Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG052720\
Data File : BG045473.D
Acq On : 27 May 2020 12:52
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 27 13:45:17 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG051520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 27 13:40:45 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	120	0.00
2		1,4-Dioxane	40.000	40.900	-2.2	123	0.00
3		Pyridine	40.000	44.898	-12.2	130	0.00
4		n-Nitrosodimethylamine	40.000	43.825	-9.6	129	0.00
5	S	2-Fluorophenol	80.000	88.573	-10.7	130	0.00
6		Aniline	40.000	43.425	-8.6	129	0.00
7	S	Phenol-d6	80.000	91.207	-14.0	133	0.00
8		2-Chlorophenol	40.000	44.832	-12.1	133	0.00
9		Benzaldehyde	40.000	40.130	-0.3	116	0.00
10	C	Phenol	40.000	45.022	-12.6	131	0.00
11		bis(2-Chloroethyl)ether	40.000	45.308	-13.3	131	0.00
12		1,3-Dichlorobenzene	40.000	44.493	-11.2	132	0.00
13	C	1,4-Dichlorobenzene	40.000	45.483	-13.7	134	0.00
14		1,2-Dichlorobenzene	40.000	44.450	-11.1	133	0.00
15		Benzyl Alcohol	40.000	46.077	-15.2	135	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	44.227	-10.6	133	0.00
17		2-Methylphenol	40.000	46.010	-15.0	134	0.00
18		Hexachloroethane	40.000	44.609	-11.5	130	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	46.016	-15.0	135	0.00
20		3+4-Methylphenols	40.000	44.576	-11.4	130	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	122	0.00
22		Acetophenone	40.000	43.876	-9.7	133	0.00
23	S	Nitrobenzene-d5	80.000	88.074	-10.1	133	0.00
24		Nitrobenzene	40.000	44.258	-10.6	137	0.00
25		Isophorone	40.000	44.962	-12.4	134	0.00
26	C	2-Nitrophenol	40.000	45.478	-13.7	138	0.00
27		2,4-Dimethylphenol	40.000	44.339	-10.8	133	0.00
28		bis(2-Chloroethoxy)methane	40.000	44.700	-11.8	137	0.00
29	C	2,4-Dichlorophenol	40.000	45.278	-13.2	139	0.00
30		1,2,4-Trichlorobenzene	40.000	44.491	-11.2	136	0.00
31		Naphthalene	40.000	44.119	-10.3	134	0.00
32		Benzoic acid	40.000	49.387	-23.5	181	0.00
33		4-Chloroaniline	40.000	45.242	-13.1	135	0.00
34	C	Hexachlorobutadiene	40.000	45.334	-13.3	139	0.00
35		Caprolactam	40.000	42.948	-7.4	126	0.00
36	C	4-Chloro-3-methylphenol	40.000	44.784	-12.0	135	0.00
37		2-Methylnaphthalene	40.000	45.188	-13.0	136	0.00
38		1-Methylnaphthalene	40.000	44.805	-12.0	135	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	125	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	44.273	-10.7	140	0.00
41	P	Hexachlorocyclopentadiene	40.000	43.682	-9.2	136	0.00
42	S	2,4,6-Tribromophenol	80.000	88.653	-10.8	136	0.00
43	C	2,4,6-Trichlorophenol	40.000	45.119	-12.8	139	0.00
44		2,4,5-Trichlorophenol	40.000	44.243	-10.6	136	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG052720\
Data File : BG045473.D
Acq On : 27 May 2020 12:52
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 27 13:45:17 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG051520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 27 13:40:45 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	Amount	Calc.	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	80.000	88.674	-10.8	135	0.00
46		1,1'-Biphenyl	40.000	43.822	-9.6	136	0.00
47		2-Chloronaphthalene	40.000	43.853	-9.6	137	0.00
48		2-Nitroaniline	40.000	43.504	-8.8	130	0.00
49		Acenaphthylene	40.000	44.362	-10.9	137	0.00
50		Dimethylphthalate	40.000	43.924	-9.8	133	0.00
51		2,6-Dinitrotoluene	40.000	45.160	-12.9	137	0.00
52	C	Acenaphthene	40.000	44.532	-11.3	136	0.00
53		3-Nitroaniline	40.000	43.722	-9.3	135	0.00
54	P	2,4-Dinitrophenol	40.000	41.411	-3.5	129	0.00
55		Dibenzofuran	40.000	43.968	-9.9	134	0.00
56	P	4-Nitrophenol	40.000	42.472	-6.2	125	0.00
57		2,4-Dinitrotoluene	40.000	43.688	-9.2	135	0.00
58		Fluorene	40.000	44.207	-10.5	135	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	43.498	-8.7	133	0.00
60		Diethylphthalate	40.000	43.162	-7.9	131	0.00
61		4-Chlorophenyl-phenylether	40.000	44.702	-11.8	137	0.00
62		4-Nitroaniline	40.000	43.575	-8.9	131	0.00
63		Azobenzene	40.000	43.007	-7.5	131	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	118	0.00
65		4,6-Dinitro-2-methylphenol	40.000	44.072	-10.2	126	0.00
66	c	n-Nitrosodiphenylamine	40.000	44.761	-11.9	133	0.00
67		4-Bromophenyl-phenylether	40.000	45.902	-14.8	136	0.00
68		Hexachlorobenzene	40.000	45.006	-12.5	133	0.00
69		Atrazine	40.000	43.439	-8.6	126	0.00
70	C	Pentachlorophenol	40.000	42.049	-5.1	121	0.00
71		Phenanthrene	40.000	44.113	-10.3	128	0.00
72		Anthracene	40.000	44.217	-10.5	129	0.00
73		Carbazole	40.000	43.387	-8.5	125	0.00
74		Di-n-butylphthalate	40.000	43.053	-7.6	122	0.00
75	C	Fluoranthene	40.000	42.743	-6.9	121	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	107	0.00
77		Benzidine	40.000	34.320	14.2	88	0.00
78		Pyrene	40.000	45.443	-13.6	118	0.00
79	S	Terphenyl-d14	80.000	91.323	-14.2	117	0.00
80		Butylbenzylphthalate	40.000	42.227	-5.6	109	0.00
81		Benzo(a)anthracene	40.000	44.687	-11.7	118	0.00
82		3,3'-Dichlorobenzidine	40.000	43.086	-7.7	113	0.00
83		Chrysene	40.000	44.823	-12.1	117	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	43.312	-8.3	114	0.00
85	c	Di-n-octyl phthalate	40.000	43.733	-9.3	116	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	44.191	-10.5	119	0.00
87	I	Perylene-d12	20.000	20.000	0.0	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG052720\
Data File : BG045473.D
Acq On : 27 May 2020 12:52
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: May 27 13:45:17 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG051520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 27 13:40:45 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	40.000	45.270	-13.2	118	0.00
89	Benzo(k)fluoranthene	40.000	46.385	-16.0	122	0.00
90 C	Benzo(a)pyrene	40.000	45.605	-14.0	120	0.00
91	Dibenzo(a,h)anthracene	40.000	45.205	-13.0	121	0.00
92	Benzo(g,h,i)perylene	40.000	45.303	-13.3	120	0.00

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