Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG031218\
Data File : BG033113.D
Acq On : 13 Mar 2018 6:38
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 13 09:01:30 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 09 15:10:14 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	67	0.00
2		1,4-Dioxane	0.543	0.495	8.8	61	0.00
3		Pyridine	1.495	1.478	1.1	63	0.00
4		n-Nitrosodimethylamine	0.600	0.685	-14.2	74	0.00
5	S	2-Fluorophenol	1.250	1.214	2.9	63	0.00
6		Aniline	2.359	2.262	4.1	64	0.00
7	S	Phenol-d6	1.742	1.711	1.8	64	0.00
8		2-Chlorophenol	1.368	1.305	4.6	62	0.00
9		Benzaldehyde	1.004	0.841	16.2	57	0.00
10	C	Phenol	1.757	1.735	1.3	65	0.00
11		bis(2-Chloroethyl)ether	1.436	1.311	8.7	61	0.00
12		1,3-Dichlorobenzene	1.603	1.479	7.7	61	0.00
13	C	1,4-Dichlorobenzene	1.613	1.510	6.4	62	0.00
14		1,2-Dichlorobenzene	1.546	1.455	5.9	63	0.00
15		Benzyl Alcohol	1.281	1.306	-2.0	68	0.00
16		2,2'-oxybis(1-Chloropropane	2.469	2.646	-7.2	72	0.00
17		2-Methylphenol	1.295	1.275	1.5	65	0.00
18		Hexachloroethane	0.549	0.543	1.1	65	0.00
19	P	n-Nitroso-di-n-propylamine	1.230	1.193	3.0	65	0.00
20		3+4-Methylphenols	1.801	1.798	0.2	67	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	71	0.00
22		Acetophenone	0.505	0.467	7.5	65	0.00
23	S	Nitrobenzene-d5	0.242	0.331	-36.8#	90	0.00
24		Nitrobenzene	0.289	0.351	-21.5	83	0.00
25		Isophorone	0.702	0.631	10.1	64	0.00
26	C	2-Nitrophenol	0.103	0.161	-56.3#	116	0.00
27		2,4-Dimethylphenol	0.305	0.270	11.5	64	0.00
28		bis(2-Chloroethoxy)methane	0.409	0.357	12.7	62	0.00
29	C	2,4-Dichlorophenol	0.277	0.270	2.5	69	0.00
30		1,2,4-Trichlorobenzene	0.324	0.294	9.3	64	0.00
31		Naphthalene	1.045	0.951	9.0	65	0.00
32		Benzoic acid	0.174	0.201	-15.5	88	0.00
33		4-Chloroaniline	0.467	0.413	11.6	63	0.00
34	C	Hexachlorobutadiene	0.182	0.171	6.0	66	0.00
35		Caprolactam	0.111	0.112	-0.9	70	0.00
36	C	4-Chloro-3-methylphenol	0.322	0.339	-5.3	73	0.00
37		2-Methylnaphthalene	0.756	0.693	8.3	65	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	78	0.00
39		1,2,4,5-Tetrachlorobenzene	0.594	0.546	8.1	72	0.00
40	P	Hexachlorocyclopentadiene	0.303	0.277	8.6	70	0.00
41	S	2,4,6-Tribromophenol	0.210	0.240	-14.3	87	0.00
42	C	2,4,6-Trichlorophenol	0.374	0.363	2.9	75	0.00
43		2,4,5-Trichlorophenol	0.433	0.423	2.3	75	0.00
44	S	2-Fluorobiphenyl	1.387	1.224	11.8	69	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG031218\
Data File : BG033113.D
Acq On : 13 Mar 2018 6:38
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 13 09:01:30 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 09 15:10:14 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	1.748	1.518	13.2	68	0.00
46		2-Chloronaphthalene	1.306	1.129	13.6	67	0.00
47		2-Nitroaniline	0.282	0.425	-50.7#	118	0.00
48		Acenaphthylene	2.137	1.926	9.9	70	0.00
49		Dimethylphthalate	1.698	1.508	11.2	69	0.00
50		2,6-Dinitrotoluene	0.244	0.329	-34.8#	104	0.00
51	C	Acenaphthene	1.331	1.197	10.1	70	0.00
52		3-Nitroaniline	0.308	0.367	-19.2	92	0.00
53	P	2,4-Dinitrophenol	0.070	0.083	-18.6	96	0.00
54		Dibenzofuran	1.895	1.701	10.2	70	0.00
55	P	4-Nitrophenol	0.233	0.217	6.9	71	0.00
56		2,4-Dinitrotoluene	0.293	0.459	-56.7#	127	0.00
57		Fluorene	1.564	1.405	10.2	71	0.00
58		2,3,4,6-Tetrachlorophenol	0.336	0.348	-3.6	79	0.00
59		Diethylphthalate	1.791	1.631	8.9	71	0.00
60		4-Chlorophenyl-phenylether	0.765	0.703	8.1	72	0.00
61		4-Nitroaniline	0.336	0.355	-5.7	80	0.00
62		Azobenzene	1.602	1.467	8.4	72	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	81	0.00
64		4,6-Dinitro-2-methylphenol	0.054	0.089	-64.8#	141	0.00
65	c	n-Nitrosodiphenylamine	0.651	0.570	12.4	71	0.00
66		4-Bromophenyl-phenylether	0.211	0.194	8.1	75	0.00
67		Hexachlorobenzene	0.229	0.214	6.6	77	0.00
68		Atrazine	0.214	0.187	12.6	70	0.00
69	C	Pentachlorophenol	0.144	0.136	5.6	76	0.00
70		Phenanthrene	1.116	1.001	10.3	73	0.00
71		Anthracene	1.120	0.998	10.9	72	0.00
72		Carbazole	1.104	0.954	13.6	70	0.00
73		Di-n-butylphthalate	1.355	1.213	10.5	71	0.00
74	C	Fluoranthene	1.233	1.125	8.8	74	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	84	0.00
76		Benzidine	0.682	0.224	67.2#	29#	0.00
77		Pyrene	1.410	1.230	12.8	74	0.00
78	S	Terphenyl-d14	0.890	0.804	9.7	77	0.00
79		Butylbenzylphthalate	0.625	0.602	3.7	78	0.00
80		Benzo(a)anthracene	1.283	1.164	9.3	77	0.00
81		3,3'-Dichlorobenzidine	0.475	0.412	13.3	71	0.00
82		Chrysene	1.225	1.119	8.7	77	0.00
83		Bis(2-ethylhexyl)phthalate	0.926	0.863	6.8	75	0.00
84	c	Di-n-octyl phthalate	1.411	1.367	3.1	76	0.00
85		Indeno(1,2,3-cd)pyrene	1.429	1.250	12.5	72	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	88	0.00
87		Benzo(b)fluoranthene	1.183	1.034	12.6	75	0.08

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG031218\
Data File : BG033113.D
Acq On : 13 Mar 2018 6:38
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 13 09:01:30 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 09 15:10:14 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	1.175	1.034	12.0	78	0.00
89 C	Benzo(a)pyrene	1.125	0.988	12.2	77	0.00
90	Dibenzo(a,h)anthracene	1.132	0.924	18.4	70	0.00
91	Benzo(g,h,i)perylene	1.116	0.923	17.3	72	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1