Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
Data File : BG041199.D
Acq On : 12 Jun 2019 8:48
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 13 07:30:41 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 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	73	0.00
2		1,4-Dioxane	40.000	38.003	5.0	70	0.00
3		Pyridine	40.000	39.976	0.1	73	0.00
4		n-Nitrosodimethylamine	40.000	42.138	-5.3	76	0.00
5	S	2-Fluorophenol	80.000	81.259	-1.6	74	0.00
6		Aniline	40.000	39.637	0.9	73	0.00
7	S	Phenol-d6	80.000	80.270	-0.3	74	0.00
8		2-Chlorophenol	40.000	40.583	-1.5	75	0.00
9		Benzaldehyde	40.000	40.606	-1.5	74	0.00
10	C	Phenol	40.000	40.502	-1.3	75	0.00
11		bis(2-Chloroethyl)ether	40.000	40.372	-0.9	73	0.00
12		1,3-Dichlorobenzene	40.000	40.906	-2.3	74	0.00
13	C	1,4-Dichlorobenzene	40.000	41.113	-2.8	74	0.00
14		1,2-Dichlorobenzene	40.000	40.470	-1.2	74	0.00
15		Benzyl Alcohol	40.000	40.537	-1.3	75	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.542	3.6	71	0.00
17		2-Methylphenol	40.000	39.253	1.9	72	0.00
18		Hexachloroethane	40.000	40.562	-1.4	75	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.970	2.6	71	0.00
20		3+4-Methylphenols	40.000	40.005	-0.0	74	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	76	0.00
22		Acetophenone	40.000	39.924	0.2	74	0.00
23	S	Nitrobenzene-d5	80.000	81.105	-1.4	76	0.00
24		Nitrobenzene	40.000	41.123	-2.8	76	0.00
25		Isophorone	40.000	38.853	2.9	72	0.00
26	C	2-Nitrophenol	40.000	43.281	-8.2	80	0.00
27		2,4-Dimethylphenol	40.000	39.074	2.3	73	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.270	4.3	71	0.00
29	C	2,4-Dichlorophenol	40.000	39.581	1.0	73	0.00
30		1,2,4-Trichlorobenzene	40.000	42.053	-5.1	78	0.00
31		Naphthalene	40.000	40.352	-0.9	75	0.00
32		Benzoic acid	40.000	35.540	11.2	66	0.00
33		4-Chloroaniline	40.000	39.120	2.2	73	0.00
34	C	Hexachlorobutadiene	40.000	41.153	-2.9	79	0.00
35		Caprolactam	40.000	37.658	5.9	69	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.559	3.6	72	0.00
37		2-Methylnaphthalene	40.000	39.325	1.7	73	0.00
38		1-Methylnaphthalene	40.000	39.727	0.7	74	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	75	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.603	-4.0	76	0.00
41	P	Hexachlorocyclopentadiene	40.000	41.194	-3.0	81	0.00
42	S	2,4,6-Tribromophenol	80.000	77.322	3.3	70	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.672	-1.7	73	0.00
44		2,4,5-Trichlorophenol	40.000	38.774	3.1	71	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
Data File : BG041199.D
Acq On : 12 Jun 2019 8:48
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 13 07:30:41 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 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	83.327	-4.2	75	0.00
46		1,1'-Biphenyl	40.000	41.190	-3.0	74	0.00
47		2-Chloronaphthalene	40.000	41.854	-4.6	76	0.00
48		2-Nitroaniline	40.000	41.648	-4.1	75	0.00
49		Acenaphthylene	40.000	40.047	-0.1	71	0.00
50		Dimethylphthalate	40.000	39.353	1.6	70	0.00
51		2,6-Dinitrotoluene	40.000	38.952	2.6	71	0.00
52	C	Acenaphthene	40.000	40.004	-0.0	72	0.00
53		3-Nitroaniline	40.000	40.106	-0.3	71	0.00
54	P	2,4-Dinitrophenol	40.000	39.359	1.6	74	0.00
55		Dibenzofuran	40.000	40.357	-0.9	72	0.00
56	P	4-Nitrophenol	40.000	34.418	14.0	61	0.00
57		2,4-Dinitrotoluene	40.000	39.581	1.0	70	0.00
58		Fluorene	40.000	39.498	1.3	71	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.021	2.4	70	0.00
60		Diethylphthalate	40.000	39.215	2.0	70	0.00
61		4-Chlorophenyl-phenylether	40.000	40.097	-0.2	73	0.00
62		4-Nitroaniline	40.000	38.706	3.2	70	0.00
63		Azobenzene	40.000	38.984	2.5	69	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	71	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.042	-2.6	77	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.759	0.6	69	0.00
67		4-Bromophenyl-phenylether	40.000	39.702	0.7	68	0.00
68		Hexachlorobenzene	40.000	39.837	0.4	69	0.00
69		Atrazine	40.000	39.585	1.0	63	0.00
70	C	Pentachlorophenol	40.000	37.704	5.7	65	0.00
71		Phenanthrene	40.000	41.255	-3.1	70	0.00
72		Anthracene	40.000	41.530	-3.8	71	0.00
73		Carbazole	40.000	41.814	-4.5	71	0.00
74		Di-n-butylphthalate	40.000	41.347	-3.4	69	0.00
75	C	Fluoranthene	40.000	42.335	-5.8	72	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	70	0.00
77		Benzidine	40.000	32.245	19.4	53	0.00
78		Pyrene	40.000	42.611	-6.5	72	0.00
79	S	Terphenyl-d14	80.000	88.130	-10.2	73	0.00
80		Butylbenzylphthalate	40.000	41.717	-4.3	70	0.00
81		Benzo(a)anthracene	40.000	40.863	-2.2	70	0.00
82		3,3'-Dichlorobenzidine	40.000	39.615	1.0	69	0.00
83		Chrysene	40.000	41.746	-4.4	71	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.788	-4.5	71	0.00
85	c	Di-n-octyl phthalate	40.000	41.594	-4.0	70	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	32.345	19.1	57	0.00
87	I	Perylene-d12	20.000	20.000	0.0	67	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061219\
Data File : BG041199.D
Acq On : 12 Jun 2019 8:48
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 13 07:30:41 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Jun 09 23:08:34 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	39.701	0.7	65	0.00
89	Benzo(k)fluoranthene	40.000	42.720	-6.8	69	0.00
90 C	Benzo(a)pyrene	40.000	40.552	-1.4	66	0.00
91	Dibenzo(a,h)anthracene	40.000	32.205	19.5	52	0.00
92	Benzo(g,h,i)perylene	40.000	32.357	19.1	53	0.00

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