Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG030118\
Data File : BG032918.D
Acq On : 1 Mar 2018 15:43
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 01 17:41:58 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 01 17:38:52 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	102	0.00
2		1,4-Dioxane	0.543	0.528	2.8	99	0.00
3		Pyridine	1.495	1.529	-2.3	100	0.00
4		n-Nitrosodimethylamine	0.600	0.635	-5.8	106	0.00
5	S	2-Fluorophenol	1.250	1.250	0.0	99	0.00
6		Aniline	2.359	2.269	3.8	98	0.00
7	S	Phenol-d6	1.742	1.750	-0.5	101	0.00
8		2-Chlorophenol	1.368	1.374	-0.4	101	0.00
9		Benzaldehyde	1.004	0.889	11.5	93	0.00
10	C	Phenol	1.757	1.771	-0.8	102	0.00
11		bis(2-Chloroethyl)ether	1.436	1.367	4.8	97	0.00
12		1,3-Dichlorobenzene	1.603	1.534	4.3	97	0.00
13	C	1,4-Dichlorobenzene	1.613	1.540	4.5	97	0.00
14		1,2-Dichlorobenzene	1.546	1.471	4.9	97	0.00
15		Benzyl Alcohol	1.281	1.266	1.2	101	0.00
16		2,2'-oxybis(1-Chloropropane	2.469	2.455	0.6	102	0.00
17		2-Methylphenol	1.295	1.288	0.5	100	0.00
18		Hexachloroethane	0.549	0.554	-0.9	102	0.00
19	P	n-Nitroso-di-n-propylamine	1.230	1.199	2.5	100	0.00
20		3+4-Methylphenols	1.801	1.826	-1.4	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
22		Acetophenone	0.505	0.482	4.6	101	0.00
23	S	Nitrobenzene-d5	0.242	0.328	-35.5#	133	0.00
24		Nitrobenzene	0.289	0.345	-19.4	123	0.00
25		Isophorone	0.702	0.661	5.8	101	0.00
26	C	2-Nitrophenol	0.103	0.160	-55.3#	174#	0.00
27		2,4-Dimethylphenol	0.305	0.292	4.3	104	0.00
28		bis(2-Chloroethoxy)methane	0.409	0.387	5.4	101	0.00
29	C	2,4-Dichlorophenol	0.277	0.277	0.0	106	0.00
30		1,2,4-Trichlorobenzene	0.324	0.305	5.9	100	0.00
31		Naphthalene	1.045	0.988	5.5	101	0.00
32		Benzoic acid	0.174	0.186	-6.9	122	0.00
33		4-Chloroaniline	0.467	0.438	6.2	100	0.00
34	C	Hexachlorobutadiene	0.182	0.169	7.1	98	0.00
35		Caprolactam	0.111	0.110	0.9	103	0.00
36	C	4-Chloro-3-methylphenol	0.322	0.328	-1.9	106	0.00
37		2-Methylnaphthalene	0.756	0.717	5.2	101	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
39		1,2,4,5-Tetrachlorobenzene	0.594	0.550	7.4	103	0.00
40	P	Hexachlorocyclopentadiene	0.303	0.294	3.0	105	0.00
41	S	2,4,6-Tribromophenol	0.210	0.208	1.0	107	0.00
42	C	2,4,6-Trichlorophenol	0.374	0.378	-1.1	111	0.00
43		2,4,5-Trichlorophenol	0.433	0.441	-1.8	111	0.00
44	S	2-Fluorobiphenyl	1.387	1.301	6.2	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG030118\
Data File : BG032918.D
Acq On : 1 Mar 2018 15:43
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 01 17:41:58 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 01 17:38:52 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	1.636	6.4	103	0.00
46		2-Chloronaphthalene	1.306	1.219	6.7	103	0.00
47		2-Nitroaniline	0.282	0.400	-41.8#	157#	0.00
48		Acenaphthylene	2.137	2.006	6.1	103	0.00
49		Dimethylphthalate	1.698	1.554	8.5	101	0.00
50		2,6-Dinitrotoluene	0.244	0.334	-36.9#	149	0.00
51	C	Acenaphthene	1.331	1.242	6.7	103	0.00
52		3-Nitroaniline	0.308	0.375	-21.8	133	0.00
53	P	2,4-Dinitrophenol	0.070	0.060	14.3	98	0.00
54		Dibenzofuran	1.895	1.748	7.8	102	0.00
55	P	4-Nitrophenol	0.233	0.237	-1.7	110	0.00
56		2,4-Dinitrotoluene	0.293	0.451	-53.9#	176#	0.00
57		Fluorene	1.564	1.440	7.9	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.336	0.334	0.6	107	0.00
59		Diethylphthalate	1.791	1.641	8.4	101	0.00
60		4-Chlorophenyl-phenylether	0.765	0.710	7.2	103	0.00
61		4-Nitroaniline	0.336	0.365	-8.6	116	0.00
62		Azobenzene	1.602	1.459	8.9	101	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
64		4,6-Dinitro-2-methylphenol	0.054	0.080	-48.1#	167#	0.00
65	c	n-Nitrosodiphenylamine	0.651	0.619	4.9	101	0.00
66		4-Bromophenyl-phenylether	0.211	0.198	6.2	101	0.00
67		Hexachlorobenzene	0.229	0.211	7.9	100	0.00
68		Atrazine	0.214	0.192	10.3	94	0.00
69	C	Pentachlorophenol	0.144	0.134	6.9	98	0.00
70		Phenanthrene	1.116	1.043	6.5	100	0.00
71		Anthracene	1.120	1.047	6.5	99	0.00
72		Carbazole	1.104	0.997	9.7	96	0.00
73		Di-n-butylphthalate	1.355	1.269	6.3	98	0.00
74	C	Fluoranthene	1.233	1.129	8.4	98	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
76		Benzidine	0.682	0.446	34.6#	71	0.00
77		Pyrene	1.410	1.306	7.4	98	0.00
78	S	Terphenyl-d14	0.890	0.814	8.5	97	0.00
79		Butylbenzylphthalate	0.625	0.650	-4.0	104	0.00
80		Benzo(a)anthracene	1.283	1.218	5.1	100	0.00
81		3,3'-Dichlorobenzidine	0.475	0.445	6.3	96	0.00
82		Chrysene	1.225	1.164	5.0	100	0.00
83		Bis(2-ethylhexyl)phthalate	0.926	0.945	-2.1	102	0.00
84	c	Di-n-octyl phthalate	1.411	1.515	-7.4	104	0.00
85		Indeno(1,2,3-cd)pyrene	1.429	1.312	8.2	94	0.00
86	I	Perylene-d12	1.000	1.000	0.0	106	0.00
87		Benzo(b)fluoranthene	1.183	1.099	7.1	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\Data\BG030118\
Data File : BG032918.D
Acq On : 1 Mar 2018 15:43
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 01 17:41:58 2018
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 01 17:38:52 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	1.125	4.3	102	0.00
89 C	Benzo(a)pyrene	1.125	1.072	4.7	100	0.00
90	Dibenzo(a,h)anthracene	1.132	1.000	11.7	92	0.00
91	Benzo(g,h,i)perylene	1.116	0.986	11.6	92	0.00

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