Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070720\
Data File : BG045953.D
Acq On : 7 Jul 2020 9:04
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 07 11:22:44 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 07 09:58:42 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	94	0.00
2		1,4-Dioxane	0.569	0.575	-1.1	97	0.00
3		Pyridine	1.720	1.769	-2.8	98	0.00
4		n-Nitrosodimethylamine	0.582	0.607	-4.3	95	0.00
5	S	2-Fluorophenol	1.233	1.223	0.8	97	0.00
6		Aniline	2.271	2.287	-0.7	96	0.00
7	S	Phenol-d6	1.776	1.791	-0.8	97	0.00
8		2-Chlorophenol	1.328	1.316	0.9	96	0.00
9		Benzaldehyde	1.125	1.075	4.4	104	0.00
10	C	Phenol	1.781	1.795	-0.8	98	0.00
11		bis(2-Chloroethyl)ether	1.473	1.503	-2.0	99	0.00
12		1,3-Dichlorobenzene	1.530	1.527	0.2	97	0.00
13	C	1,4-Dichlorobenzene	1.505	1.524	-1.3	97	0.00
14		1,2-Dichlorobenzene	1.449	1.444	0.3	96	0.00
15		Benzyl Alcohol	1.392	1.398	-0.4	96	0.00
16		2,2'-oxybis(1-Chloropropane	1.922	1.978	-2.9	98	0.00
17		2-Methylphenol	1.336	1.321	1.1	94	0.00
18		Hexachloroethane	0.579	0.578	0.2	96	0.00
19	P	n-Nitroso-di-n-propylamine	1.247	1.276	-2.3	98	0.00
20		3+4-Methylphenols	1.821	1.818	0.2	94	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22		Acetophenone	0.534	0.530	0.7	95	0.00
23	S	Nitrobenzene-d5	0.347	0.374	-7.8	102	0.00
24		Nitrobenzene	0.384	0.419	-9.1	101	0.00
25		Isophorone	0.848	0.866	-2.1	96	0.00
26	C	2-Nitrophenol	0.125	0.138	-10.4	107	0.00
27		2,4-Dimethylphenol	0.302	0.305	-1.0	95	0.00
28		bis(2-Chloroethoxy)methane	0.476	0.488	-2.5	96	0.00
29	C	2,4-Dichlorophenol	0.313	0.312	0.3	94	0.00
30		1,2,4-Trichlorobenzene	0.346	0.349	-0.9	96	0.00
31		Naphthalene	1.072	1.072	0.0	96	0.00
32		Benzoic acid	0.163	0.183	-12.3	102	0.00
33		4-Chloroaniline	0.479	0.474	1.0	94	0.00
34	C	Hexachlorobutadiene	0.202	0.208	-3.0	99	0.00
35		Caprolactam	0.119	0.116	2.5	91	0.00
36	C	4-Chloro-3-methylphenol	0.356	0.363	-2.0	95	0.00
37		2-Methylnaphthalene	0.762	0.769	-0.9	95	0.00
38		1-Methylnaphthalene	0.719	0.718	0.1	94	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
40		1,2,4,5-Tetrachlorobenzene	0.564	0.569	-0.9	96	0.00
41	P	Hexachlorocyclopentadiene	0.335	0.335	0.0	94	0.00
42	S	2,4,6-Tribromophenol	0.186	0.196	-5.4	95	0.00
43	C	2,4,6-Trichlorophenol	0.381	0.393	-3.1	96	0.00
44		2,4,5-Trichlorophenol	0.431	0.444	-3.0	95	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070720\
Data File : BG045953.D
Acq On : 7 Jul 2020 9:04
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 07 11:22:44 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 07 09:58:42 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	1.282	1.293	-0.9	96	0.00
46		1,1'-Biphenyl	1.505	1.526	-1.4	96	0.00
47		2-Chloronaphthalene	1.187	1.179	0.7	94	0.00
48		2-Nitroaniline	0.278	0.357	-28.4#	104	0.00
49		Acenaphthylene	1.915	1.901	0.7	93	0.00
50		Dimethylphthalate	1.501	1.492	0.6	94	0.00
51		2,6-Dinitrotoluene	0.233	0.278	-19.3	102	0.00
52	C	Acenaphthene	1.208	1.217	-0.7	94	0.00
53		3-Nitroaniline	0.289	0.338	-17.0	98	0.00
54	P	2,4-Dinitrophenol	0.082	0.094	-14.6	99	0.00
55		Dibenzofuran	1.771	1.742	1.6	93	0.00
56	P	4-Nitrophenol	0.245	0.269	-9.8	101	0.00
57		2,4-Dinitrotoluene	0.288	0.363	-26.0#	103	0.00
58		Fluorene	1.453	1.420	2.3	92	0.00
59		2,3,4,6-Tetrachlorophenol	0.333	0.348	-4.5	95	0.00
60		Diethylphthalate	1.574	1.542	2.0	92	0.00
61		4-Chlorophenyl-phenylether	0.723	0.721	0.3	95	0.00
62		4-Nitroaniline	0.310	0.344	-11.0	99	0.00
63		Azobenzene	1.689	1.680	0.5	94	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
65		4,6-Dinitro-2-methylphenol	0.059	0.071	-20.3	108	0.00
66	c	n-Nitrosodiphenylamine	0.620	0.625	-0.8	93	0.00
67		4-Bromophenyl-phenylether	0.223	0.223	0.0	90	0.00
68		Hexachlorobenzene	0.224	0.223	0.4	92	0.00
69		Atrazine	0.183	0.183	0.0	90	0.00
70	C	Pentachlorophenol	0.125	0.134	-7.2	93	0.00
71		Phenanthrene	1.057	1.065	-0.8	93	0.00
72		Anthracene	1.073	1.059	1.3	91	0.00
73		Carbazole	1.076	1.061	1.4	91	0.00
74		Di-n-butylphthalate	1.300	1.265	2.7	91	0.00
75	C	Fluoranthene	1.222	1.202	1.6	92	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
77		Benzidine	0.572	0.641	-12.1	99	0.00
78		Pyrene	1.365	1.417	-3.8	92	0.00
79	S	Terphenyl-d14	0.899	0.921	-2.4	93	0.00
80		Butylbenzylphthalate	0.645	0.680	-5.4	92	0.00
81		Benzo(a)anthracene	1.315	1.322	-0.5	91	0.00
82		3,3'-Dichlorobenzidine	0.486	0.500	-2.9	92	0.00
83		Chrysene	1.253	1.259	-0.5	90	0.00
84		Bis(2-ethylhexyl)phthalate	0.936	0.953	-1.8	92	0.00
85	c	Di-n-octyl phthalate	1.625	1.667	-2.6	92	0.00
86	I	Perylene-d12	1.000	1.000	0.0	91	0.00
87		Indeno(1,2,3-cd)pyrene	1.463	1.506	-2.9	92	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070720\
Data File : BG045953.D
Acq On : 7 Jul 2020 9:04
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 07 11:22:44 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG070620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 07 09:58:42 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	1.270	1.319	-3.9	93	0.00
89	Benzo(k)fluoranthene	1.200	1.216	-1.3	90	0.00
90 C	Benzo(a)pyrene	1.194	1.229	-2.9	93	0.00
91	Dibenzo(a,h)anthracene	1.180	1.186	-0.5	91	-0.01
92	Benzo(g,h,i)perylene	1.188	1.202	-1.2	91	-0.01

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