Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040219\
Data File : BG040318.D
Acq On : 3 Apr 2019 7:27
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 03 08:39:33 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 28 05:47:00 2019
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	106	-0.01
2		1,4-Dioxane	0.566	0.557	1.6	104	-0.03
3		Pyridine	1.523	1.628	-6.9	106	-0.02
4		n-Nitrosodimethylamine	0.705	0.686	2.7	102	-0.01
5	S	2-Fluorophenol	1.203	1.245	-3.5	107	-0.01
6		Aniline	2.160	2.258	-4.5	107	-0.02
7	S	Phenol-d6	1.663	1.810	-8.8	112	-0.01
8		2-Chlorophenol	1.408	1.448	-2.8	106	-0.01
9		Benzaldehyde	1.140	1.161	-1.8	101	-0.01
10	C	Phenol	1.805	1.976	-9.5	113	-0.02
11		bis(2-Chloroethyl)ether	1.445	1.468	-1.6	104	-0.01
12		1,3-Dichlorobenzene	1.531	1.505	1.7	101	-0.02
13	C	1,4-Dichlorobenzene	1.525	1.547	-1.4	105	-0.01
14		1,2-Dichlorobenzene	1.458	1.477	-1.3	105	-0.01
15		Benzyl Alcohol	1.339	1.427	-6.6	109	-0.02
16		2,2'-oxybis(1-Chloropropane	2.774	2.815	-1.5	104	-0.01
17		2-Methylphenol	1.249	1.323	-5.9	108	-0.01
18		Hexachloroethane	0.580	0.567	2.2	102	-0.02
19	P	n-Nitroso-di-n-propylamine	1.192	1.283	-7.6	112	-0.02
20		3+4-Methylphenols	1.692	1.881	-11.2	113	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	111	-0.01
22		Acetophenone	0.499	0.506	-1.4	113	-0.01
23	S	Nitrobenzene-d5	0.376	0.376	0.0	110	-0.01
24		Nitrobenzene	0.390	0.391	-0.3	111	-0.02
25		Isophorone	0.774	0.812	-4.9	116	-0.02
26	C	2-Nitrophenol	0.185	0.199	-7.6	118	-0.01
27		2,4-Dimethylphenol	0.293	0.309	-5.5	116	-0.02
28		bis(2-Chloroethoxy)methane	0.458	0.478	-4.4	113	-0.02
29	C	2,4-Dichlorophenol	0.318	0.349	-9.7	119	-0.02
30		1,2,4-Trichlorobenzene	0.350	0.349	0.3	111	-0.01
31		Naphthalene	1.025	1.063	-3.7	113	-0.02
32		Benzoic acid	0.177	0.027	84.7#	18#	0.03
33		4-Chloroaniline	0.437	0.464	-6.2	115	-0.01
34	C	Hexachlorobutadiene	0.224	0.214	4.5	106	-0.01
35		Caprolactam	0.126	0.145	-15.1	125	-0.02
36	C	4-Chloro-3-methylphenol	0.366	0.412	-12.6	123	-0.02
37		2-Methylnaphthalene	0.758	0.805	-6.2	117	-0.02
38		1-Methylnaphthalene	0.735	0.782	-6.4	117	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	124	-0.02
40		1,2,4,5-Tetrachlorobenzene	0.636	0.631	0.8	120	-0.01
41	P	Hexachlorocyclopentadiene	0.349	0.314	10.0	113	-0.01
42	S	2,4,6-Tribromophenol	0.216	0.240	-11.1	135	-0.01
43	C	2,4,6-Trichlorophenol	0.410	0.427	-4.1	127	-0.01
44		2,4,5-Trichlorophenol	0.447	0.478	-6.9	132	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040219\
Data File : BG040318.D
Acq On : 3 Apr 2019 7:27
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 03 08:39:33 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 28 05:47:00 2019
