Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081018\
Data File : BG036269.D
Acq On : 10 Aug 2018 16:06
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 10 16:57:02 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 31 12:30:06 2018
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	65	-0.02
2		1,4-Dioxane	0.613	0.532	13.2	61	0.00
3		Pyridine	1.601	1.384	13.6	54	0.00
4		n-Nitrosodimethylamine	0.687	0.578	15.9	55	0.00
5	S	2-Fluorophenol	1.205	1.122	6.9	60	0.00
6		Aniline	2.330	2.028	13.0	57	-0.02
7	S	Phenol-d6	1.797	1.665	7.3	60	-0.01
8		2-Chlorophenol	1.225	1.156	5.6	61	-0.02
9		Benzaldehyde	1.103	0.928	15.9	54	-0.01
10	C	Phenol	1.816	1.714	5.6	61	-0.01
11		bis(2-Chloroethyl)ether	1.422	1.279	10.1	59	-0.02
12		1,3-Dichlorobenzene	1.480	1.417	4.3	63	-0.03
13	C	1,4-Dichlorobenzene	1.503	1.465	2.5	64	-0.03
14		1,2-Dichlorobenzene	1.444	1.395	3.4	64	-0.03
15		Benzyl Alcohol	1.632	1.396	14.5	55	-0.01
16		2,2'-oxybis(1-Chloropropane	1.192	1.115	6.5	61	-0.02
17		2-Methylphenol	1.235	1.097	11.2	57	-0.01
18		Hexachloroethane	0.575	0.559	2.8	65	-0.02
19	P	n-Nitroso-di-n-propylamine	1.513	1.244	17.8	55	-0.02
20		3+4-Methylphenols	1.713	1.594	6.9	58	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	61	-0.03
22		Acetophenone	0.622	0.566	9.0	57	-0.02
23	S	Nitrobenzene-d5	0.479	0.454	5.2	58	-0.02
24		Nitrobenzene	0.481	0.435	9.6	57	-0.01
25		Isophorone	0.889	0.752	15.4	53	-0.02
26	C	2-Nitrophenol	0.171	0.180	-5.3	63	-0.02
27		2,4-Dimethylphenol	0.289	0.282	2.4	59	-0.02
28		bis(2-Chloroethoxy)methane	0.490	0.420	14.3	53	-0.03
29	C	2,4-Dichlorophenol	0.330	0.351	-6.4	63	-0.01
30		1,2,4-Trichlorobenzene	0.405	0.420	-3.7	65	-0.02
31		Naphthalene	1.042	0.976	6.3	60	-0.02
32		Benzoic acid	0.195	0.141	27.7#	44#	-0.01
33		4-Chloroaniline	0.454	0.407	10.4	57	-0.02
34	C	Hexachlorobutadiene	0.316	0.344	-8.9	68	-0.03
35		Caprolactam	0.142	0.117	17.6	53	-0.02
36	C	4-Chloro-3-methylphenol	0.443	0.434	2.0	60	-0.01
37		2-Methylnaphthalene	0.776	0.739	4.8	59	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	61	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.755	0.768	-1.7	67	-0.02
40	P	Hexachlorocyclopentadiene	0.396	0.442	-11.6	70	-0.02
41	S	2,4,6-Tribromophenol	0.264	0.279	-5.7	68	-0.02
42	C	2,4,6-Trichlorophenol	0.441	0.445	-0.9	62	-0.01
43		2,4,5-Trichlorophenol	0.494	0.484	2.0	65	-0.01
44	S	2-Fluorobiphenyl	1.493	1.423	4.7	63	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081018\
Data File : BG036269.D
Acq On : 10 Aug 2018 16:06
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 10 16:57:02 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 31 12:30:06 2018
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		1,1'-Biphenyl	1.551	1.438	7.3	60	-0.03
46		2-Chloronaphthalene	1.206	1.140	5.5	62	-0.02
47		2-Nitroaniline	0.426	0.378	11.3	56	-0.02
48		Acenaphthylene	2.020	1.861	7.9	60	-0.02
49		Dimethylphthalate	1.752	1.581	9.8	60	-0.02
50		2,6-Dinitrotoluene	0.357	0.358	-0.3	63	-0.02
51	C	Acenaphthene	1.259	1.152	8.5	60	-0.02
52		3-Nitroaniline	0.365	0.314	14.0	54	-0.01
53	P	2,4-Dinitrophenol	0.158	0.139	12.0	56	0.00
54		Dibenzofuran	2.002	1.882	6.0	61	-0.02
55	P	4-Nitrophenol	0.260	0.218	16.2	53	0.00
56		2,4-Dinitrotoluene	0.512	0.517	-1.0	62	-0.01
57		Fluorene	1.678	1.593	5.1	63	-0.02
58		2,3,4,6-Tetrachlorophenol	0.473	0.462	2.3	62	-0.02
59		Diethylphthalate	1.802	1.603	11.0	58	-0.02
60		4-Chlorophenyl-phenylether	0.986	0.958	2.8	64	-0.03
61		4-Nitroaniline	0.382	0.311	18.6	51	-0.02
62		Azobenzene	1.777	1.531	13.8	56	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	61	-0.02
64		4,6-Dinitro-2-methylphenol	0.111	0.105	5.4	59	-0.01
65	c	n-Nitrosodiphenylamine	0.569	0.546	4.0	60	-0.02
66		4-Bromophenyl-phenylether	0.232	0.238	-2.6	64	-0.02
67		Hexachlorobenzene	0.239	0.244	-2.1	64	-0.01
68		Atrazine	0.234	0.206	12.0	53	-0.01
69	C	Pentachlorophenol	0.146	0.135	7.5	56	-0.01
70		Phenanthrene	0.998	0.956	4.2	61	-0.02
71		Anthracene	0.994	0.955	3.9	61	-0.02
72		Carbazole	0.944	0.857	9.2	57	-0.01
73		Di-n-butylphthalate	1.115	1.024	8.2	57	-0.02
74	C	Fluoranthene	1.344	1.273	5.3	60	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	58	-0.02
76		Benzidine	0.526	0.258	51.0#	31#	-0.01
77		Pyrene	1.265	1.258	0.6	61	-0.02
78	S	Terphenyl-d14	0.907	0.944	-4.1	63	-0.02
79		Butylbenzylphthalate	0.452	0.458	-1.3	59	-0.01
80		Benzo(a)anthracene	1.246	1.211	2.8	59	-0.03
81		3,3'-Dichlorobenzidine	0.435	0.416	4.4	57	-0.03
82		Chrysene	1.155	1.152	0.3	61	-0.02
83		Bis(2-ethylhexyl)phthalate	0.615	0.630	-2.4	60	-0.03
84	c	Di-n-octyl phthalate	1.029	1.055	-2.5	60	-0.03
85		Indeno(1,2,3-cd)pyrene	1.279	1.229	3.9	57	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	59	-0.04
87		Benzo(b)fluoranthene	1.216	1.188	2.3	61	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081018\
Data File : BG036269.D
Acq On : 10 Aug 2018 16:06
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 10 16:57:02 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 31 12:30:06 2018
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(k)fluoranthene	1.188	1.155	2.8	59	-0.03
89 C	Benzo(a)pyrene	1.131	1.099	2.8	59	-0.04
90	Dibenzo(a,h)anthracene	1.110	1.015	8.6	55	-0.06
91	Benzo(g,h,i)perylene	1.086	1.007	7.3	56	-0.05

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