Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081120\
Data File : BF121237.D
Acq On : 11 Aug 2020 13:33
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 11 16:13:46 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 29 14:55:43 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	90	-0.01
2		1,4-Dioxane	40.000	39.086	2.3	90	-0.02
3		Pyridine	40.000	39.629	0.9	92	-0.02
4		n-Nitrosodimethylamine	40.000	37.684	5.8	87	-0.02
5	S	2-Fluorophenol	80.000	78.993	1.3	92	0.00
6		Aniline	40.000	36.456	8.9	85	-0.01
7	S	Phenol-d6	80.000	77.521	3.1	91	0.00
8		2-Chlorophenol	40.000	39.664	0.8	91	0.00
9		Benzaldehyde	40.000	41.642	-4.1	94	0.00
10	C	Phenol	40.000	37.444	6.4	87	0.00
11		bis(2-Chloroethyl)ether	40.000	39.334	1.7	91	-0.01
12		1,3-Dichlorobenzene	40.000	39.692	0.8	93	-0.01
13	C	1,4-Dichlorobenzene	40.000	39.811	0.5	93	0.00
14		1,2-Dichlorobenzene	40.000	39.169	2.1	90	0.00
15		Benzyl Alcohol	40.000	37.051	7.4	86	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.403	1.5	92	-0.01
17		2-Methylphenol	40.000	39.046	2.4	92	-0.01
18		Hexachloroethane	40.000	40.450	-1.1	92	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	37.855	5.4	91	-0.01
20		3+4-Methylphenols	40.000	36.796	8.0	93	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	90	0.00
22		Acetophenone	40.000	39.948	0.1	94	0.00
23	S	Nitrobenzene-d5	80.000	82.110	-2.6	94	-0.01
24		Nitrobenzene	40.000	40.707	-1.8	94	-0.01
25		Isophorone	40.000	39.000	2.5	91	-0.01
26	C	2-Nitrophenol	40.000	44.285	-10.7	98	0.00
27		2,4-Dimethylphenol	40.000	39.676	0.8	93	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.654	-1.6	96	-0.01
29	C	2,4-Dichlorophenol	40.000	41.310	-3.3	94	0.00
30		1,2,4-Trichlorobenzene	40.000	40.642	-1.6	93	0.00
31		Naphthalene	40.000	39.842	0.4	93	0.00
32		Benzoic acid	40.000	32.778	18.1	67	0.00
33		4-Chloroaniline	40.000	39.491	1.3	93	-0.01
34	C	Hexachlorobutadiene	40.000	40.793	-2.0	94	-0.01
35		Caprolactam	40.000	40.123	-0.3	93	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.039	-2.6	95	0.00
37		2-Methylnaphthalene	40.000	40.492	-1.2	93	0.00
38		1-Methylnaphthalene	40.000	40.438	-1.1	94	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	91	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	42.524	-6.3	97	0.00
41	P	Hexachlorocyclopentadiene	40.000	37.727	5.7	82	0.00
42	S	2,4,6-Tribromophenol	80.000	80.489	-0.6	91	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.364	4.1	87	0.00
44		2,4,5-Trichlorophenol	40.000	38.997	2.5	89	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081120\
Data File : BF121237.D
Acq On : 11 Aug 2020 13:33
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 11 16:13:46 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 29 14:55:43 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	72.382	9.5	94	0.00
46		1,1'-Biphenyl	40.000	40.435	-1.1	93	0.00
47		2-Chloronaphthalene	40.000	40.483	-1.2	93	0.00
48		2-Nitroaniline	40.000	43.012	-7.5	98	0.00
49		Acenaphthylene	40.000	40.200	-0.5	93	0.00
50		Dimethylphthalate	40.000	40.700	-1.8	95	0.00
51		2,6-Dinitrotoluene	40.000	42.504	-6.3	95	0.00
52	C	Acenaphthene	40.000	37.470	6.3	86	0.00
53		3-Nitroaniline	40.000	40.430	-1.1	92	0.00
54	P	2,4-Dinitrophenol	40.000	35.502	11.2	85	0.01
55		Dibenzofuran	40.000	38.645	3.4	90	0.00
56	P	4-Nitrophenol	40.000	36.594	8.5	82	0.01
57		2,4-Dinitrotoluene	40.000	41.357	-3.4	91	0.00
58		Fluorene	40.000	39.610	1.0	95	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.052	4.9	87	0.00
60		Diethylphthalate	40.000	39.247	1.9	91	0.00
61		4-Chlorophenyl-phenylether	40.000	37.754	5.6	95	0.00
62		4-Nitroaniline	40.000	41.050	-2.6	94	0.00
63		Azobenzene	40.000	39.045	2.4	90	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	89	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.753	-4.4	100	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.268	-3.2	92	0.00
67		4-Bromophenyl-phenylether	40.000	42.311	-5.8	95	0.00
68		Hexachlorobenzene	40.000	42.153	-5.4	95	0.00
69		Atrazine	40.000	43.026	-7.6	98	0.00
70	C	Pentachlorophenol	40.000	35.392	11.5	74	0.01
71		Phenanthrene	40.000	40.317	-0.8	92	0.00
72		Anthracene	40.000	40.709	-1.8	91	0.00
73		Carbazole	40.000	40.707	-1.8	92	0.00
74		Di-n-butylphthalate	40.000	39.927	0.2	90	0.00
75	C	Fluoranthene	40.000	39.681	0.8	89	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	79	0.00
77		Benzidine	40.000	36.624	8.4	67	0.00
78		Pyrene	40.000	44.173	-10.4	88	0.00
79	S	Terphenyl-d14	80.000	83.074	-3.8	89	0.00
80		Butylbenzylphthalate	40.000	42.230	-5.6	82	0.00
81		Benzo(a)anthracene	40.000	42.777	-6.9	88	0.00
82		3,3'-Dichlorobenzidine	40.000	40.782	-2.0	81	0.00
83		Chrysene	40.000	40.313	-0.8	80	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	44.712	-11.8	89	0.00
85	c	Di-n-octyl phthalate	40.000	36.018	10.0	70	0.00
86	I	Perylene-d12	20.000	20.000	0.0	62	0.01
87		Indeno(1,2,3-cd)pyrene	40.000	35.266	11.8	54	0.03

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081120\
Data File : BF121237.D
Acq On : 11 Aug 2020 13:33
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 11 16:13:46 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 29 14:55:43 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	44.936	-12.3	71	0.01
89	Benzo(k)fluoranthene	40.000	43.782	-9.5	66	0.00
90 C	Benzo(a)pyrene	40.000	42.321	-5.8	65	0.02
91	Dibenzo(a,h)anthracene	40.000	35.669	10.8	55	0.02
92	Benzo(g,h,i)perylene	40.000	34.974	12.6	54	0.04

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