Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031219\
Data File : BG039889.D
Acq On : 13 Mar 2019 1:30
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 13 05:50:43 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 13 05:34:05 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	103	-0.02
2		1,4-Dioxane	0.541	0.553	-2.2	110	-0.01
3		Pyridine	1.449	1.519	-4.8	101	-0.01
4		n-Nitrosodimethylamine	0.687	0.722	-5.1	106	-0.01
5	S	2-Fluorophenol	1.172	1.216	-3.8	106	-0.01
6		Aniline	2.290	2.156	5.9	97	-0.02
7	S	Phenol-d6	1.748	1.735	0.7	100	-0.02
8		2-Chlorophenol	1.422	1.392	2.1	100	-0.01
9		Benzaldehyde	1.128	1.006	10.8	91	-0.01
10	C	Phenol	1.894	1.864	1.6	99	-0.01
11		bis(2-Chloroethyl)ether	1.476	1.394	5.6	98	-0.01
12		1,3-Dichlorobenzene	1.609	1.590	1.2	102	-0.01
13	C	1,4-Dichlorobenzene	1.638	1.591	2.9	101	-0.01
14		1,2-Dichlorobenzene	1.583	1.507	4.8	99	-0.01
15		Benzyl Alcohol	1.387	1.300	6.3	93	-0.01
16		2,2'-oxybis(1-Chloropropane	2.953	2.753	6.8	97	-0.01
17		2-Methylphenol	1.207	1.149	4.8	95	-0.02
18		Hexachloroethane	0.588	0.565	3.9	102	-0.02
19	P	n-Nitroso-di-n-propylamine	1.258	1.133	9.9	90	-0.01
20		3+4-Methylphenols	1.677	1.605	4.3	96	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	96	-0.01
22		Acetophenone	0.537	0.523	2.6	92	-0.02
23	S	Nitrobenzene-d5	0.370	0.381	-3.0	97	-0.01
24		Nitrobenzene	0.388	0.392	-1.0	95	-0.01
25		Isophorone	0.775	0.737	4.9	89	-0.01
26	C	2-Nitrophenol	0.187	0.195	-4.3	95	-0.01
27		2,4-Dimethylphenol	0.291	0.293	-0.7	93	-0.01
28		bis(2-Chloroethoxy)methane	0.471	0.450	4.5	90	-0.01
29	C	2,4-Dichlorophenol	0.328	0.341	-4.0	95	-0.01
30		1,2,4-Trichlorobenzene	0.369	0.363	1.6	95	-0.01
31		Naphthalene	1.059	1.032	2.5	92	-0.01
32		Benzoic acid	0.181	0.180	0.6	93	-0.01
33		4-Chloroaniline	0.449	0.451	-0.4	93	-0.01
34	C	Hexachlorobutadiene	0.252	0.252	0.0	95	-0.02
35		Caprolactam	0.125	0.118	5.6	87	-0.02
36	C	4-Chloro-3-methylphenol	0.376	0.375	0.3	91	-0.01
37		2-Methylnaphthalene	0.792	0.763	3.7	91	-0.01
38		1-Methylnaphthalene	0.769	0.744	3.3	91	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	94	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.678	0.660	2.7	92	-0.01
41	P	Hexachlorocyclopentadiene	0.363	0.324	10.7	83	-0.01
42	S	2,4,6-Tribromophenol	0.233	0.239	-2.6	92	-0.01
43	C	2,4,6-Trichlorophenol	0.397	0.408	-2.8	94	-0.01
44		2,4,5-Trichlorophenol	0.447	0.447	0.0	89	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031219\
Data File : BG039889.D
Acq On : 13 Mar 2019 1:30
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 13 05:50:43 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 13 05:34:05 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	1.341	4.6	90	-0.01
46		1,1'-Biphenyl	1.645	1.567	4.7	90	-0.01
47		2-Chloronaphthalene	1.238	1.185	4.3	91	-0.01
48		2-Nitroaniline	0.398	0.414	-4.0	94	0.00
49		Acenaphthylene	2.031	1.977	2.7	90	-0.01
50		Dimethylphthalate	1.672	1.562	6.6	87	-0.01
51		2,6-Dinitrotoluene	0.355	0.353	0.6	89	-0.01
52	C	Acenaphthene	1.223	1.156	5.5	89	-0.01
53		3-Nitroaniline	0.356	0.360	-1.1	91	-0.01
54	P	2,4-Dinitrophenol	0.197	0.157	20.3	71	-0.01
55		Dibenzofuran	2.006	1.924	4.1	90	0.00
56	P	4-Nitrophenol	0.319	0.309	3.1	87	-0.01
57		2,4-Dinitrotoluene	0.498	0.508	-2.0	91	-0.01
58		Fluorene	1.577	1.506	4.5	90	-0.01
59		2,3,4,6-Tetrachlorophenol	0.437	0.435	0.5	90	0.00
60		Diethylphthalate	1.656	1.585	4.3	88	-0.01
61		4-Chlorophenyl-phenylether	0.842	0.801	4.9	89	-0.01
62		4-Nitroaniline	0.386	0.401	-3.9	92	-0.01
63		Azobenzene	1.501	1.471	2.0	92	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	93	-0.01
65		4,6-Dinitro-2-methylphenol	0.118	0.111	5.9	87	0.00
66	c	n-Nitrosodiphenylamine	0.579	0.565	2.4	88	-0.01
67		4-Bromophenyl-phenylether	0.224	0.223	0.4	92	-0.01
68		Hexachlorobenzene	0.232	0.230	0.9	92	-0.01
69		Atrazine	0.228	0.222	2.6	88	0.00
70	C	Pentachlorophenol	0.147	0.139	5.4	86	0.00
71		Phenanthrene	1.102	1.068	3.1	90	0.00
72		Anthracene	1.074	1.063	1.0	92	-0.01
73		Carbazole	0.968	0.973	-0.5	93	-0.01
74		Di-n-butylphthalate	1.139	1.166	-2.4	94	-0.01
75	C	Fluoranthene	1.302	1.321	-1.5	95	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	100	-0.02
77		Benzidine	0.635	0.567	10.7	81	0.00
78		Pyrene	1.300	1.260	3.1	95	-0.01
79	S	Terphenyl-d14	0.908	0.882	2.9	97	-0.01
80		Butylbenzylphthalate	0.504	0.524	-4.0	100	-0.01
81		Benzo(a)anthracene	1.253	1.231	1.8	97	-0.01
82		3,3'-Dichlorobenzidine	0.458	0.443	3.3	93	-0.01
83		Chrysene	1.215	1.165	4.1	96	-0.01
84		Bis(2-ethylhexyl)phthalate	0.708	0.711	-0.4	97	-0.02
85	c	Di-n-octyl phthalate	1.172	1.203	-2.6	97	-0.02
86		Indeno(1,2,3-cd)pyrene	1.379	1.322	4.1	93	-0.04
87	I	Perylene-d12	1.000	1.000	0.0	97	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG031219\
Data File : BG039889.D
Acq On : 13 Mar 2019 1:30
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 13 05:50:43 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG030419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 13 05:34:05 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	1.182	0.9	96	-0.02
89	Benzo(k)fluoranthene	1.110	1.086	2.2	96	-0.02
90 C	Benzo(a)pyrene	1.102	1.104	-0.2	97	-0.02
91	Dibenzo(a,h)anthracene	1.080	1.023	5.3	93	-0.04
92	Benzo(g,h,i)perylene	1.085	1.035	4.6	92	-0.04

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