Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011023\
Data File : BF131988.D
Acq On : 10 Jan 2023 09:56
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jan 11 00:33:00 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 10 00:32:25 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	102	0.00
2		1,4-Dioxane	40.000	38.140	4.6	96	0.00
3		Pyridine	40.000	39.699	0.8	97	0.00
4		n-Nitrosodimethylamine	40.000	40.611	-1.5	99	0.00
5	S	2-Fluorophenol	80.000	77.989	2.5	98	0.00
6		Aniline	40.000	38.852	2.9	97	0.00
7	S	Phenol-d6	80.000	77.829	2.7	98	0.00
8		2-Chlorophenol	40.000	40.116	-0.3	100	0.00
9		Benzaldehyde	40.000	37.406	6.5	97	0.00
10	C	Phenol	40.000	39.128	2.2	99	0.00
11		bis(2-Chloroethyl)ether	40.000	39.704	0.7	99	0.00
12		1,3-Dichlorobenzene	40.000	39.217	2.0	98	0.00
13	C	1,4-Dichlorobenzene	40.000	39.565	1.1	99	0.00
14		1,2-Dichlorobenzene	40.000	39.221	1.9	99	0.00
15		Benzyl Alcohol	40.000	39.878	0.3	98	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.860	2.9	97	0.00
17		2-Methylphenol	40.000	39.414	1.5	99	0.00
18		Hexachloroethane	40.000	40.102	-0.3	100	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.911	2.7	99	0.00
20		3+4-Methylphenols	40.000	38.982	2.5	97	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	100	0.00
22		Acetophenone	40.000	39.626	0.9	97	0.00
23	S	Nitrobenzene-d5	80.000	80.426	-0.5	98	0.00
24		Nitrobenzene	40.000	39.364	1.6	96	0.00
25		Isophorone	40.000	40.069	-0.2	98	0.00
26	C	2-Nitrophenol	40.000	41.146	-2.9	101	0.00
27		2,4-Dimethylphenol	40.000	40.517	-1.3	100	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.665	0.8	97	0.00
29	C	2,4-Dichlorophenol	40.000	40.028	-0.1	99	0.00
30		1,2,4-Trichlorobenzene	40.000	40.257	-0.6	98	0.00
31		Naphthalene	40.000	40.115	-0.3	99	0.00
32		Benzoic acid	40.000	39.861	0.3	93	0.00
33		4-Chloroaniline	40.000	41.060	-2.7	101	0.00
34	C	Hexachlorobutadiene	40.000	40.615	-1.5	99	0.00
35		Caprolactam	40.000	39.171	2.1	97	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.469	1.3	97	0.00
37		2-Methylnaphthalene	40.000	40.286	-0.7	99	0.00
38		1-Methylnaphthalene	40.000	40.286	-0.7	100	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	95	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.333	-3.3	98	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.053	9.9	89	0.00
42	S	2,4,6-Tribromophenol	80.000	82.003	-2.5	99	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.396	-3.5	97	0.00
44		2,4,5-Trichlorophenol	40.000	40.356	-0.9	96	0.00
45	S	2-Fluorobiphenyl	80.000	80.283	-0.4	98	0.00
46		1,1'-Biphenyl	40.000	40.645	-1.6	99	0.00
47		2-Chloronaphthalene	40.000	40.435	-1.1	97	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011023\
Data File : BF131988.D
Acq On : 10 Jan 2023 09:56
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jan 11 00:33:00 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 10 00:32:25 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	41.156	-2.9	99	0.00
49		Acenaphthylene	40.000	40.391	-1.0	98	0.00
50		Dimethylphthalate	40.000	40.307	-0.8	98	0.00
51		2,6-Dinitrotoluene	40.000	40.710	-1.8	97	0.00
52	C	Acenaphthene	40.000	40.176	-0.4	97	0.00
53		3-Nitroaniline	40.000	41.699	-4.2	101	0.00
54	P	2,4-Dinitrophenol	40.000	41.382	-3.5	95	0.00
55		Dibenzofuran	40.000	40.137	-0.3	98	0.00
56	P	4-Nitrophenol	40.000	39.749	0.6	92	0.00
57		2,4-Dinitrotoluene	40.000	40.011	-0.0	97	0.00
58		Fluorene	40.000	39.951	0.1	99	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.305	-3.3	96	0.00
60		Diethylphthalate	40.000	40.741	-1.9	99	0.00
61		4-Chlorophenyl-phenylether	40.000	40.183	-0.5	98	0.00
62		4-Nitroaniline	40.000	41.508	-3.8	99	0.00
63		Azobenzene	40.000	40.452	-1.1	99	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	99	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.731	-4.3	97	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.285	-0.7	99	0.00
67		4-Bromophenyl-phenylether	40.000	40.491	-1.2	99	0.00
68		Hexachlorobenzene	40.000	39.580	1.1	97	0.00
69		Atrazine	40.000	40.530	-1.3	102	0.00
70	C	Pentachlorophenol	40.000	35.811	10.5	89	0.00
71		Phenanthrene	40.000	40.022	-0.1	98	0.00
72		Anthracene	40.000	40.144	-0.4	99	0.00
73		Carbazole	40.000	39.880	0.3	99	0.00
74		Di-n-butylphthalate	40.000	40.517	-1.3	101	0.00
75	C	Fluoranthene	40.000	39.352	1.6	98	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	94	0.00
77		Benzidine	40.000	42.183	-5.5	107	0.00
78		Pyrene	40.000	42.588	-6.5	98	0.00
79	S	Terphenyl-d14	80.000	84.053	-5.1	97	0.00
80		Butylbenzylphthalate	40.000	42.459	-6.1	97	0.00
81		Benzo(a)anthracene	40.000	40.034	-0.1	93	0.00
82		3,3'-Dichlorobenzidine	40.000	40.897	-2.2	95	0.00
83		Chrysene	40.000	39.952	0.1	93	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	42.066	-5.2	96	0.00
85	c	Di-n-octyl phthalate	40.000	40.950	-2.4	96	0.00
86	I	Perylene-d12	20.000	20.000	0.0	97	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	42.931	-7.3	97	0.00
88		Benzo(b)fluoranthene	40.000	42.149	-5.4	100	0.00
89		Benzo(k)fluoranthene	40.000	37.473	6.3	90	0.00
90	C	Benzo(a)pyrene	40.000	40.138	-0.3	96	0.00
91		Dibenzo(a,h)anthracene	40.000	42.930	-7.3	97	0.00
92		Benzo(g,h,i)perylene	40.000	39.360	1.6	97	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011023\
Data File : BF131988.D
Acq On : 10 Jan 2023 09:56
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jan 11 00:33:00 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 10 00:32:25 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