Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121121\
Data File : BG051478.D
Acq On : 11 Dec 2021 12:17
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Dec 11 23:05:48 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 11 23:03:39 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	107	0.00
2		1,4-Dioxane	0.543	0.578	-6.4	116	0.00
3		Pyridine	1.614	1.650	-2.2	109	0.00
4		n-Nitrosodimethylamine	0.694	0.795	-14.6	123	0.00
5	S	2-Fluorophenol	1.294	1.281	1.0	109	0.00
6		Aniline	2.165	2.145	0.9	107	0.00
7	S	Phenol-d6	1.822	1.848	-1.4	110	0.00
8		2-Chlorophenol	1.358	1.351	0.5	110	0.00
9		Benzaldehyde	1.318	1.129	14.3	102	0.00
10	C	Phenol	1.871	1.904	-1.8	111	0.00
11		bis(2-Chloroethyl)ether	1.414	1.521	-7.6	118	0.00
12		1,3-Dichlorobenzene	1.453	1.455	-0.1	110	0.00
13	C	1,4-Dichlorobenzene	1.506	1.496	0.7	109	0.00
14		1,2-Dichlorobenzene	1.429	1.419	0.7	109	0.00
15		Benzyl Alcohol	1.456	1.624	-11.5	118	0.00
16		2,2'-oxybis(1-Chloropropane	2.044	2.248	-10.0	121	0.00
17		2-Methylphenol	1.263	1.273	-0.8	110	0.00
18		Hexachloroethane	0.563	0.589	-4.6	116	0.00
19	P	n-Nitroso-di-n-propylamine	1.345	1.412	-5.0	114	0.00
20		3+4-Methylphenols	1.800	1.789	0.6	106	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.540	0.547	-1.3	113	0.00
23	S	Nitrobenzene-d5	0.438	0.463	-5.7	117	0.00
24		Nitrobenzene	0.425	0.454	-6.8	118	0.00
25		Isophorone	0.789	0.814	-3.2	113	0.00
26	C	2-Nitrophenol	0.192	0.192	0.0	109	0.00
27		2,4-Dimethylphenol	0.284	0.280	1.4	111	0.00
28		bis(2-Chloroethoxy)methane	0.470	0.475	-1.1	112	0.00
29	C	2,4-Dichlorophenol	0.320	0.310	3.1	107	0.00
30		1,2,4-Trichlorobenzene	0.360	0.350	2.8	108	0.00
31		Naphthalene	1.055	1.053	0.2	110	0.00
32		Benzoic acid	0.184	0.156	15.2	88	0.00
33		4-Chloroaniline	0.454	0.439	3.3	107	0.00
34	C	Hexachlorobutadiene	0.222	0.212	4.5	108	0.00
35		Caprolactam	0.086	0.079	8.1	99	0.00
36	C	4-Chloro-3-methylphenol	0.396	0.384	3.0	106	0.00
37		2-Methylnaphthalene	0.748	0.728	2.7	108	0.00
38		1-Methylnaphthalene	0.730	0.709	2.9	108	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
40		1,2,4,5-Tetrachlorobenzene	0.608	0.590	3.0	106	0.00
41	P	Hexachlorocyclopentadiene	0.124	0.289	-133.1#	230#	0.00
42	S	2,4,6-Tribromophenol	0.213	0.198	7.0	98	0.00
43	C	2,4,6-Trichlorophenol	0.419	0.417	0.5	107	0.00
44		2,4,5-Trichlorophenol	0.453	0.418	7.7	96	0.00
45	S	2-Fluorobiphenyl	1.325	1.269	4.2	106	0.00
46		1,1'-Biphenyl	1.438	1.397	2.9	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121121\
Data File : BG051478.D
Acq On : 11 Dec 2021 12:17
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 11 23:05:48 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 11 23:03:39 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)
47		2-Chloronaphthalene	1.138	1.108	2.6	106	0.00
48		2-Nitroaniline	0.411	0.439	-6.8	113	0.00
49		Acenaphthylene	1.815	1.771	2.4	106	0.00
50		Dimethylphthalate	1.541	1.516	1.6	105	0.00
51		2,6-Dinitrotoluene	0.322	0.322	0.0	107	0.00
52	C	Acenaphthene	1.059	1.042	1.6	106	0.00
53		3-Nitroaniline	0.361	0.351	2.8	101	0.00
54	P	2,4-Dinitrophenol	0.162	0.162	0.0	97	0.00
55		Dibenzofuran	1.839	1.764	4.1	105	0.00
56	P	4-Nitrophenol	0.254	0.221	13.0	84	0.00
57		2,4-Dinitrotoluene	0.461	0.454	1.5	103	0.00
58		Fluorene	1.444	1.403	2.8	105	0.00
59		2,3,4,6-Tetrachlorophenol	0.389	0.374	3.9	100	0.00
60		Diethylphthalate	1.625	1.624	0.1	108	0.00
61		4-Chlorophenyl-phenylether	0.789	0.767	2.8	105	0.00
62		4-Nitroaniline	0.386	0.347	10.1	94	0.00
63		Azobenzene	1.627	1.711	-5.2	114	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	0.119	0.130	-9.2	106	0.00
66	c	n-Nitrosodiphenylamine	0.574	0.555	3.3	103	0.00
67		4-Bromophenyl-phenylether	0.216	0.208	3.7	103	0.00
68		Hexachlorobenzene	0.220	0.206	6.4	102	0.00
69		Atrazine	0.153	0.140	8.5	98	0.00
70	C	Pentachlorophenol	0.100	0.106	-6.0	103	0.00
71		Phenanthrene	1.075	1.056	1.8	105	0.00
72		Anthracene	1.066	1.039	2.5	104	0.00
73		Carbazole	1.049	1.021	2.7	103	0.00
74		Di-n-butylphthalate	1.270	1.250	1.6	104	0.00
75	C	Fluoranthene	1.325	1.288	2.8	103	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
77		Benzidine	0.328	0.191	41.8#	52	0.00
78		Pyrene	1.450	1.505	-3.8	103	0.00
79	S	Terphenyl-d14	1.094	1.083	1.0	99	0.00
80		Butylbenzylphthalate	0.618	0.637	-3.1	104	0.00
81		Benzo(a)anthracene	1.364	1.357	0.5	101	0.00
82		3,3'-Dichlorobenzidine	0.610	0.578	5.2	95	0.00
83		Chrysene	1.281	1.291	-0.8	102	0.00
84		Bis(2-ethylhexyl)phthalate	0.861	0.883	-2.6	104	0.00
85	c	Di-n-octyl phthalate	1.443	1.498	-3.8	106	0.00
86	I	Perylene-d12	1.000	1.000	0.0	98	0.00
87		Indeno(1,2,3-cd)pyrene	1.435	1.383	3.6	98	0.00
88		Benzo(b)fluoranthene	1.233	1.256	-1.9	105	0.00
89		Benzo(k)fluoranthene	1.251	1.227	1.9	99	0.00
90	C	Benzo(a)pyrene	1.070	1.066	0.4	102	0.00
91		Dibenzo(a,h)anthracene	1.182	1.120	5.2	97	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121121\
Data File : BG051478.D
Acq On : 11 Dec 2021 12:17
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Dec 11 23:05:48 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 11 23:03:39 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)
92	Benzo(g,h,i)perylene	1.169	1.120	4.2	97	0.00

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