Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031919\
Data File : BG040002.D
Acq On : 19 Mar 2019 14:42
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 20 04:07:17 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 04 14:38:15 2019
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	114	-0.02
2		1,4-Dioxane	0.541	0.498	7.9	110	-0.02
3		Pyridine	1.449	1.444	0.3	107	-0.02
4		n-Nitrosodimethylamine	0.687	0.738	-7.4	120	-0.02
5	S	2-Fluorophenol	1.172	1.146	2.2	111	-0.02
6		Aniline	2.290	2.093	8.6	104	-0.02
7	S	Phenol-d6	1.748	1.641	6.1	105	-0.02
8		2-Chlorophenol	1.422	1.327	6.7	106	-0.02
9		Benzaldehyde	1.128	0.916	18.8	92	-0.02
10	C	Phenol	1.894	1.750	7.6	103	-0.02
11		bis(2-Chloroethyl)ether	1.476	1.371	7.1	107	-0.02
12		1,3-Dichlorobenzene	1.609	1.517	5.7	108	-0.02
13	C	1,4-Dichlorobenzene	1.638	1.540	6.0	109	-0.02
14		1,2-Dichlorobenzene	1.583	1.468	7.3	107	-0.02
15		Benzyl Alcohol	1.387	1.264	8.9	101	-0.02
16		2,2'-oxybis(1-Chloropropane	2.953	2.720	7.9	106	-0.01
17		2-Methylphenol	1.207	1.103	8.6	102	-0.02
18		Hexachloroethane	0.588	0.562	4.4	113	-0.02
19	P	n-Nitroso-di-n-propylamine	1.258	1.098	12.7	97	-0.02
20		3+4-Methylphenols	1.677	1.487	11.3	99	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	106	-0.02
22		Acetophenone	0.537	0.509	5.2	99	-0.02
23	S	Nitrobenzene-d5	0.370	0.374	-1.1	105	-0.02
24		Nitrobenzene	0.388	0.383	1.3	103	-0.02
25		Isophorone	0.775	0.729	5.9	97	-0.02
26	C	2-Nitrophenol	0.187	0.184	1.6	99	-0.02
27		2,4-Dimethylphenol	0.291	0.277	4.8	98	-0.02
28		bis(2-Chloroethoxy)methane	0.471	0.435	7.6	96	-0.02
29	C	2,4-Dichlorophenol	0.328	0.316	3.7	98	-0.02
30		1,2,4-Trichlorobenzene	0.369	0.353	4.3	102	-0.02
31		Naphthalene	1.059	0.987	6.8	98	-0.02
32		Benzoic acid	0.181	0.181	0.0	104	0.00
33		4-Chloroaniline	0.449	0.417	7.1	95	-0.02
34	C	Hexachlorobutadiene	0.252	0.245	2.8	103	-0.02
35		Caprolactam	0.125	0.115	8.0	94	-0.02
36	C	4-Chloro-3-methylphenol	0.376	0.349	7.2	94	-0.02
37		2-Methylnaphthalene	0.792	0.720	9.1	95	-0.02
38		1-Methylnaphthalene	0.769	0.697	9.4	94	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	101	-0.02
40		1,2,4,5-Tetrachlorobenzene	0.678	0.642	5.3	96	-0.02
41	P	Hexachlorocyclopentadiene	0.363	0.380	-4.7	104	-0.02
42	S	2,4,6-Tribromophenol	0.233	0.230	1.3	95	-0.02
43	C	2,4,6-Trichlorophenol	0.397	0.395	0.5	97	-0.02
44		2,4,5-Trichlorophenol	0.447	0.424	5.1	91	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031919\
Data File : BG040002.D
Acq On : 19 Mar 2019 14:42
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 20 04:07:17 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 04 14:38:15 2019
