Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030520\
Data File : BF119209.D
Acq On : 5 Mar 2020 12:10
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 06 06:22:00 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 2020
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	63	-0.02
2		1,4-Dioxane	40.000	36.656	8.4	61	-0.05
3		Pyridine	40.000	37.910	5.2	63	-0.05
4		n-Nitrosodimethylamine	40.000	40.987	-2.5	68	-0.07
5	S	2-Fluorophenol	80.000	82.814	-3.5	67	-0.01
6		Aniline	40.000	39.637	0.9	64	-0.02
7	S	Phenol-d6	80.000	80.837	-1.0	66	-0.02
8		2-Chlorophenol	40.000	41.288	-3.2	67	-0.02
9		Benzaldehyde	40.000	35.027	12.4	57	-0.02
10	C	Phenol	40.000	41.609	-4.0	68	-0.02
11		bis(2-Chloroethyl)ether	40.000	40.272	-0.7	66	-0.03
12		1,3-Dichlorobenzene	40.000	40.121	-0.3	66	-0.02
13	C	1,4-Dichlorobenzene	40.000	40.455	-1.1	66	-0.02
14		1,2-Dichlorobenzene	40.000	41.145	-2.9	67	-0.02
15		Benzyl Alcohol	40.000	43.076	-7.7	69	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	39.655	0.9	66	-0.02
17		2-Methylphenol	40.000	40.225	-0.6	66	-0.01
18		Hexachloroethane	40.000	40.945	-2.4	68	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	43.651	-9.1	73	-0.03
20		3+4-Methylphenols	40.000	44.160	-10.4	74	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	66	-0.02
22		Acetophenone	40.000	41.342	-3.4	72	-0.02
23	S	Nitrobenzene-d5	80.000	80.052	-0.1	70	-0.02
24		Nitrobenzene	40.000	39.507	1.2	69	-0.02
25		Isophorone	40.000	39.085	2.3	69	-0.02
26	C	2-Nitrophenol	40.000	39.304	1.7	69	-0.02
27		2,4-Dimethylphenol	40.000	39.431	1.4	69	-0.02
28		bis(2-Chloroethoxy)methane	40.000	38.578	3.6	68	-0.02
29	C	2,4-Dichlorophenol	40.000	38.869	2.8	67	-0.02
30		1,2,4-Trichlorobenzene	40.000	38.223	4.4	66	-0.02
31		Naphthalene	40.000	39.564	1.1	68	-0.02
32		Benzoic acid	40.000	40.045	-0.1	74	-0.02
33		4-Chloroaniline	40.000	39.257	1.9	68	-0.02
34	C	Hexachlorobutadiene	40.000	38.876	2.8	68	-0.02
35		Caprolactam	40.000	40.529	-1.3	70	-0.04
36	C	4-Chloro-3-methylphenol	40.000	40.418	-1.0	70	-0.01
37		2-Methylnaphthalene	40.000	39.759	0.6	69	-0.02
38		1-Methylnaphthalene	40.000	39.918	0.2	69	-0.02
39	I	Acenaphthene-d10	20.000	20.000	0.0	67	-0.02
40		1,2,4,5-Tetrachlorobenzene	40.000	38.436	3.9	69	-0.02
41	P	Hexachlorocyclopentadiene	40.000	32.625	18.4	59	-0.02
42	S	2,4,6-Tribromophenol	80.000	78.157	2.3	70	-0.02
43	C	2,4,6-Trichlorophenol	40.000	37.959	5.1	68	-0.01
44		2,4,5-Trichlorophenol	40.000	37.475	6.3	67	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030520\
Data File : BF119209.D
Acq On : 5 Mar 2020 12:10
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 06 06:22:00 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 2020
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	78.638	1.7	72	-0.02
46		1,1'-Biphenyl	40.000	39.263	1.8	69	-0.02
47		2-Chloronaphthalene	40.000	39.519	1.2	71	-0.02
48		2-Nitroaniline	40.000	41.815	-4.5	74	-0.02
49		Acenaphthylene	40.000	39.460	1.3	70	-0.02
50		Dimethylphthalate	40.000	39.301	1.7	71	-0.02
51		2,6-Dinitrotoluene	40.000	39.090	2.3	70	-0.02
52	C	Acenaphthene	40.000	39.492	1.3	70	-0.02
53		3-Nitroaniline	40.000	39.641	0.9	70	-0.02
54	P	2,4-Dinitrophenol	40.000	37.695	5.8	69	-0.01
55		Dibenzofuran	40.000	39.198	2.0	68	-0.02
56	P	4-Nitrophenol	40.000	39.184	2.0	68	0.00
57		2,4-Dinitrotoluene	40.000	38.914	2.7	68	-0.02
58		Fluorene	40.000	41.862	-4.7	74	-0.02
59		2,3,4,6-Tetrachlorophenol	40.000	38.012	5.0	69	-0.02
60		Diethylphthalate	40.000	40.564	-1.4	72	-0.02
61		4-Chlorophenyl-phenylether	40.000	40.736	-1.8	72	-0.02
62		4-Nitroaniline	40.000	40.585	-1.5	72	-0.02
63		Azobenzene	40.000	40.661	-1.7	73	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	70	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	39.183	2.0	73	-0.01
66	c	n-Nitrosodiphenylamine	40.000	38.651	3.4	71	-0.02
67		4-Bromophenyl-phenylether	40.000	37.874	5.3	71	-0.02
68		Hexachlorobenzene	40.000	37.983	5.0	70	-0.02
69		Atrazine	40.000	35.504	11.2	66	-0.02
70	C	Pentachlorophenol	40.000	33.063	17.3	62	-0.01
71		Phenanthrene	40.000	39.337	1.7	72	-0.02
72		Anthracene	40.000	39.486	1.3	72	-0.02
73		Carbazole	40.000	39.778	0.6	73	-0.01
74		Di-n-butylphthalate	40.000	40.944	-2.4	75	-0.02
75	C	Fluoranthene	40.000	40.480	-1.2	74	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	72	-0.01
77		Benzidine	40.000	34.646	13.4	65	-0.02
78		Pyrene	40.000	36.473	8.8	72	-0.02
79	S	Terphenyl-d14	80.000	76.975	3.8	77	-0.02
80		Butylbenzylphthalate	40.000	38.924	2.7	76	-0.02
81		Benzo(a)anthracene	40.000	40.049	-0.1	77	-0.01
82		3,3'-Dichlorobenzidine	40.000	40.043	-0.1	74	-0.02
83		Chrysene	40.000	38.751	3.1	75	-0.02
84		Bis(2-ethylhexyl)phthalate	40.000	41.408	-3.5	80	-0.02
85	c	Di-n-octyl phthalate	40.000	41.693	-4.2	76	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	35.662	10.8	53	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	51	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030520\
Data File : BF119209.D
Acq On : 5 Mar 2020 12:10
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 06 06:22:00 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 2020
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	38.386	4.0	74	-0.01
89	Benzo(k)fluoranthene	40.000	40.992	-2.5	72	-0.01
90 C	Benzo(a)pyrene	40.000	38.643	3.4	53	-0.01
91	Dibenzo(a,h)anthracene	40.000	36.865	7.8	54	-0.02
92	Benzo(g,h,i)perylene	40.000	34.751	13.1	51	-0.01

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