Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG091720\
Data File : BG046658.D
Acq On : 17 Sep 2020 11:12
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 17 14:36:23 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 17 14:27:03 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	64	0.00
2		1,4-Dioxane	0.470	0.459	2.3	68	0.00
3		Pyridine	1.375	1.384	-0.7	67	0.00
4		n-Nitrosodimethylamine	0.507	0.511	-0.8	65	0.00
5	S	2-Fluorophenol	1.129	1.110	1.7	67	0.00
6		Aniline	1.801	1.860	-3.3	69	0.00
7	S	Phenol-d6	1.601	1.663	-3.9	69	0.00
8		2-Chlorophenol	1.193	1.203	-0.8	69	0.00
9		Benzaldehyde	0.897	0.857	4.5	65	0.00
10	C	Phenol	1.474	1.510	-2.4	69	0.00
11		bis(2-Chloroethyl)ether	1.185	1.188	-0.3	68	0.00
12		1,3-Dichlorobenzene	1.515	1.518	-0.2	69	0.00
13	C	1,4-Dichlorobenzene	1.517	1.536	-1.3	70	0.00
14		1,2-Dichlorobenzene	1.455	1.470	-1.0	70	0.00
15		Benzyl Alcohol	1.028	1.114	-8.4	71	0.00
16		2,2'-oxybis(1-Chloropropane	1.838	1.900	-3.4	70	0.00
17		2-Methylphenol	1.046	1.091	-4.3	70	0.00
18		Hexachloroethane	0.514	0.515	-0.2	69	0.00
19	P	n-Nitroso-di-n-propylamine	0.958	1.044	-9.0	72	0.00
20		3+4-Methylphenols	1.443	1.530	-6.0	70	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	64	0.00
22		Acetophenone	0.487	0.499	-2.5	72	0.00
23	S	Nitrobenzene-d5	0.350	0.354	-1.1	71	0.00
24		Nitrobenzene	0.346	0.346	0.0	70	0.00
25		Isophorone	0.642	0.680	-5.9	72	0.00
26	C	2-Nitrophenol	0.182	0.186	-2.2	69	0.00
27		2,4-Dimethylphenol	0.251	0.260	-3.6	71	0.00
28		bis(2-Chloroethoxy)methane	0.413	0.433	-4.8	72	0.00
29	C	2,4-Dichlorophenol	0.334	0.344	-3.0	70	0.00
30		1,2,4-Trichlorobenzene	0.397	0.396	0.3	71	0.00
31		Naphthalene	0.977	1.006	-3.0	72	0.00
32		Benzoic acid	0.156	0.095	39.1#	38#	0.00
33		4-Chloroaniline	0.431	0.457	-6.0	72	0.00
34	C	Hexachlorobutadiene	0.258	0.261	-1.2	71	0.00
35		Caprolactam	0.125	0.140	-12.0	75	0.00
36	C	4-Chloro-3-methylphenol	0.327	0.354	-8.3	73	0.00
37		2-Methylnaphthalene	0.770	0.813	-5.6	73	0.00
38		1-Methylnaphthalene	0.730	0.761	-4.2	72	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	67	0.00
40		1,2,4,5-Tetrachlorobenzene	0.660	0.649	1.7	72	0.00
41	P	Hexachlorocyclopentadiene	0.287	0.291	-1.4	69	0.00
42	S	2,4,6-Tribromophenol	0.271	0.287	-5.9	76	0.00
43	C	2,4,6-Trichlorophenol	0.445	0.437	1.8	70	0.00
44		2,4,5-Trichlorophenol	0.468	0.457	2.4	70	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG091720\
Data File : BG046658.D
Acq On : 17 Sep 2020 11:12
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 17 14:36:23 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 17 14:27:03 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	1.330	-1.5	76	0.00
46		1,1'-Biphenyl	1.391	1.407	-1.2	74	0.00
47		2-Chloronaphthalene	1.180	1.180	0.0	74	0.00
48		2-Nitroaniline	0.328	0.346	-5.5	75	0.00
49		Acenaphthylene	1.790	1.866	-4.2	76	0.00
50		Dimethylphthalate	1.547	1.637	-5.8	77	0.00
51		2,6-Dinitrotoluene	0.341	0.359	-5.3	76	0.00
52	C	Acenaphthene	1.109	1.137	-2.5	75	0.00
53		3-Nitroaniline	0.369	0.392	-6.2	75	0.00
54	P	2,4-Dinitrophenol	0.199	0.350	-75.9#	115	0.00
55		Dibenzofuran	1.780	1.870	-5.1	78	0.00
56	P	4-Nitrophenol	0.272	0.274	-0.7	69	0.00
57		2,4-Dinitrotoluene	0.486	0.527	-8.4	77	0.00
58		Fluorene	1.494	1.564	-4.7	77	0.00
59		2,3,4,6-Tetrachlorophenol	0.454	0.482	-6.2	76	0.00
60		Diethylphthalate	1.509	1.626	-7.8	79	0.00
61		4-Chlorophenyl-phenylether	0.823	0.856	-4.0	76	0.00
62		4-Nitroaniline	0.390	0.424	-8.7	78	0.00
63		Azobenzene	1.227	1.299	-5.9	78	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	69	0.00
65		4,6-Dinitro-2-methylphenol	0.113	0.121	-7.1	74	0.00
66	c	n-Nitrosodiphenylamine	0.540	0.562	-4.1	78	0.00
67		4-Bromophenyl-phenylether	0.233	0.245	-5.2	78	0.00
68		Hexachlorobenzene	0.258	0.264	-2.3	77	0.00
69		Atrazine	0.220	0.232	-5.5	77	0.00
70	C	Pentachlorophenol	0.135	0.126	6.7	64	0.00
71		Phenanthrene	1.014	1.051	-3.6	79	0.00
72		Anthracene	1.000	1.035	-3.5	79	0.00
73		Carbazole	0.944	0.976	-3.4	78	0.00
74		Di-n-butylphthalate	1.086	1.130	-4.1	79	0.00
75	C	Fluoranthene	1.305	1.319	-1.1	78	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	71	0.00
77		Benzidine	0.465	0.394	15.3	56	0.00
78		Pyrene	1.274	1.318	-3.5	79	0.00
79	S	Terphenyl-d14	0.940	0.945	-0.5	79	0.00
80		Butylbenzylphthalate	0.497	0.512	-3.0	76	0.00
81		Benzo(a)anthracene	1.247	1.270	-1.8	79	0.00
82		3,3'-Dichlorobenzidine	0.463	0.472	-1.9	77	0.00
83		Chrysene	1.199	1.232	-2.8	80	0.00
84		Bis(2-ethylhexyl)phthalate	0.678	0.710	-4.7	79	0.00
85	c	Di-n-octyl phthalate	1.161	1.235	-6.4	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	1.486	-3.5	79	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG091720\
Data File : BG046658.D
Acq On : 17 Sep 2020 11:12
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Sep 17 14:36:23 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082520.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 17 14:27:03 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	1.277	-2.2	79	0.00
89	Benzo(k)fluoranthene	1.207	1.249	-3.5	80	0.00
90 C	Benzo(a)pyrene	1.165	1.215	-4.3	80	0.00
91	Dibenzo(a,h)anthracene	1.159	1.200	-3.5	79	0.00
92	Benzo(g,h,i)perylene	1.158	1.192	-2.9	79	0.00

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