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	S	2-Fluorobiphenyl	1.405	1.311	6.7	95	-0.02
46		1,1'-Biphenyl	1.645	1.526	7.2	94	-0.02
47		2-Chloronaphthalene	1.238	1.142	7.8	94	-0.02
48		2-Nitroaniline	0.398	0.408	-2.5	99	-0.01
49		Acenaphthylene	2.031	1.905	6.2	93	-0.02
50		Dimethylphthalate	1.672	1.537	8.1	92	-0.02
51		2,6-Dinitrotoluene	0.355	0.354	0.3	95	-0.02
52	C	Acenaphthene	1.223	1.125	8.0	93	-0.02
53		3-Nitroaniline	0.356	0.352	1.1	96	-0.02
54	P	2,4-Dinitrophenol	0.197	0.219	-11.2	107	-0.01
55		Dibenzofuran	2.006	1.838	8.4	92	-0.02
56	P	4-Nitrophenol	0.319	0.327	-2.5	99	0.00
57		2,4-Dinitrotoluene	0.498	0.504	-1.2	97	-0.01
58		Fluorene	1.577	1.470	6.8	94	-0.02
59		2,3,4,6-Tetrachlorophenol	0.437	0.420	3.9	93	-0.01
60		Diethylphthalate	1.656	1.548	6.5	92	-0.02
61		4-Chlorophenyl-phenylether	0.842	0.768	8.8	92	-0.02
62		4-Nitroaniline	0.386	0.393	-1.8	96	-0.01
63		Azobenzene	1.501	1.437	4.3	96	-0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	104	-0.02
65		4,6-Dinitro-2-methylphenol	0.118	0.121	-2.5	105	-0.01
66	c	n-Nitrosodiphenylamine	0.579	0.528	8.8	92	-0.02
67		4-Bromophenyl-phenylether	0.224	0.202	9.8	93	-0.02
68		Hexachlorobenzene	0.232	0.212	8.6	95	-0.02
69		Atrazine	0.228	0.213	6.6	95	-0.01
70	C	Pentachlorophenol	0.147	0.143	2.7	98	-0.01
71		Phenanthrene	1.102	1.014	8.0	95	-0.01
72		Anthracene	1.074	1.004	6.5	97	-0.02
73		Carbazole	0.968	0.940	2.9	101	-0.01
74		Di-n-butylphthalate	1.139	1.159	-1.8	105	-0.02
75	C	Fluoranthene	1.302	1.287	1.2	104	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	92	-0.02
77		Benzidine	0.635	0.638	-0.5	83	-0.01
78		Pyrene	1.300	1.503	-15.6	104	-0.02
79	S	Terphenyl-d14	0.908	1.034	-13.9	105	-0.02
80		Butylbenzylphthalate	0.504	0.521	-3.4	91	-0.02
81		Benzo(a)anthracene	1.253	1.178	6.0	85	-0.02
82		3,3'-Dichlorobenzidine	0.458	0.399	12.9	76	-0.02
83		Chrysene	1.215	1.124	7.5	85	-0.02
84		Bis(2-ethylhexyl)phthalate	0.708	0.657	7.2	82	-0.02
85	c	Di-n-octyl phthalate	1.172	1.115	4.9	82	-0.03
86		Indeno(1,2,3-cd)pyrene	1.379	1.233	10.6	80	-0.05
87	I	Perylene-d12	1.000	1.000	0.0	83	-0.04

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031919\
Data File : BG040002.D
Acq On : 19 Mar 2019 14:42
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 20 04:07:17 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 04 14:38:15 2019
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(b)fluoranthene	1.193	1.130	5.3	80	-0.03
89	Benzo(k)fluoranthene	1.110	1.037	6.6	79	-0.03
90 C	Benzo(a)pyrene	1.102	1.074	2.5	81	-0.04
91	Dibenzo(a,h)anthracene	1.080	1.009	6.6	79	-0.06
92	Benzo(g,h,i)perylene	1.085	1.028	5.3	79	-0.06

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