Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070219\
Data File : BG041565.D
Acq On : 2 Jul 2019 9:22
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 03 07:19:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG062619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 27 13:25:55 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	101	0.00
2		1,4-Dioxane	0.508	0.520	-2.4	98	0.00
3		Pyridine	1.288	1.383	-7.4	107	0.00
4		n-Nitrosodimethylamine	0.717	0.701	2.2	97	0.00
5	S	2-Fluorophenol	1.122	1.104	1.6	98	0.00
6		Aniline	2.036	1.984	2.6	98	0.00
7	S	Phenol-d6	1.579	1.561	1.1	101	0.00
8		2-Chlorophenol	1.258	1.235	1.8	97	0.00
9		Benzaldehyde	0.954	0.861	9.7	93	0.00
10	C	Phenol	1.592	1.588	0.3	101	0.00
11		bis(2-Chloroethyl)ether	1.262	1.177	6.7	96	0.00
12		1,3-Dichlorobenzene	1.627	1.592	2.2	98	0.00
13	C	1,4-Dichlorobenzene	1.641	1.628	0.8	101	0.00
14		1,2-Dichlorobenzene	1.571	1.549	1.4	99	0.00
15		Benzyl Alcohol	1.258	1.255	0.2	99	0.00
16		2,2'-oxybis(1-Chloropropane	1.772	1.562	11.9	89	0.00
17		2-Methylphenol	1.159	1.090	6.0	95	0.00
18		Hexachloroethane	0.647	0.621	4.0	97	0.00
19	P	n-Nitroso-di-n-propylamine	1.178	1.111	5.7	97	0.00
20		3+4-Methylphenols	1.544	1.517	1.7	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
22		Acetophenone	0.600	0.571	4.8	96	0.00
23	S	Nitrobenzene-d5	0.480	0.467	2.7	97	0.00
24		Nitrobenzene	0.481	0.465	3.3	99	0.00
25		Isophorone	0.861	0.808	6.2	95	0.00
26	C	2-Nitrophenol	0.201	0.197	2.0	100	0.00
27		2,4-Dimethylphenol	0.280	0.266	5.0	92	0.00
28		bis(2-Chloroethoxy)methane	0.455	0.424	6.8	95	0.00
29	C	2,4-Dichlorophenol	0.378	0.378	0.0	99	0.00
30		1,2,4-Trichlorobenzene	0.484	0.481	0.6	101	0.00
31		Naphthalene	1.106	1.078	2.5	100	0.00
32		Benzoic acid	0.154	0.158	-2.6	112	0.00
33		4-Chloroaniline	0.466	0.463	0.6	99	0.00
34	C	Hexachlorobutadiene	0.384	0.393	-2.3	104	0.00
35		Caprolactam	0.118	0.111	5.9	96	0.00
36	C	4-Chloro-3-methylphenol	0.409	0.395	3.4	100	0.00
37		2-Methylnaphthalene	0.817	0.787	3.7	97	0.00
38		1-Methylnaphthalene	0.780	0.749	4.0	98	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
40		1,2,4,5-Tetrachlorobenzene	0.881	0.880	0.1	100	0.00
41	P	Hexachlorocyclopentadiene	0.436	0.395	9.4	86	0.00
42	S	2,4,6-Tribromophenol	0.341	0.344	-0.9	101	0.00
43	C	2,4,6-Trichlorophenol	0.523	0.517	1.1	99	0.00
44		2,4,5-Trichlorophenol	0.521	0.522	-0.2	104	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070219\
Data File : BG041565.D
Acq On : 2 Jul 2019 9:22
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 03 07:19:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG062619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 27 13:25:55 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	1.559	1.564	-0.3	100	0.00
46		1,1'-Biphenyl	1.579	1.540	2.5	98	0.00
47		2-Chloronaphthalene	1.355	1.341	1.0	99	0.00
48		2-Nitroaniline	0.459	0.441	3.9	97	0.00
49		Acenaphthylene	2.026	1.978	2.4	99	0.00
50		Dimethylphthalate	1.820	1.767	2.9	99	0.00
51		2,6-Dinitrotoluene	0.386	0.386	0.0	102	0.00
52	C	Acenaphthene	1.201	1.167	2.8	98	0.00
53		3-Nitroaniline	0.369	0.336	8.9	90	0.00
54	P	2,4-Dinitrophenol	0.208	0.202	2.9	101	0.00
55		Dibenzofuran	2.092	2.063	1.4	99	0.00
56	P	4-Nitrophenol	0.238	0.221	7.1	90	0.00
57		2,4-Dinitrotoluene	0.560	0.543	3.0	97	0.00
58		Fluorene	1.664	1.651	0.8	100	0.00
59		2,3,4,6-Tetrachlorophenol	0.549	0.545	0.7	98	0.00
60		Diethylphthalate	1.863	1.798	3.5	97	0.00
61		4-Chlorophenyl-phenylether	1.073	1.071	0.2	101	0.00
62		4-Nitroaniline	0.359	0.348	3.1	96	0.00
63		Azobenzene	1.571	1.514	3.6	98	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
65		4,6-Dinitro-2-methylphenol	0.124	0.121	2.4	99	0.00
66	c	n-Nitrosodiphenylamine	0.550	0.526	4.4	98	0.00
67		4-Bromophenyl-phenylether	0.273	0.268	1.8	102	0.00
68		Hexachlorobenzene	0.294	0.290	1.4	101	0.00
69		Atrazine	0.233	0.209	10.3	95	0.00
70	C	Pentachlorophenol	0.133	0.118	11.3	89	0.00
71		Phenanthrene	1.102	1.084	1.6	100	0.00
72		Anthracene	1.103	1.073	2.7	99	0.00
73		Carbazole	0.937	0.910	2.9	98	0.00
74		Di-n-butylphthalate	1.178	1.127	4.3	96	0.00
75	C	Fluoranthene	1.469	1.442	1.8	98	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
77		Benzidine	0.444	0.423	4.7	84	0.00
78		Pyrene	1.365	1.336	2.1	98	0.00
79	S	Terphenyl-d14	0.941	0.918	2.4	97	0.00
80		Butylbenzylphthalate	0.510	0.474	7.1	93	0.00
81		Benzo(a)anthracene	1.390	1.336	3.9	96	0.00
82		3,3'-Dichlorobenzidine	0.481	0.475	1.2	98	0.00
83		Chrysene	1.346	1.291	4.1	97	0.00
84		Bis(2-ethylhexyl)phthalate	0.723	0.663	8.3	93	0.00
85	c	Di-n-octyl phthalate	1.223	1.122	8.3	92	0.00
86		Indeno(1,2,3-cd)pyrene	1.630	1.582	2.9	98	0.00
87	I	Perylene-d12	1.000	1.000	0.0	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070219\
Data File : BG041565.D
Acq On : 2 Jul 2019 9:22
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 03 07:19:26 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG062619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 27 13:25:55 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	1.276	1.201	5.9	94	0.00
89	Benzo(k)fluoranthene	1.252	1.229	1.8	100	0.00
90 C	Benzo(a)pyrene	1.212	1.149	5.2	97	0.00
91	Dibenzo(a,h)anthracene	1.214	1.161	4.4	97	0.00
92	Benzo(g,h,i)perylene	1.207	1.152	4.6	97	0.00

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