Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG110419\
Data File : BG043498.D
Acq On : 5 Nov 2019 2:50
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 05 04:54:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 04 15:22:15 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	135	0.00
2		1,4-Dioxane	40.000	37.846	5.4	122	0.00
3		Pyridine	40.000	39.796	0.5	115	0.00
4		n-Nitrosodimethylamine	40.000	37.653	5.9	122	0.00
5	S	2-Fluorophenol	80.000	79.512	0.6	129	0.00
6		Aniline	40.000	37.575	6.1	119	0.00
7	S	Phenol-d6	80.000	77.245	3.4	125	0.00
8		2-Chlorophenol	40.000	38.994	2.5	126	0.00
9		Benzaldehyde	40.000	41.496	-3.7	139	0.00
10	C	Phenol	40.000	38.125	4.7	121	0.00
11		bis(2-Chloroethyl)ether	40.000	36.876	7.8	118	0.00
12		1,3-Dichlorobenzene	40.000	39.741	0.6	130	0.00
13	C	1,4-Dichlorobenzene	40.000	38.645	3.4	125	0.00
14		1,2-Dichlorobenzene	40.000	37.986	5.0	123	0.00
15		Benzyl Alcohol	40.000	38.645	3.4	121	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	35.213	12.0	111	0.00
17		2-Methylphenol	40.000	37.886	5.3	125	0.00
18		Hexachloroethane	40.000	38.078	4.8	124	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	35.002	12.5	114	0.00
20		3+4-Methylphenols	40.000	36.925	7.7	120	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	124	0.00
22		Acetophenone	40.000	39.290	1.8	116	0.00
23	S	Nitrobenzene-d5	80.000	77.845	2.7	117	0.00
24		Nitrobenzene	40.000	39.397	1.5	117	0.00
25		Isophorone	40.000	36.679	8.3	109	0.00
26	C	2-Nitrophenol	40.000	40.562	-1.4	124	0.00
27		2,4-Dimethylphenol	40.000	40.188	-0.5	120	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.436	6.4	108	0.00
29	C	2,4-Dichlorophenol	40.000	41.616	-4.0	122	0.00
30		1,2,4-Trichlorobenzene	40.000	40.879	-2.2	123	0.00
31		Naphthalene	40.000	39.727	0.7	120	0.00
32		Benzoic acid	40.000	37.708	5.7	128	0.00
33		4-Chloroaniline	40.000	40.048	-0.1	116	0.00
34	C	Hexachlorobutadiene	40.000	40.131	-0.3	122	0.00
35		Caprolactam	40.000	38.700	3.2	112	0.02
36	C	4-Chloro-3-methylphenol	40.000	38.889	2.8	115	0.00
37		2-Methylnaphthalene	40.000	38.588	3.5	115	0.00
38		1-Methylnaphthalene	40.000	38.651	3.4	115	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	118	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.040	-2.6	116	0.00
41	P	Hexachlorocyclopentadiene	40.000	13.494	66.3#	38	0.00
42	S	2,4,6-Tribromophenol	80.000	76.209	4.7	106	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.071	-0.2	111	0.00
44		2,4,5-Trichlorophenol	40.000	39.137	2.2	109	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG110419\
Data File : BG043498.D
Acq On : 5 Nov 2019 2:50
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 05 04:54:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 04 15:22:15 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	79.078	1.2	111	0.00
46		1,1'-Biphenyl	40.000	39.747	0.6	111	0.00
47		2-Chloronaphthalene	40.000	40.226	-0.6	114	0.00
48		2-Nitroaniline	40.000	39.153	2.1	107	0.00
49		Acenaphthylene	40.000	40.055	-0.1	112	0.00
50		Dimethylphthalate	40.000	37.062	7.3	105	0.00
51		2,6-Dinitrotoluene	40.000	38.812	3.0	108	0.00
52	C	Acenaphthene	40.000	39.228	1.9	111	0.00
53		3-Nitroaniline	40.000	39.542	1.1	110	0.00
54	P	2,4-Dinitrophenol	40.000	19.951	50.1#	46	0.00
55		Dibenzofuran	40.000	38.840	2.9	109	0.00
56	P	4-Nitrophenol	40.000	37.059	7.4	102	0.00
57		2,4-Dinitrotoluene	40.000	38.438	3.9	109	0.00
58		Fluorene	40.000	38.514	3.7	107	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	37.746	5.6	101	0.00
60		Diethylphthalate	40.000	36.672	8.3	103	0.00
61		4-Chlorophenyl-phenylether	40.000	38.127	4.7	108	0.00
62		4-Nitroaniline	40.000	41.505	-3.8	114	0.00
63		Azobenzene	40.000	37.512	6.2	105	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	117	0.00
65		4,6-Dinitro-2-methylphenol	40.000	19.561	51.1#	56	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.250	4.4	106	0.00
67		4-Bromophenyl-phenylether	40.000	38.105	4.7	107	0.00
68		Hexachlorobenzene	40.000	37.898	5.3	107	0.00
69		Atrazine	40.000	35.441	11.4	101	0.00
70	C	Pentachlorophenol	40.000	34.882	12.8	94	0.00
71		Phenanthrene	40.000	38.168	4.6	106	0.00
72		Anthracene	40.000	38.477	3.8	105	0.00
73		Carbazole	40.000	38.534	3.7	107	0.00
74		Di-n-butylphthalate	40.000	38.493	3.8	106	0.00
75	C	Fluoranthene	40.000	37.163	7.1	101	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	112	0.00
77		Benzidine	40.000	42.568	-6.4	106	0.00
78		Pyrene	40.000	39.126	2.2	104	0.00
79	S	Terphenyl-d14	80.000	78.029	2.5	103	0.00
80		Butylbenzylphthalate	40.000	40.803	-2.0	109	0.00
81		Benzo(a)anthracene	40.000	40.273	-0.7	109	0.00
82		3,3'-Dichlorobenzidine	40.000	42.575	-6.4	113	0.00
83		Chrysene	40.000	39.864	0.3	106	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.832	-2.1	110	0.00
85	c	Di-n-octyl phthalate	40.000	40.964	-2.4	111	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	38.808	3.0	105	0.00
87	I	Perylene-d12	20.000	20.000	0.0	117	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG110419\
Data File : BG043498.D
Acq On : 5 Nov 2019 2:50
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 05 04:54:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 04 15:22:15 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	39.719	0.7	111	0.00
89	Benzo(k)fluoranthene	40.000	38.503	3.7	107	0.00
90 C	Benzo(a)pyrene	40.000	39.660	0.9	111	0.00
91	Dibenzo(a,h)anthracene	40.000	37.626	5.9	105	0.00
92	Benzo(g,h,i)perylene	40.000	33.752	15.6	94	0.02

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