Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062620\
Data File : BG045874.D
Acq On : 26 Jun 2020 15:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 26 16:50:37 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 15:44:32 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	123	0.00
2		1,4-Dioxane	0.542	0.483	10.9	104	-0.01
3		Pyridine	1.607	1.533	4.6	109	0.00
4		n-Nitrosodimethylamine	0.539	0.586	-8.7	126	0.00
5	S	2-Fluorophenol	1.184	1.209	-2.1	117	0.00
6		Aniline	2.121	2.200	-3.7	120	0.00
7	S	Phenol-d6	1.801	1.948	-8.2	124	0.00
8		2-Chlorophenol	1.267	1.338	-5.6	125	0.00
9		Benzaldehyde	1.082	1.132	-4.6	123	0.00
10	C	Phenol	1.745	1.869	-7.1	124	0.00
11		bis(2-Chloroethyl)ether	1.325	1.323	0.2	117	0.00
12		1,3-Dichlorobenzene	1.510	1.576	-4.4	121	0.00
13	C	1,4-Dichlorobenzene	1.509	1.561	-3.4	119	0.00
14		1,2-Dichlorobenzene	1.444	1.497	-3.7	121	0.00
15		Benzyl Alcohol	1.394	1.556	-11.6	128	0.00
16		2,2'-oxybis(1-Chloropropane	1.249	1.275	-2.1	121	0.00
17		2-Methylphenol	1.299	1.389	-6.9	123	0.00
18		Hexachloroethane	0.573	0.614	-7.2	124	0.00
19	P	n-Nitroso-di-n-propylamine	1.184	1.281	-8.2	125	0.00
20		3+4-Methylphenols	1.708	1.932	-13.1	129	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	129	0.00
22		Acetophenone	0.549	0.560	-2.0	121	0.00
23	S	Nitrobenzene-d5	0.438	0.447	-2.1	122	0.00
24		Nitrobenzene	0.467	0.477	-2.1	121	0.00
25		Isophorone	0.848	0.892	-5.2	124	0.00
26	C	2-Nitrophenol	0.194	0.209	-7.7	124	0.00
27		2,4-Dimethylphenol	0.297	0.322	-8.4	131	0.00
28		bis(2-Chloroethoxy)methane	0.443	0.465	-5.0	126	0.00
29	C	2,4-Dichlorophenol	0.346	0.388	-12.1	133	0.00
30		1,2,4-Trichlorobenzene	0.410	0.431	-5.1	127	0.00
31		Naphthalene	1.093	1.123	-2.7	122	0.00
32		Benzoic acid	0.170	0.247	-45.3#	179#	0.00
33		4-Chloroaniline	0.458	0.494	-7.9	128	0.00
34	C	Hexachlorobutadiene	0.292	0.310	-6.2	126	0.00
35		Caprolactam	0.112	0.133	-18.8	143	0.00
36	C	4-Chloro-3-methylphenol	0.400	0.466	-16.5	140	0.00
37		2-Methylnaphthalene	0.814	0.887	-9.0	130	0.00
38		1-Methylnaphthalene	0.771	0.837	-8.6	129	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	144	0.00
40		1,2,4,5-Tetrachlorobenzene	0.659	0.667	-1.2	133	0.00
41	P	Hexachlorocyclopentadiene	0.299	0.149	50.2#	64	0.00
42	S	2,4,6-Tribromophenol	0.302	0.320	-6.0	143	0.00
43	C	2,4,6-Trichlorophenol	0.446	0.465	-4.3	139	0.00
44		2,4,5-Trichlorophenol	0.509	0.528	-3.7	137	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062620\
Data File : BG045874.D
Acq On : 26 Jun 2020 15:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 26 16:50:37 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 15:44:32 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.361	1.349	0.9	131	0.00
46		1,1'-Biphenyl	1.432	1.419	0.9	132	0.00
47		2-Chloronaphthalene	1.157	1.156	0.1	133	0.00
48		2-Nitroaniline	0.385	0.413	-7.3	141	0.00
49		Acenaphthylene	1.853	1.860	-0.4	133	0.00
50		Dimethylphthalate	1.567	1.567	0.0	134	0.00
51		2,6-Dinitrotoluene	0.352	0.359	-2.0	136	0.00
52	C	Acenaphthene	1.123	1.142	-1.7	137	0.00
53		3-Nitroaniline	0.351	0.374	-6.6	143	0.00
54	P	2,4-Dinitrophenol	0.152	0.190	-25.0#	172#	0.00
55		Dibenzofuran	1.824	1.871	-2.6	136	0.00
56	P	4-Nitrophenol	0.226	0.244	-8.0	143	0.00
57		2,4-Dinitrotoluene	0.504	0.521	-3.4	137	0.00
58		Fluorene	1.517	1.580	-4.2	138	0.00
59		2,3,4,6-Tetrachlorophenol	0.442	0.465	-5.2	140	0.00
60		Diethylphthalate	1.602	1.659	-3.6	137	0.00
61		4-Chlorophenyl-phenylether	0.874	0.910	-4.1	139	0.00
62		4-Nitroaniline	0.361	0.385	-6.6	140	0.00
63		Azobenzene	1.538	1.576	-2.5	138	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	144	0.00
65		4,6-Dinitro-2-methylphenol	0.124	0.132	-6.5	143	0.00
66	c	n-Nitrosodiphenylamine	0.566	0.584	-3.2	139	0.00
67		4-Bromophenyl-phenylether	0.260	0.269	-3.5	138	0.00
68		Hexachlorobenzene	0.269	0.285	-5.9	145	0.00
69		Atrazine	0.220	0.206	6.4	127	0.00
70	C	Pentachlorophenol	0.111	0.116	-4.5	136	0.00
71		Phenanthrene	1.031	1.045	-1.4	136	0.00
72		Anthracene	1.027	1.043	-1.6	136	0.00
73		Carbazole	0.978	1.000	-2.2	136	0.00
74		Di-n-butylphthalate	1.158	1.177	-1.6	135	0.00
75	C	Fluoranthene	1.268	1.249	1.5	131	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	139	0.00
77		Benzidine	0.520	0.554	-6.5	129	0.00
78		Pyrene	1.339	1.365	-1.9	129	0.00
79	S	Terphenyl-d14	0.987	1.000	-1.3	127	0.00
80		Butylbenzylphthalate	0.561	0.572	-2.0	131	0.00
81		Benzo(a)anthracene	1.297	1.341	-3.4	134	0.00
82		3,3'-Dichlorobenzidine	0.490	0.504	-2.9	131	0.00
83		Chrysene	1.254	1.266	-1.0	128	0.00
84		Bis(2-ethylhexyl)phthalate	0.780	0.794	-1.8	131	0.00
85	c	Di-n-octyl phthalate	1.345	1.349	-0.3	129	0.00
86	I	Perylene-d12	1.000	1.000	0.0	134	0.00
87		Indeno(1,2,3-cd)pyrene	1.450	1.533	-5.7	132	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG062620\
Data File : BG045874.D
Acq On : 26 Jun 2020 15:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 26 16:50:37 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 15:44:32 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.254	1.284	-2.4	127	0.00
89	Benzo(k)fluoranthene	1.200	1.245	-3.8	127	0.00
90 C	Benzo(a)pyrene	1.171	1.219	-4.1	130	0.00
91	Dibenzo(a,h)anthracene	1.163	1.239	-6.5	133	0.00
92	Benzo(g,h,i)perylene	1.164	1.239	-6.4	133	0.00

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