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	1.350	1.338	0.9	119	-0.01
46		1,1'-Biphenyl	1.580	1.583	-0.2	122	-0.01
47		2-Chloronaphthalene	1.212	1.216	-0.3	122	-0.02
48		2-Nitroaniline	0.409	0.432	-5.6	127	-0.02
49		Acenaphthylene	2.055	2.121	-3.2	124	-0.01
50		Dimethylphthalate	1.565	1.629	-4.1	126	-0.01
51		2,6-Dinitrotoluene	0.351	0.376	-7.1	131	-0.02
52	C	Acenaphthene	1.178	1.201	-2.0	124	-0.02
53		3-Nitroaniline	0.372	0.389	-4.6	127	-0.01
54	P	2,4-Dinitrophenol	0.187	0.040#	78.6#	27#	-0.01
55		Dibenzofuran	1.892	1.993	-5.3	128	-0.01
56	P	4-Nitrophenol	0.300	0.261	13.0	107	-0.01
57		2,4-Dinitrotoluene	0.488	0.526	-7.8	130	-0.01
58		Fluorene	1.488	1.558	-4.7	128	-0.02
59		2,3,4,6-Tetrachlorophenol	0.405	0.440	-8.6	131	-0.01
60		Diethylphthalate	1.602	1.655	-3.3	126	-0.01
61		4-Chlorophenyl-phenylether	0.795	0.835	-5.0	127	-0.01
62		4-Nitroaniline	0.399	0.421	-5.5	129	-0.01
63		Azobenzene	1.511	1.584	-4.8	128	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	131	-0.01
65		4,6-Dinitro-2-methylphenol	0.112	0.047	58.0#	56	-0.01
66	c	n-Nitrosodiphenylamine	0.585	0.597	-2.1	129	-0.02
67		4-Bromophenyl-phenylether	0.225	0.235	-4.4	133	-0.02
68		Hexachlorobenzene	0.226	0.238	-5.3	135	-0.01
69		Atrazine	0.218	0.220	-0.9	128	-0.02
70	C	Pentachlorophenol	0.131	0.129	1.5	129	-0.01
71		Phenanthrene	1.051	1.081	-2.9	130	-0.02
72		Anthracene	1.052	1.068	-1.5	128	-0.02
73		Carbazole	0.996	1.008	-1.2	128	-0.02
74		Di-n-butylphthalate	1.180	1.171	0.8	125	-0.02
75	C	Fluoranthene	1.245	1.275	-2.4	129	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	132	-0.01
77		Benzidine	0.722	0.840	-16.3	148	0.00
78		Pyrene	1.315	1.302	1.0	127	-0.02
79	S	Terphenyl-d14	0.889	0.854	3.9	124	-0.02
80		Butylbenzylphthalate	0.563	0.550	2.3	126	-0.01
81		Benzo(a)anthracene	1.232	1.234	-0.2	129	-0.01
82		3,3'-Dichlorobenzidine	0.469	0.456	2.8	123	-0.01
83		Chrysene	1.184	1.218	-2.9	132	-0.01
84		Bis(2-ethylhexyl)phthalate	0.805	0.785	2.5	127	-0.01
85	c	Di-n-octyl phthalate	1.371	1.319	3.8	123	0.00
86		Indeno(1,2,3-cd)pyrene	1.448	1.465	-1.2	132	0.06
87	I	Perylene-d12	1.000	1.000	0.0	123	0.03

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040219\
Data File : BG040318.D
Acq On : 3 Apr 2019 7:27
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Apr 03 08:39:33 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 28 05:47:00 2019
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	1.224	1.313	-7.3	129	0.00
89	Benzo(k)fluoranthene	1.126	1.170	-3.9	127	0.00
90 C	Benzo(a)pyrene	1.149	1.217	-5.9	128	0.03
91	Dibenzo(a,h)anthracene	1.067	1.180	-10.6	134	0.13
92	Benzo(g,h,i)perylene	1.097	1.197	-9.1	133	0.20

(#) = Out of Range SPCC's out = 1 CCC's out = 0