Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082819\
Data File : BG042525.D
Acq On : 28 Aug 2019 8:44
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 28 15:16:52 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 22 14:29:15 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	78	0.00
2		1,4-Dioxane	40.000	43.229	-8.1	83	0.00
3		Pyridine	40.000	41.313	-3.3	77	0.00
4		n-Nitrosodimethylamine	40.000	40.745	-1.9	80	0.00
5	S	2-Fluorophenol	80.000	82.546	-3.2	79	0.00
6		Aniline	40.000	39.749	0.6	77	0.00
7	S	Phenol-d6	80.000	79.640	0.4	76	0.00
8		2-Chlorophenol	40.000	41.047	-2.6	79	0.00
9		Benzaldehyde	40.000	40.642	-1.6	94	0.00
10	C	Phenol	40.000	40.340	-0.9	77	0.00
11		bis(2-Chloroethyl)ether	40.000	41.000	-2.5	79	0.00
12		1,3-Dichlorobenzene	40.000	41.339	-3.3	80	0.00
13	C	1,4-Dichlorobenzene	40.000	40.673	-1.7	79	0.00
14		1,2-Dichlorobenzene	40.000	40.679	-1.7	79	0.00
15		Benzyl Alcohol	40.000	40.503	-1.3	78	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.471	1.3	77	-0.01
17		2-Methylphenol	40.000	40.354	-0.9	79	0.00
18		Hexachloroethane	40.000	40.588	-1.5	79	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.805	-2.0	79	0.00
20		3+4-Methylphenols	40.000	39.912	0.2	78	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	79	0.00
22		Acetophenone	40.000	41.289	-3.2	80	0.00
23	S	Nitrobenzene-d5	80.000	82.255	-2.8	80	0.00
24		Nitrobenzene	40.000	40.567	-1.4	79	0.00
25		Isophorone	40.000	40.286	-0.7	77	-0.01
26	C	2-Nitrophenol	40.000	41.114	-2.8	81	0.00
27		2,4-Dimethylphenol	40.000	39.619	1.0	76	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.589	-1.5	78	0.00
29	C	2,4-Dichlorophenol	40.000	40.875	-2.2	78	0.00
30		1,2,4-Trichlorobenzene	40.000	41.621	-4.1	81	0.00
31		Naphthalene	40.000	40.605	-1.5	78	0.00
32		Benzoic acid	40.000	35.825	10.4	74	0.00
33		4-Chloroaniline	40.000	40.527	-1.3	77	0.00
34	C	Hexachlorobutadiene	40.000	42.266	-5.7	83	0.00
35		Caprolactam	40.000	38.371	4.1	72	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.915	-2.3	78	0.00
37		2-Methylnaphthalene	40.000	40.753	-1.9	77	0.00
38		1-Methylnaphthalene	40.000	41.027	-2.6	78	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	77	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	42.560	-6.4	81	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.902	2.7	77	0.00
42	S	2,4,6-Tribromophenol	80.000	81.799	-2.2	75	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.238	-3.1	78	0.00
44		2,4,5-Trichlorophenol	40.000	41.876	-4.7	80	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082819\
Data File : BG042525.D
Acq On : 28 Aug 2019 8:44
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 28 15:16:52 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 22 14:29:15 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	Amount	Calc.	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	80.000	86.498	-8.1	80	0.00
46		1,1'-Biphenyl	40.000	41.352	-3.4	78	0.00
47		2-Chloronaphthalene	40.000	41.297	-3.2	78	0.00
48		2-Nitroaniline	40.000	40.367	-0.9	76	0.00
49		Acenaphthylene	40.000	41.604	-4.0	77	0.00
50		Dimethylphthalate	40.000	41.675	-4.2	76	0.00
51		2,6-Dinitrotoluene	40.000	41.190	-3.0	76	0.00
52	C	Acenaphthene	40.000	40.654	-1.6	76	0.00
53		3-Nitroaniline	40.000	40.031	-0.1	74	0.00
54	P	2,4-Dinitrophenol	40.000	35.034	12.4	67	0.00
55		Dibenzofuran	40.000	42.103	-5.3	77	0.00
56	P	4-Nitrophenol	40.000	39.333	1.7	73	0.00
57		2,4-Dinitrotoluene	40.000	40.887	-2.2	75	0.00
58		Fluorene	40.000	41.700	-4.3	76	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.413	-3.5	77	0.00
60		Diethylphthalate	40.000	41.003	-2.5	74	0.00
61		4-Chlorophenyl-phenylether	40.000	41.404	-3.5	77	0.00
62		4-Nitroaniline	40.000	40.335	-0.8	73	0.00
63		Azobenzene	40.000	39.923	0.2	74	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	76	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.979	0.1	73	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.003	-2.5	75	0.00
67		4-Bromophenyl-phenylether	40.000	41.249	-3.1	77	0.00
68		Hexachlorobenzene	40.000	40.940	-2.3	76	0.00
69		Atrazine	40.000	36.365	9.1	70	0.00
70	C	Pentachlorophenol	40.000	37.739	5.7	69	0.00
71		Phenanthrene	40.000	42.115	-5.3	76	0.00
72		Anthracene	40.000	42.114	-5.3	75	0.00
73		Carbazole	40.000	41.651	-4.1	75	0.00
74		Di-n-butylphthalate	40.000	42.536	-6.3	75	0.00
75	C	Fluoranthene	40.000	44.249	-10.6	78	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	78	0.00
77		Benzidine	40.000	34.735	13.2	77	0.00
78		Pyrene	40.000	42.390	-6.0	78	0.00
79	S	Terphenyl-d14	80.000	84.299	-5.4	79	0.00
80		Butylbenzylphthalate	40.000	41.028	-2.6	77	0.00
81		Benzo(a)anthracene	40.000	42.311	-5.8	78	0.00
82		3,3'-Dichlorobenzidine	40.000	40.325	-0.8	76	0.00
83		Chrysene	40.000	41.993	-5.0	77	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.155	-2.9	76	0.00
85	c	Di-n-octyl phthalate	40.000	41.091	-2.7	76	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	40.027	-0.1	76	-0.02
87	I	Perylene-d12	20.000	20.000	0.0	76	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082819\
Data File : BG042525.D
Acq On : 28 Aug 2019 8:44
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 28 15:16:52 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 22 14:29:15 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	40.000	41.336	-3.3	74	0.00
89	Benzo(k)fluoranthene	40.000	42.675	-6.7	78	0.00
90 C	Benzo(a)pyrene	40.000	42.269	-5.7	77	0.00
91	Dibenzo(a,h)anthracene	40.000	40.970	-2.4	76	-0.01
92	Benzo(g,h,i)perylene	40.000	40.904	-2.3	76	-0.01

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