Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061422\
Data File : BG053899.D
Acq On : 14 Jun 2022 12:21
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jun 15 02:58:18 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 13 15:49:31 2022
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	103	0.00
2		1,4-Dioxane	0.457	0.473	-3.5	105	0.00
3		Pyridine	1.220	1.307	-7.1	106	0.00
4		n-Nitrosodimethylamine	0.537	0.561	-4.5	102	0.00
5	S	2-Fluorophenol	1.046	1.085	-3.7	105	0.00
6		Aniline	1.682	1.723	-2.4	103	0.00
7	S	Phenol-d6	1.419	1.479	-4.2	106	0.00
8		2-Chlorophenol	1.117	1.159	-3.8	105	0.00
9		Benzaldehyde	0.838	0.787	6.1	104	0.00
10	C	Phenol	1.438	1.483	-3.1	102	0.00
11		bis(2-Chloroethyl)ether	1.104	1.124	-1.8	105	0.00
12		1,3-Dichlorobenzene	1.400	1.391	0.6	104	0.00
13	C	1,4-Dichlorobenzene	1.406	1.410	-0.3	106	0.00
14		1,2-Dichlorobenzene	1.350	1.362	-0.9	104	0.00
15		Benzyl Alcohol	1.061	1.100	-3.7	103	0.00
16		2,2'-oxybis(1-Chloropropane	1.650	1.681	-1.9	104	0.00
17		2-Methylphenol	0.999	1.005	-0.6	102	0.00
18		Hexachloroethane	0.466	0.484	-3.9	106	0.00
19	P	n-Nitroso-di-n-propylamine	0.953	0.960	-0.7	102	0.00
20		3+4-Methylphenols	1.401	1.417	-1.1	102	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.486	0.491	-1.0	102	0.00
23	S	Nitrobenzene-d5	0.356	0.367	-3.1	104	0.00
24		Nitrobenzene	0.350	0.362	-3.4	104	0.00
25		Isophorone	0.665	0.672	-1.1	103	0.00
26	C	2-Nitrophenol	0.155	0.159	-2.6	101	0.00
27		2,4-Dimethylphenol	0.251	0.255	-1.6	102	0.00
28		bis(2-Chloroethoxy)methane	0.358	0.361	-0.8	103	0.00
29	C	2,4-Dichlorophenol	0.336	0.355	-5.7	106	0.00
30		1,2,4-Trichlorobenzene	0.416	0.419	-0.7	102	0.00
31		Naphthalene	1.010	1.006	0.4	101	0.00
32		Benzoic acid	0.087	0.081	6.9	90	0.00
33		4-Chloroaniline	0.419	0.427	-1.9	102	0.00
34	C	Hexachlorobutadiene	0.320	0.327	-2.2	102	0.00
35		Caprolactam	0.092	0.098	-6.5	101	0.00
36	C	4-Chloro-3-methylphenol	0.327	0.331	-1.2	100	0.00
37		2-Methylnaphthalene	0.747	0.752	-0.7	102	0.00
38		1-Methylnaphthalene	0.727	0.725	0.3	102	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
40		1,2,4,5-Tetrachlorobenzene	0.771	0.771	0.0	102	0.00
41	P	Hexachlorocyclopentadiene	0.359	0.369	-2.8	102	0.00
42	S	2,4,6-Tribromophenol	0.321	0.322	-0.3	100	0.00
43	C	2,4,6-Trichlorophenol	0.433	0.440	-1.6	100	0.00
44		2,4,5-Trichlorophenol	0.483	0.493	-2.1	101	0.00
45	S	2-Fluorobiphenyl	1.375	1.367	0.6	102	0.00
46		1,1'-Biphenyl	1.393	1.362	2.2	99	0.00
47		2-Chloronaphthalene	1.127	1.128	-0.1	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061422\
Data File : BG053899.D
Acq On : 14 Jun 2022 12:21
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jun 15 02:58:18 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 13 15:49:31 2022
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	0.300	0.310	-3.3	101	0.00
49		Acenaphthylene	1.753	1.753	0.0	101	0.00
50		Dimethylphthalate	1.523	1.516	0.5	102	0.00
51		2,6-Dinitrotoluene	0.308	0.324	-5.2	104	0.00
52	C	Acenaphthene	1.118	1.118	0.0	101	0.00
53		3-Nitroaniline	0.298	0.307	-3.0	100	0.00
54	P	2,4-Dinitrophenol	0.140	0.139	0.7	99	0.00
55		Dibenzofuran	1.801	1.796	0.3	102	0.00
56	P	4-Nitrophenol	0.212	0.233	-9.9	108	0.00
57		2,4-Dinitrotoluene	0.435	0.455	-4.6	103	0.00
58		Fluorene	1.478	1.489	-0.7	103	0.00
59		2,3,4,6-Tetrachlorophenol	0.473	0.474	-0.2	99	0.00
60		Diethylphthalate	1.475	1.487	-0.8	104	0.00
61		4-Chlorophenyl-phenylether	0.885	0.869	1.8	101	0.00
62		4-Nitroaniline	0.311	0.320	-2.9	103	0.00
63		Azobenzene	1.187	1.212	-2.1	105	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
65		4,6-Dinitro-2-methylphenol	0.100	0.103	-3.0	103	0.00
66	c	n-Nitrosodiphenylamine	0.521	0.525	-0.8	102	0.00
67		4-Bromophenyl-phenylether	0.250	0.251	-0.4	102	0.00
68		Hexachlorobenzene	0.286	0.282	1.4	101	0.00
69		Atrazine	0.216	0.222	-2.8	102	0.00
70	C	Pentachlorophenol	0.133	0.132	0.8	98	0.00
71		Phenanthrene	1.028	1.031	-0.3	102	0.00
72		Anthracene	1.003	1.015	-1.2	103	0.00
73		Carbazole	0.861	0.879	-2.1	104	0.00
74		Di-n-butylphthalate	0.996	1.029	-3.3	104	0.00
75	C	Fluoranthene	1.349	1.363	-1.0	103	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
77		Benzidine	0.414	0.459	-10.9	114	0.00
78		Pyrene	1.208	1.200	0.7	103	0.00
79	S	Terphenyl-d14	0.993	0.978	1.5	102	0.00
80		Butylbenzylphthalate	0.387	0.402	-3.9	106	0.00
81		Benzo(a)anthracene	1.295	1.294	0.1	105	0.00
82		3,3'-Dichlorobenzidine	0.387	0.396	-2.3	105	0.00
83		Chrysene	1.234	1.208	2.1	104	0.00
84		Bis(2-ethylhexyl)phthalate	0.554	0.571	-3.1	106	0.00
85	c	Di-n-octyl phthalate	0.952	0.992	-4.2	107	0.00
86	I	Perylene-d12	1.000	1.000	0.0	105	0.01
87		Indeno(1,2,3-cd)pyrene	1.509	1.535	-1.7	105	0.00
88		Benzo(b)fluoranthene	1.216	1.199	1.4	101	0.00
89		Benzo(k)fluoranthene	1.204	1.230	-2.2	109	0.00
90	C	Benzo(a)pyrene	1.026	1.041	-1.5	105	0.00
91		Dibenzo(a,h)anthracene	1.248	1.257	-0.7	105	0.00
92		Benzo(g,h,i)perylene	1.232	1.248	-1.3	105	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061422\
Data File : BG053899.D
Acq On : 14 Jun 2022 12:21
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jun 15 02:58:18 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 13 15:49:31 2022
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