthene	1.046	1.028	1.7	130	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	132	0.00
77		Benzidine	0.624	0.656	-5.1	147	0.00
78		Pyrene	1.861	1.760	5.4	127	0.00
79	S	Terphenyl-d14	1.223	1.193	2.5	134	0.00
80		Butylbenzylphthalate	0.605	0.731	-20.8	153#	0.00
81		Benzo(a)anthracene	1.346	1.306	3.0	132	0.00
82		3,3'-Dichlorobenzidine	0.427	0.452	-5.9	145	0.00
83		Chrysene	1.362	1.347	1.1	138	0.00
84		Bis(2-ethylhexyl)phthalate	0.656	0.806	-22.9	164#	0.00
85	c	Di-n-octyl phthalate	1.016	1.290	-27.0#	169#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	138	0.00
87		Indeno(1,2,3-cd)pyrene	1.441	1.460	-1.3	129	0.00
88		Benzo(b)fluoranthene	1.320	1.370	-3.8	142	0.00
89		Benzo(k)fluoranthene	1.268	1.247	1.7	141	0.00
90	C	Benzo(a)pyrene	1.066	1.079	-1.2	137	0.00
91		Dibenzo(a,h)anthracene	1.167	1.183	-1.4	130	0.00
92		Benzo(g,h,i)perylene	1.220	1.216	0.3	126	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125935.D
Acq On : 26 Oct 2021 15:33
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Oct 26 17:09:53 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 25 18:49:16 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 = 2