Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091820\
Data File : BG046669.D
Acq On : 18 Sep 2020 13:44
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 18 14:17:37 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 15:06:53 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	68	0.00
2		1,4-Dioxane	0.470	0.484	-3.0	77	0.00
3		Pyridine	1.375	1.458	-6.0	76	0.00
4		n-Nitrosodimethylamine	0.507	0.531	-4.7	73	0.00
5	S	2-Fluorophenol	1.129	1.165	-3.2	76	0.00
6		Aniline	1.801	1.950	-8.3	78	0.00
7	S	Phenol-d6	1.601	1.740	-8.7	78	0.00
8		2-Chlorophenol	1.193	1.236	-3.6	77	0.00
9		Benzaldehyde	0.897	0.936	-4.3	76	0.00
10	C	Phenol	1.474	1.590	-7.9	78	0.00
11		bis(2-Chloroethyl)ether	1.185	1.265	-6.8	78	0.00
12		1,3-Dichlorobenzene	1.515	1.538	-1.5	75	0.00
13	C	1,4-Dichlorobenzene	1.517	1.566	-3.2	76	0.00
14		1,2-Dichlorobenzene	1.455	1.483	-1.9	76	0.00
15		Benzyl Alcohol	1.028	1.143	-11.2	78	0.00
16		2,2'-oxybis(1-Chloropropane	1.838	1.959	-6.6	77	0.00
17		2-Methylphenol	1.046	1.138	-8.8	79	0.00
18		Hexachloroethane	0.514	0.520	-1.2	75	0.00
19	P	n-Nitroso-di-n-propylamine	0.958	1.083	-13.0	81	0.00
20		3+4-Methylphenols	1.443	1.599	-10.8	79	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	71	0.00
22		Acetophenone	0.487	0.497	-2.1	79	0.00
23	S	Nitrobenzene-d5	0.350	0.358	-2.3	79	0.00
24		Nitrobenzene	0.346	0.351	-1.4	79	0.00
25		Isophorone	0.642	0.687	-7.0	81	0.00
26	C	2-Nitrophenol	0.182	0.194	-6.6	79	0.00
27		2,4-Dimethylphenol	0.251	0.263	-4.8	80	0.00
28		bis(2-Chloroethoxy)methane	0.413	0.434	-5.1	80	0.00
29	C	2,4-Dichlorophenol	0.334	0.352	-5.4	80	0.00
30		1,2,4-Trichlorobenzene	0.397	0.392	1.3	78	0.00
31		Naphthalene	0.977	1.004	-2.8	80	0.00
32		Benzoic acid	0.156	0.096	38.5#	42#	0.00
33		4-Chloroaniline	0.431	0.462	-7.2	81	0.00
34	C	Hexachlorobutadiene	0.258	0.253	1.9	76	0.00
35		Caprolactam	0.125	0.139	-11.2	83	0.00
36	C	4-Chloro-3-methylphenol	0.327	0.362	-10.7	83	0.00
37		2-Methylnaphthalene	0.770	0.812	-5.5	81	0.00
38		1-Methylnaphthalene	0.730	0.773	-5.9	81	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	73	0.00
40		1,2,4,5-Tetrachlorobenzene	0.660	0.655	0.8	81	0.00
41	P	Hexachlorocyclopentadiene	0.287	0.291	-1.4	76	0.00
42	S	2,4,6-Tribromophenol	0.271	0.275	-1.5	80	0.00
43	C	2,4,6-Trichlorophenol	0.445	0.442	0.7	78	0.00
44		2,4,5-Trichlorophenol	0.468	0.467	0.2	79	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091820\
Data File : BG046669.D
Acq On : 18 Sep 2020 13:44
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 18 14:17:37 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 15:06:53 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	1.310	1.331	-1.6	83	0.00
46		1,1'-Biphenyl	1.391	1.430	-2.8	83	0.00
47		2-Chloronaphthalene	1.180	1.194	-1.2	82	0.00
48		2-Nitroaniline	0.328	0.352	-7.3	85	0.00
49		Acenaphthylene	1.790	1.862	-4.0	84	0.00
50		Dimethylphthalate	1.547	1.605	-3.7	83	0.00
51		2,6-Dinitrotoluene	0.341	0.353	-3.5	83	0.00
52	C	Acenaphthene	1.109	1.136	-2.4	83	0.00
53		3-Nitroaniline	0.369	0.384	-4.1	81	0.00
54	P	2,4-Dinitrophenol	0.199	0.186	6.5	67	0.00
55		Dibenzofuran	1.780	1.851	-4.0	85	0.00
56	P	4-Nitrophenol	0.272	0.265	2.6	74	0.00
57		2,4-Dinitrotoluene	0.486	0.508	-4.5	81	0.00
58		Fluorene	1.494	1.528	-2.3	83	0.00
59		2,3,4,6-Tetrachlorophenol	0.454	0.471	-3.7	82	0.00
60		Diethylphthalate	1.509	1.547	-2.5	83	0.00
61		4-Chlorophenyl-phenylether	0.823	0.836	-1.6	82	0.00
62		4-Nitroaniline	0.390	0.395	-1.3	80	0.00
63		Azobenzene	1.227	1.279	-4.2	85	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	73	0.00
65		4,6-Dinitro-2-methylphenol	0.113	0.122	-8.0	79	0.00
66	c	n-Nitrosodiphenylamine	0.540	0.569	-5.4	84	0.00
67		4-Bromophenyl-phenylether	0.233	0.247	-6.0	83	0.00
68		Hexachlorobenzene	0.258	0.267	-3.5	82	0.00
69		Atrazine	0.220	0.225	-2.3	79	0.00
70	C	Pentachlorophenol	0.135	0.125	7.4	67	0.00
71		Phenanthrene	1.014	1.044	-3.0	83	0.00
72		Anthracene	1.000	1.023	-2.3	82	0.00
73		Carbazole	0.944	0.951	-0.7	80	0.00
74		Di-n-butylphthalate	1.086	1.082	0.4	80	0.00
75	C	Fluoranthene	1.305	1.277	2.1	80	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	72	0.00
77		Benzidine	0.465	0.454	2.4	66	0.00
78		Pyrene	1.274	1.309	-2.7	79	0.00
79	S	Terphenyl-d14	0.940	0.935	0.5	79	0.00
80		Butylbenzylphthalate	0.497	0.505	-1.6	76	0.00
81		Benzo(a)anthracene	1.247	1.258	-0.9	79	0.00
82		3,3'-Dichlorobenzidine	0.463	0.468	-1.1	77	0.00
83		Chrysene	1.199	1.217	-1.5	80	0.00
84		Bis(2-ethylhexyl)phthalate	0.678	0.695	-2.5	78	0.00
85	c	Di-n-octyl phthalate	1.161	1.210	-4.2	80	0.00
86	I	Perylene-d12	1.000	1.000	0.0	68	0.00
87		Indeno(1,2,3-cd)pyrene	1.436	1.484	-3.3	79	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091820\
Data File : BG046669.D
Acq On : 18 Sep 2020 13:44
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 18 14:17:37 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 15:06:53 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	1.249	1.269	-1.6	78	0.00
89	Benzo(k)fluoranthene	1.207	1.262	-4.6	80	0.00
90 C	Benzo(a)pyrene	1.165	1.197	-2.7	78	0.00
91	Dibenzo(a,h)anthracene	1.159	1.200	-3.5	79	0.00
92	Benzo(g,h,i)perylene	1.158	1.196	-3.3	79	0.00

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