Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021022\
Data File : BG052348.D
Acq On : 10 Feb 2022 10:25
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Feb 10 12:06:41 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 08 01:21:20 2022
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	74	0.01
2		1,4-Dioxane	0.528	0.460	12.9	68	0.00
3		Pyridine	1.532	1.424	7.0	68	0.00
4		n-Nitrosodimethylamine	0.791	0.711	10.1	68	0.00
5	S	2-Fluorophenol	1.148	1.108	3.5	72	0.01
6		Aniline	2.057	2.110	-2.6	74	0.00
7	S	Phenol-d6	1.771	1.805	-1.9	74	0.00
8		2-Chlorophenol	1.262	1.261	0.1	76	0.00
9		Benzaldehyde	1.087	0.956	12.1	69	0.00
10	C	Phenol	1.759	1.766	-0.4	74	0.01
11		bis(2-Chloroethyl)ether	1.345	1.285	4.5	73	0.00
12		1,3-Dichlorobenzene	1.438	1.400	2.6	73	0.00
13	C	1,4-Dichlorobenzene	1.466	1.457	0.6	75	0.00
14		1,2-Dichlorobenzene	1.446	1.403	3.0	73	0.00
15		Benzyl Alcohol	1.469	1.527	-3.9	74	0.00
16		2,2'-oxybis(1-Chloropropane	1.942	1.797	7.5	71	0.01
17		2-Methylphenol	1.161	1.213	-4.5	77	0.01
18		Hexachloroethane	0.510	0.490	3.9	74	0.01
19	P	n-Nitroso-di-n-propylamine	1.335	1.358	-1.7	76	0.01
20		3+4-Methylphenols	1.661	1.793	-7.9	79	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	78	0.00
22		Acetophenone	0.529	0.500	5.5	76	0.01
23	S	Nitrobenzene-d5	0.460	0.433	5.9	75	0.01
24		Nitrobenzene	0.437	0.408	6.6	73	0.00
25		Isophorone	0.783	0.791	-1.0	80	0.00
26	C	2-Nitrophenol	0.185	0.189	-2.2	78	0.01
27		2,4-Dimethylphenol	0.274	0.288	-5.1	82	0.00
28		bis(2-Chloroethoxy)methane	0.444	0.429	3.4	77	0.00
29	C	2,4-Dichlorophenol	0.328	0.333	-1.5	80	0.00
30		1,2,4-Trichlorobenzene	0.383	0.363	5.2	76	0.00
31		Naphthalene	1.048	1.011	3.5	77	0.01
32		Benzoic acid	0.161	0.190	-18.0	92	0.00
33		4-Chloroaniline	0.438	0.449	-2.5	80	0.01
34	C	Hexachlorobutadiene	0.259	0.237	8.5	75	0.01
35		Caprolactam	0.120	0.134	-11.7	87	0.01
36	C	4-Chloro-3-methylphenol	0.374	0.405	-8.3	84	0.00
37		2-Methylnaphthalene	0.779	0.775	0.5	80	0.00
38		1-Methylnaphthalene	0.755	0.758	-0.4	79	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	84	0.00
40		1,2,4,5-Tetrachlorobenzene	0.658	0.596	9.4	79	0.00
41	P	Hexachlorocyclopentadiene	0.278	0.246	11.5	75	0.00
42	S	2,4,6-Tribromophenol	0.287	0.302	-5.2	89	0.00
43	C	2,4,6-Trichlorophenol	0.430	0.428	0.5	84	0.00
44		2,4,5-Trichlorophenol	0.466	0.461	1.1	84	0.00
45	S	2-Fluorobiphenyl	1.439	1.352	6.0	81	0.01
46		1,1'-Biphenyl	1.467	1.393	5.0	82	0.01
47		2-Chloronaphthalene	1.130	1.066	5.7	81	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021022\
Data File : BG052348.D
Acq On : 10 Feb 2022 10:25
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Feb 10 12:06:41 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 08 01:21:20 2022
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)
48		2-Nitroaniline	0.402	0.403	-0.2	85	0.00
49		Acenaphthylene	1.839	1.783	3.0	84	0.00
50		Dimethylphthalate	1.532	1.534	-0.1	87	0.00
51		2,6-Dinitrotoluene	0.331	0.325	1.8	83	0.00
52	C	Acenaphthene	1.205	1.180	2.1	84	0.00
53		3-Nitroaniline	0.344	0.358	-4.1	87	0.00
54	P	2,4-Dinitrophenol	0.179	0.186	-3.9	85	0.00
55		Dibenzofuran	1.894	1.826	3.6	85	0.00
56	P	4-Nitrophenol	0.258	0.270	-4.7	86	0.00
57		2,4-Dinitrotoluene	0.471	0.493	-4.7	88	0.01
58		Fluorene	1.512	1.517	-0.3	87	0.00
59		2,3,4,6-Tetrachlorophenol	0.445	0.463	-4.0	89	0.00
60		Diethylphthalate	1.548	1.548	0.0	87	0.00
61		4-Chlorophenyl-phenylether	0.861	0.831	3.5	85	0.00
62		4-Nitroaniline	0.362	0.367	-1.4	86	0.00
63		Azobenzene	1.584	1.529	3.5	84	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
65		4,6-Dinitro-2-methylphenol	0.120	0.127	-5.8	91	0.00
66	c	n-Nitrosodiphenylamine	0.551	0.520	5.6	85	0.00
67		4-Bromophenyl-phenylether	0.243	0.229	5.8	86	0.00
68		Hexachlorobenzene	0.264	0.241	8.7	84	0.00
69		Atrazine	0.223	0.217	2.7	88	0.00
70	C	Pentachlorophenol	0.129	0.128	0.8	87	0.00
71		Phenanthrene	1.031	0.973	5.6	87	0.00
72		Anthracene	1.010	0.945	6.4	86	0.00
73		Carbazole	0.985	0.944	4.2	88	0.00
74		Di-n-butylphthalate	1.074	1.047	2.5	89	0.00
75	C	Fluoranthene	1.313	1.224	6.8	87	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	90	0.00
77		Benzidine	0.427	0.488	-14.3	89	0.00
78		Pyrene	1.338	1.312	1.9	86	0.00
79	S	Terphenyl-d14	1.133	1.101	2.8	87	0.00
80		Butylbenzylphthalate	0.465	0.484	-4.1	90	0.00
81		Benzo(a)anthracene	1.323	1.262	4.6	85	0.00
82		3,3'-Dichlorobenzidine	0.462	0.439	5.0	83	0.00
83		Chrysene	1.254	1.196	4.6	86	0.00
84		Bis(2-ethylhexyl)phthalate	0.645	0.655	-1.6	86	0.00
85	c	Di-n-octyl phthalate	1.081	1.095	-1.3	87	0.00
86	I	Perylene-d12	1.000	1.000	0.0	88	0.01
87		Indeno(1,2,3-cd)pyrene	1.464	1.385	5.4	83	0.03
88		Benzo(b)fluoranthene	1.246	1.186	4.8	84	0.01
89		Benzo(k)fluoranthene	1.231	1.176	4.5	83	0.02
90	C	Benzo(a)pyrene	1.053	1.001	4.9	83	0.01
91		Dibenzo(a,h)anthracene	1.209	1.137	6.0	83	0.02
92		Benzo(g,h,i)perylene	1.187	1.125	5.2	84	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG021022\
Data File : BG052348.D
Acq On : 10 Feb 2022 10:25
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Feb 10 12:06:41 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 08 01:21:20 2022
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)

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