Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG102317\
Data File : BG029401.D
Acq On : 23 Oct 2017 16:14
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 24 05:50:49 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 17:32:15 2017
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	90	0.00
2		1,4-Dioxane	0.486	0.544	-11.9	93	0.00
3		Pyridine	1.415	1.539	-8.8	90	-0.01
4		n-Nitrosodimethylamine	0.661	0.700	-5.9	89	0.00
5	S	2-Fluorophenol	1.197	1.288	-7.6	89	0.00
6		Aniline	2.081	2.143	-3.0	84	0.00
7	S	Phenol-d6	1.680	1.766	-5.1	86	-0.01
8		2-Chlorophenol	1.372	1.415	-3.1	87	-0.01
9		Benzaldehyde	0.845	0.824	2.5	80	0.00
10	C	Phenol	1.812	1.899	-4.8	86	-0.01
11		bis(2-Chloroethyl)ether	1.320	1.397	-5.8	86	0.00
12		1,3-Dichlorobenzene	1.541	1.576	-2.3	86	-0.01
13	C	1,4-Dichlorobenzene	1.573	1.623	-3.2	85	-0.01
14		1,2-Dichlorobenzene	1.517	1.541	-1.6	85	0.00
15		Benzyl Alcohol	1.184	1.235	-4.3	86	-0.01
16		2,2'-oxybis(1-Chloropropane	2.242	2.312	-3.1	85	0.00
17		2-Methylphenol	1.204	1.239	-2.9	85	-0.01
18		Hexachloroethane	0.536	0.556	-3.7	86	0.00
19	P	n-Nitroso-di-n-propylamine	1.143	1.200	-5.0	87	0.00
20		3+4-Methylphenols	1.696	1.783	-5.1	87	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	85	-0.01
22		Acetophenone	0.482	0.516	-7.1	86	0.00
23	S	Nitrobenzene-d5	0.315	0.338	-7.3	85	-0.01
24		Nitrobenzene	0.354	0.379	-7.1	85	0.00
25		Isophorone	0.657	0.710	-8.1	86	-0.01
26	C	2-Nitrophenol	0.169	0.190	-12.4	86	-0.01
27		2,4-Dimethylphenol	0.306	0.332	-8.5	87	-0.01
28		bis(2-Chloroethoxy)methane	0.403	0.436	-8.2	87	-0.01
29	C	2,4-Dichlorophenol	0.326	0.355	-8.9	87	0.00
30		1,2,4-Trichlorobenzene	0.353	0.373	-5.7	85	-0.01
31		Naphthalene	0.997	1.027	-3.0	83	-0.01
32		Benzoic acid	0.256	0.328	-28.1#	96	-0.01
33		4-Chloroaniline	0.419	0.456	-8.8	87	-0.01
34	C	Hexachlorobutadiene	0.219	0.242	-10.5	89	0.00
35		Caprolactam	0.108	0.139	-28.7#	103	-0.01
36	C	4-Chloro-3-methylphenol	0.359	0.409	-13.9	91	-0.01
37		2-Methylnaphthalene	0.726	0.759	-4.5	84	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	95	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.730	0.730	0.0	89	-0.01
40	P	Hexachlorocyclopentadiene	0.247	0.311	-25.9#	109	0.00
41	S	2,4,6-Tribromophenol	0.258	0.311	-20.5	104	0.00
42	C	2,4,6-Trichlorophenol	0.458	0.479	-4.6	92	0.00
43		2,4,5-Trichlorophenol	0.471	0.523	-11.0	98	-0.01
44	S	2-Fluorobiphenyl	1.364	1.353	0.8	89	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG102317\
Data File : BG029401.D
Acq On : 23 Oct 2017 16:14
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 24 05:50:49 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 17:32:15 2017
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	1.719	1.695	1.4	88	-0.01
46		2-Chloronaphthalene	1.254	1.226	2.2	87	-0.01
47		2-Nitroaniline	0.344	0.390	-13.4	96	0.00
48		Acenaphthylene	1.962	1.975	-0.7	89	-0.01
49		Dimethylphthalate	1.560	1.754	-12.4	98	-0.01
50		2,6-Dinitrotoluene	0.327	0.383	-17.1	98	0.00
51	C	Acenaphthene	1.277	1.295	-1.4	89	-0.01
52		3-Nitroaniline	0.353	0.415	-17.6	100	0.00
53	P	2,4-Dinitrophenol	0.153	0.139	9.2	80	0.00
54		Dibenzofuran	1.823	1.901	-4.3	93	0.00
55	P	4-Nitrophenol	0.291	0.336	-15.5	98	0.00
56		2,4-Dinitrotoluene	0.443	0.561	-26.6#	107	-0.01
57		Fluorene	1.506	1.603	-6.4	95	0.00
58		2,3,4,6-Tetrachlorophenol	0.393	0.472	-20.1	104	0.00
59		Diethylphthalate	1.555	1.780	-14.5	101	0.00
60		4-Chlorophenyl-phenylether	0.803	0.902	-12.3	99	-0.01
61		4-Nitroaniline	0.362	0.425	-17.4	102	-0.01
62		Azobenzene	1.311	1.403	-7.0	95	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	-0.01
64		4,6-Dinitro-2-methylphenol	0.105	0.121	-15.2	116	-0.01
65	c	n-Nitrosodiphenylamine	0.599	0.582	2.8	100	0.00
66		4-Bromophenyl-phenylether	0.229	0.230	-0.4	103	0.00
67		Hexachlorobenzene	0.260	0.263	-1.2	105	0.00
68		Atrazine	0.214	0.236	-10.3	111	0.00
69	C	Pentachlorophenol	0.149	0.160	-7.4	105	-0.01
70		Phenanthrene	1.033	1.056	-2.2	106	0.00
71		Anthracene	1.037	1.072	-3.4	106	-0.01
72		Carbazole	0.997	1.053	-5.6	108	0.00
73		Di-n-butylphthalate	1.146	1.232	-7.5	110	-0.01
74	C	Fluoranthene	1.208	1.320	-9.3	112	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	116	-0.01
76		Benzidine	0.604	0.542	10.3	95	-0.01
77		Pyrene	1.213	1.236	-1.9	112	0.00
78	S	Terphenyl-d14	0.879	0.892	-1.5	111	-0.01
79		Butylbenzylphthalate	0.517	0.524	-1.4	110	0.00
80		Benzo(a)anthracene	1.174	1.248	-6.3	116	-0.01
81		3,3'-Dichlorobenzidine	0.485	0.485	0.0	110	-0.01
82		Chrysene	1.125	1.189	-5.7	117	-0.01
83		Bis(2-ethylhexyl)phthalate	0.742	0.751	-1.2	110	-0.02
84	c	Di-n-octyl phthalate	1.293	1.305	-0.9	110	-0.02
85		Indeno(1,2,3-cd)pyrene	1.383	1.504	-8.7	119	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	114	-0.02
87		Benzo(b)fluoranthene	1.145	1.220	-6.6	114	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG102317\
Data File : BG029401.D
Acq On : 23 Oct 2017 16:14
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Oct 24 05:50:49 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 17:32:15 2017
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	1.141	1.241	-8.8	117	-0.02
89 C	Benzo(a)pyrene	1.101	1.182	-7.4	115	-0.02
90	Dibenzo(a,h)anthracene	1.114	1.213	-8.9	115	-0.03
91	Benzo(g,h,i)perylene	1.035	1.212	-17.1	123	-0.04

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