Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090119\
Data File : BG042666.D
Acq On : 1 Sep 2019 17:58
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 02 07:55:17 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 22 14:29:15 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	67	-0.01
2		1,4-Dioxane	0.412	0.357	13.3	58	0.00
3		Pyridine	1.057	0.923	12.7	57	0.00
4		n-Nitrosodimethylamine	0.377	0.359	4.8	65	0.00
5	S	2-Fluorophenol	0.988	0.926	6.3	63	-0.01
6		Aniline	1.779	1.606	9.7	61	0.00
7	S	Phenol-d6	1.334	1.247	6.5	62	-0.01
8		2-Chlorophenol	1.197	1.143	4.5	64	-0.01
9		Benzaldehyde	0.668	0.617	7.6	74	-0.01
10	C	Phenol	1.356	1.269	6.4	62	0.00
11		bis(2-Chloroethyl)ether	1.065	0.962	9.7	60	0.00
12		1,3-Dichlorobenzene	1.522	1.466	3.7	65	-0.01
13	C	1,4-Dichlorobenzene	1.542	1.497	2.9	65	-0.01
14		1,2-Dichlorobenzene	1.477	1.452	1.7	67	-0.01
15		Benzyl Alcohol	0.960	0.956	0.4	67	-0.01
16		2,2'-oxybis(1-Chloropropane	1.444	1.260	12.7	59	-0.02
17		2-Methylphenol	0.999	0.954	4.5	65	-0.01
18		Hexachloroethane	0.528	0.499	5.5	64	0.00
19	P	n-Nitroso-di-n-propylamine	0.864	0.808	6.5	63	-0.01
20		3+4-Methylphenols	1.382	1.338	3.2	66	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	69	-0.01
22		Acetophenone	0.428	0.420	1.9	67	0.00
23	S	Nitrobenzene-d5	0.289	0.286	1.0	66	-0.01
24		Nitrobenzene	0.293	0.277	5.5	64	-0.01
25		Isophorone	0.571	0.550	3.7	64	-0.01
26	C	2-Nitrophenol	0.184	0.180	2.2	67	0.00
27		2,4-Dimethylphenol	0.253	0.255	-0.8	67	-0.01
28		bis(2-Chloroethoxy)methane	0.360	0.342	5.0	64	-0.01
29	C	2,4-Dichlorophenol	0.331	0.348	-5.1	70	-0.01
30		1,2,4-Trichlorobenzene	0.406	0.428	-5.4	72	-0.01
31		Naphthalene	0.929	0.925	0.4	67	-0.01
32		Benzoic acid	0.169	0.138	18.3	58	0.00
33		4-Chloroaniline	0.429	0.433	-0.9	67	0.00
34	C	Hexachlorobutadiene	0.273	0.297	-8.8	74	-0.01
35		Caprolactam	0.110	0.110	0.0	64	0.00
36	C	4-Chloro-3-methylphenol	0.314	0.327	-4.1	69	0.00
37		2-Methylnaphthalene	0.746	0.777	-4.2	69	-0.01
38		1-Methylnaphthalene	0.708	0.739	-4.4	69	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	74	0.00
40		1,2,4,5-Tetrachlorobenzene	0.642	0.652	-1.6	74	0.00
41	P	Hexachlorocyclopentadiene	0.337	0.320	5.0	72	-0.01
42	S	2,4,6-Tribromophenol	0.284	0.302	-6.3	75	-0.01
43	C	2,4,6-Trichlorophenol	0.434	0.429	1.2	72	-0.01
44		2,4,5-Trichlorophenol	0.450	0.438	2.7	72	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090119\
Data File : BG042666.D
Acq On : 1 Sep 2019 17:58
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 02 07:55:17 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 22 14:29:15 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	1.183	1.222	-3.3	74	0.00
46		1,1'-Biphenyl	1.293	1.271	1.7	72	0.00
47		2-Chloronaphthalene	1.115	1.100	1.3	72	-0.01
48		2-Nitroaniline	0.269	0.255	5.2	69	0.00
49		Acenaphthylene	1.677	1.671	0.4	71	0.00
50		Dimethylphthalate	1.443	1.473	-2.1	72	-0.01
51		2,6-Dinitrotoluene	0.334	0.341	-2.1	73	0.00
52	C	Acenaphthene	1.018	0.999	1.9	70	0.00
53		3-Nitroaniline	0.335	0.329	1.8	70	0.00
54	P	2,4-Dinitrophenol	0.198	0.187	5.6	68	0.00
55		Dibenzofuran	1.686	1.739	-3.1	73	0.00
56	P	4-Nitrophenol	0.238	0.216	9.2	65	0.00
57		2,4-Dinitrotoluene	0.479	0.499	-4.2	73	0.00
58		Fluorene	1.378	1.426	-3.5	73	0.00
59		2,3,4,6-Tetrachlorophenol	0.445	0.466	-4.7	75	-0.01
60		Diethylphthalate	1.411	1.452	-2.9	72	-0.01
61		4-Chlorophenyl-phenylether	0.831	0.867	-4.3	74	0.00
62		4-Nitroaniline	0.358	0.362	-1.1	71	0.00
63		Azobenzene	0.967	0.911	5.8	67	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	76	-0.01
65		4,6-Dinitro-2-methylphenol	0.125	0.114	8.8	66	0.00
66	c	n-Nitrosodiphenylamine	0.550	0.547	0.5	73	-0.01
67		4-Bromophenyl-phenylether	0.254	0.260	-2.4	76	-0.01
68		Hexachlorobenzene	0.276	0.281	-1.8	75	-0.01
69		Atrazine	0.195	0.212	-8.7	83	0.00
70	C	Pentachlorophenol	0.143	0.122	14.7	62	0.00
71		Phenanthrene	0.993	1.020	-2.7	73	0.00
72		Anthracene	0.992	1.015	-2.3	73	-0.01
73		Carbazole	0.884	0.891	-0.8	72	0.00
74		Di-n-butylphthalate	1.045	1.057	-1.1	71	-0.01
75	C	Fluoranthene	1.212	1.309	-8.0	76	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	79	0.00
77		Benzidine	0.573	0.785	-37.0#	123	0.00
78		Pyrene	1.226	1.231	-0.4	75	0.00
79	S	Terphenyl-d14	0.925	0.948	-2.5	78	0.00
80		Butylbenzylphthalate	0.482	0.465	3.5	73	-0.01
81		Benzo(a)anthracene	1.217	1.260	-3.5	78	-0.01
82		3,3'-Dichlorobenzidine	0.489	0.501	-2.5	79	-0.01
83		Chrysene	1.173	1.209	-3.1	77	-0.01
84		Bis(2-ethylhexyl)phthalate	0.670	0.653	2.5	74	-0.01
85	c	Di-n-octyl phthalate	1.152	1.127	2.2	74	-0.02
86		Indeno(1,2,3-cd)pyrene	1.595	1.583	0.8	76	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	78	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090119\
Data File : BG042666.D
Acq On : 1 Sep 2019 17:58
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 02 07:55:17 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 22 14:29:15 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	1.100	1.122	-2.0	75	-0.01
89	Benzo(k)fluoranthene	1.057	1.108	-4.8	79	-0.01
90 C	Benzo(a)pyrene	1.043	1.074	-3.0	77	-0.01
91	Dibenzo(a,h)anthracene	1.056	1.069	-1.2	77	-0.02
92	Benzo(g,h,i)perylene	1.057	1.042	1.4	75	-0.02

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