Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080123\
Data File : BF134547.D
Acq On : 01 Aug 2023 14:27
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 02 02:43:05 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 02 02:20:34 2023
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	103	0.00
2		1,4-Dioxane	40.000	36.984	7.5	96	0.00
3		Pyridine	40.000	37.584	6.0	98	0.00
4		n-Nitrosodimethylamine	40.000	38.319	4.2	98	0.00
5	S	2-Fluorophenol	80.000	74.381	7.0	97	0.00
6		Aniline	40.000	37.921	5.2	99	0.00
7	S	Phenol-d6	80.000	75.709	5.4	99	0.00
8		2-Chlorophenol	40.000	38.931	2.7	100	0.00
9		Benzaldehyde	40.000	35.734	10.7	97	0.00
10	C	Phenol	40.000	37.795	5.5	96	0.00
11		bis(2-Chloroethyl)ether	40.000	38.335	4.2	100	0.00
12		1,3-Dichlorobenzene	40.000	37.898	5.3	99	0.00
13	C	1,4-Dichlorobenzene	40.000	37.539	6.2	97	0.00
14		1,2-Dichlorobenzene	40.000	37.809	5.5	98	0.00
15		Benzyl Alcohol	40.000	38.812	3.0	100	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.463	3.8	100	0.00
17		2-Methylphenol	40.000	38.580	3.6	100	0.00
18		Hexachloroethane	40.000	39.467	1.3	103	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.295	6.8	99	0.00
20		3+4-Methylphenols	40.000	38.023	4.9	98	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	107	0.00
22		Acetophenone	40.000	36.484	8.8	99	0.00
23	S	Nitrobenzene-d5	80.000	83.256	-4.1	105	0.00
24		Nitrobenzene	40.000	40.478	-1.2	103	0.00
25		Isophorone	40.000	37.507	6.2	102	0.00
26	C	2-Nitrophenol	40.000	40.707	-1.8	114	0.00
27		2,4-Dimethylphenol	40.000	37.874	5.3	103	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.011	5.0	103	0.00
29	C	2,4-Dichlorophenol	40.000	38.217	4.5	100	0.00
30		1,2,4-Trichlorobenzene	40.000	37.981	5.0	102	0.00
31		Naphthalene	40.000	36.933	7.7	98	0.00
32		Benzoic acid	40.000	40.774	-1.9	121	0.00
33		4-Chloroaniline	40.000	36.487	8.8	99	0.00
34	C	Hexachlorobutadiene	40.000	37.905	5.2	102	0.00
35		Caprolactam	40.000	38.158	4.6	98	0.00
36	C	4-Chloro-3-methylphenol	40.000	37.988	5.0	101	0.00
37		2-Methylnaphthalene	40.000	37.265	6.8	101	0.00
38		1-Methylnaphthalene	40.000	37.431	6.4	101	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	109	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	37.332	6.7	101	0.00
41	P	Hexachlorocyclopentadiene	40.000	41.199	-3.0	103	0.00
42	S	2,4,6-Tribromophenol	80.000	80.907	-1.1	104	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.177	2.1	103	0.00
44		2,4,5-Trichlorophenol	40.000	39.535	1.2	104	0.00
45	S	2-Fluorobiphenyl	80.000	75.275	5.9	101	0.00
46		1,1'-Biphenyl	40.000	37.700	5.7	102	0.00
47		2-Chloronaphthalene	40.000	37.518	6.2	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080123\
Data File : BF134547.D
Acq On : 01 Aug 2023 14:27
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 02 02:43:05 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 02 02:20:34 2023
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	40.378	-0.9	109	0.00
49		Acenaphthylene	40.000	37.524	6.2	101	0.00
50		Dimethylphthalate	40.000	38.014	5.0	103	0.00
51		2,6-Dinitrotoluene	40.000	40.431	-1.1	108	0.00
52	C	Acenaphthene	40.000	38.053	4.9	103	0.00
53		3-Nitroaniline	40.000	43.533	-8.8	105	0.00
54	P	2,4-Dinitrophenol	40.000	41.901	-4.8	125	0.00
55		Dibenzofuran	40.000	37.424	6.4	99	0.00
56	P	4-Nitrophenol	40.000	40.307	-0.8	105	0.00
57		2,4-Dinitrotoluene	40.000	41.370	-3.4	110	0.00
58		Fluorene	40.000	36.332	9.2	101	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.257	1.9	102	0.00
60		Diethylphthalate	40.000	38.377	4.1	102	0.00
61		4-Chlorophenyl-phenylether	40.000	36.992	7.5	103	0.00
62		4-Nitroaniline	40.000	42.341	-5.9	103	0.00
63		Azobenzene	40.000	37.342	6.6	100	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	108	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.641	0.9	117	0.00
66	c	n-Nitrosodiphenylamine	40.000	37.951	5.1	102	0.00
67		4-Bromophenyl-phenylether	40.000	38.449	3.9	103	0.00
68		Hexachlorobenzene	40.000	37.857	5.4	100	0.00
69		Atrazine	40.000	39.579	1.1	107	0.00
70	C	Pentachlorophenol	40.000	42.417	-6.0	105	0.00
71		Phenanthrene	40.000	37.476	6.3	100	0.00
72		Anthracene	40.000	37.016	7.5	99	0.00
73		Carbazole	40.000	36.990	7.5	99	0.00
74		Di-n-butylphthalate	40.000	39.313	1.7	103	0.00
75	C	Fluoranthene	40.000	36.989	7.5	97	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
77		Benzidine	40.000	43.482	-8.7	100	0.00
78		Pyrene	40.000	38.655	3.4	98	0.00
79	S	Terphenyl-d14	80.000	76.921	3.8	98	0.00
80		Butylbenzylphthalate	40.000	42.714	-6.8	102	0.00
81		Benzo(a)anthracene	40.000	38.738	3.2	98	0.00
82		3,3'-Dichlorobenzidine	40.000	39.254	1.9	100	0.00
83		Chrysene	40.000	38.435	3.9	98	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.571	-3.9	101	0.00
85	c	Di-n-octyl phthalate	40.000	41.835	-4.6	102	0.00
86	I	Perylene-d12	20.000	20.000	0.0	103	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	39.587	1.0	95	0.00
88		Benzo(b)fluoranthene	40.000	38.414	4.0	99	0.00
89		Benzo(k)fluoranthene	40.000	39.336	1.7	95	0.00
90	C	Benzo(a)pyrene	40.000	38.721	3.2	96	0.00
91		Dibenzo(a,h)anthracene	40.000	40.459	-1.1	97	0.00
92		Benzo(g,h,i)perylene	40.000	39.729	0.7	95	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080123\
Data File : BF134547.D
Acq On : 01 Aug 2023 14:27
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 02 02:43:05 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 02 02:20:34 2023
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