Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110420\
Data File : BP003799.D
Acq On : 04 Nov 2020 09:32
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Nov 05 14:51:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 05 14:31:45 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	100	0.01
2		1,4-Dioxane	40.000	37.255	6.9	99	0.02
3		Pyridine	40.000	37.596	6.0	98	0.02
4		n-Nitrosodimethylamine	40.000	37.077	7.3	97	0.02
5	S	2-Fluorophenol	80.000	76.351	4.6	100	0.00
6		Aniline	40.000	37.995	5.0	98	0.01
7	S	Phenol-d6	80.000	75.064	6.2	97	0.00
8		2-Chlorophenol	40.000	37.764	5.6	98	0.01
9		Benzaldehyde	40.000	36.753	8.1	100	0.01
10	C	Phenol	40.000	37.651	5.9	99	0.00
11		bis(2-Chloroethyl)ether	40.000	37.362	6.6	98	0.02
12		1,3-Dichlorobenzene	40.000	37.070	7.3	98	0.02
13	C	1,4-Dichlorobenzene	40.000	37.248	6.9	99	0.02
14		1,2-Dichlorobenzene	40.000	37.006	7.5	97	0.01
15		Benzyl Alcohol	40.000	38.115	4.7	96	0.01
16		2,2'-oxybis(1-Chloropropane	40.000	36.718	8.2	97	0.01
17		2-Methylphenol	40.000	37.518	6.2	97	0.00
18		Hexachloroethane	40.000	38.280	4.3	98	0.02
19	P	n-Nitroso-di-n-propylamine	40.000	37.944	5.1	96	0.01
20		3+4-Methylphenols	40.000	37.942	5.1	97	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.01
22		Acetophenone	40.000	36.951	7.6	97	0.01
23	S	Nitrobenzene-d5	80.000	75.459	5.7	98	0.01
24		Nitrobenzene	40.000	37.307	6.7	98	0.01
25		Isophorone	40.000	37.972	5.1	97	0.00
26	C	2-Nitrophenol	40.000	39.688	0.8	98	0.00
27		2,4-Dimethylphenol	40.000	37.777	5.6	99	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.047	7.4	97	0.01
29	C	2,4-Dichlorophenol	40.000	37.971	5.1	98	0.00
30		1,2,4-Trichlorobenzene	40.000	36.959	7.6	98	0.01
31		Naphthalene	40.000	36.651	8.4	98	0.01
32		Benzoic acid	40.000	36.426	8.9	97	-0.01
33		4-Chloroaniline	40.000	37.767	5.6	98	0.00
34	C	Hexachlorobutadiene	40.000	37.166	7.1	99	0.02
35		Caprolactam	40.000	40.233	-0.6	100	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.226	4.4	99	0.00
37		2-Methylnaphthalene	40.000	37.075	7.3	99	0.00
38		1-Methylnaphthalene	40.000	36.834	7.9	99	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	98	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	37.084	7.3	98	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.513	3.7	98	0.01
42	S	2,4,6-Tribromophenol	80.000	78.269	2.2	100	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.911	2.7	100	0.00
44		2,4,5-Trichlorophenol	40.000	38.272	4.3	99	0.00
45	S	2-Fluorobiphenyl	80.000	73.444	8.2	98	0.00
46		1,1'-Biphenyl	40.000	36.719	8.2	99	0.00
47		2-Chloronaphthalene	40.000	36.941	7.6	98	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110420\
Data File : BP003799.D
Acq On : 04 Nov 2020 09:32
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Nov 05 14:51:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 05 14:31:45 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)
48		2-Nitroaniline	40.000	39.744	0.6	99	0.00
49		Acenaphthylene	40.000	37.346	6.6	99	0.00
50		Dimethylphthalate	40.000	37.513	6.2	100	0.00
51		2,6-Dinitrotoluene	40.000	38.909	2.7	101	0.00
52	C	Acenaphthene	40.000	37.117	7.2	99	0.00
53		3-Nitroaniline	40.000	38.952	2.6	101	0.00
54	P	2,4-Dinitrophenol	40.000	39.216	2.0	100	-0.02
55		Dibenzofuran	40.000	36.983	7.5	99	0.00
56	P	4-Nitrophenol	40.000	40.210	-0.5	101	-0.02
57		2,4-Dinitrotoluene	40.000	39.662	0.8	101	-0.01
58		Fluorene	40.000	36.990	7.5	100	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.826	2.9	100	0.00
60		Diethylphthalate	40.000	37.794	5.5	100	0.00
61		4-Chlorophenyl-phenylether	40.000	37.318	6.7	101	0.00
62		4-Nitroaniline	40.000	39.510	1.2	102	0.00
63		Azobenzene	40.000	37.572	6.1	99	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	101	0.00
65		4,6-Dinitro-2-methylphenol	40.000	36.969	7.6	101	-0.01
66	c	n-Nitrosodiphenylamine	40.000	36.889	7.8	100	0.00
67		4-Bromophenyl-phenylether	40.000	37.348	6.6	101	0.00
68		Hexachlorobenzene	40.000	36.688	8.3	101	0.00
69		Atrazine	40.000	38.390	4.0	101	0.00
70	C	Pentachlorophenol	40.000	39.077	2.3	101	-0.01
71		Phenanthrene	40.000	36.728	8.2	101	0.00
72		Anthracene	40.000	36.862	7.8	101	0.00
73		Carbazole	40.000	37.049	7.4	101	-0.01
74		Di-n-butylphthalate	40.000	39.177	2.1	102	0.00
75	C	Fluoranthene	40.000	37.348	6.6	101	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	101	-0.02
77		Benzidine	40.000	41.706	-4.3	104	-0.02
78		Pyrene	40.000	37.627	5.9	101	-0.02
79	S	Terphenyl-d14	80.000	74.328	7.1	102	-0.01
80		Butylbenzylphthalate	40.000	40.379	-0.9	102	-0.01
81		Benzo(a)anthracene	40.000	37.451	6.4	102	-0.02
82		3,3'-Dichlorobenzidine	40.000	38.892	2.8	104	-0.02
83		Chrysene	40.000	36.905	7.7	101	-0.03
84		Bis(2-ethylhexyl)phthalate	40.000	39.341	1.6	102	-0.02
85	c	Di-n-octyl phthalate	40.000	39.559	1.1	102	-0.03
86	I	Perylene-d12	20.000	20.000	0.0	102	-0.05
87		Indeno(1,2,3-cd)pyrene	40.000	36.161	9.6	102	-0.07
88		Benzo(b)fluoranthene	40.000	36.382	9.0	105	-0.04
89		Benzo(k)fluoranthene	40.000	35.617	11.0	100	-0.04
90	C	Benzo(a)pyrene	40.000	36.122	9.7	102	-0.05
91		Dibenzo(a,h)anthracene	40.000	35.910	10.2	101	-0.08
92		Benzo(g,h,i)perylene	40.000	36.354	9.1	103	-0.08

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110420\
Data File : BP003799.D
Acq On : 04 Nov 2020 09:32
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Nov 05 14:51:41 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 05 14:31:45 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)

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