Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131963.D
Acq On : 09 Jan 2023 15:05
Operator : CG\JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF010923

Quant Time: Jan 10 00:35:05 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	100	0.00
2		1,4-Dioxane	40.000	39.982	0.0	97	-0.01
3		Pyridine	40.000	40.956	-2.4	97	0.00
4		n-Nitrosodimethylamine	40.000	40.216	-0.5	95	-0.01
5	S	2-Fluorophenol	80.000	78.741	1.6	96	0.00
6		Aniline	40.000	39.653	0.9	97	0.00
7	S	Phenol-d6	80.000	78.954	1.3	97	0.00
8		2-Chlorophenol	40.000	40.433	-1.1	98	0.00
9		Benzaldehyde	40.000	38.371	4.1	97	0.00
10	C	Phenol	40.000	39.879	0.3	98	0.00
11		bis(2-Chloroethyl)ether	40.000	39.969	0.1	97	0.00
12		1,3-Dichlorobenzene	40.000	40.036	-0.1	97	0.00
13	C	1,4-Dichlorobenzene	40.000	39.961	0.1	97	0.00
14		1,2-Dichlorobenzene	40.000	39.362	1.6	97	0.00
15		Benzyl Alcohol	40.000	40.102	-0.3	96	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.087	-0.2	97	0.00
17		2-Methylphenol	40.000	40.007	-0.0	98	0.00
18		Hexachloroethane	40.000	39.582	1.0	96	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.839	0.4	99	0.00
20		3+4-Methylphenols	40.000	39.479	1.3	96	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22		Acetophenone	40.000	40.291	-0.7	97	0.00
23	S	Nitrobenzene-d5	80.000	80.711	-0.9	97	0.00
24		Nitrobenzene	40.000	40.558	-1.4	97	0.00
25		Isophorone	40.000	40.837	-2.1	98	0.00
26	C	2-Nitrophenol	40.000	41.258	-3.1	99	0.00
27		2,4-Dimethylphenol	40.000	40.747	-1.9	99	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.173	-0.4	97	0.00
29	C	2,4-Dichlorophenol	40.000	40.630	-1.6	98	0.00
30		1,2,4-Trichlorobenzene	40.000	40.394	-1.0	97	0.00
31		Naphthalene	40.000	40.174	-0.4	97	0.00
32		Benzoic acid	40.000	41.522	-3.8	94	0.00
33		4-Chloroaniline	40.000	40.861	-2.2	99	0.00
34	C	Hexachlorobutadiene	40.000	39.815	0.5	95	0.00
35		Caprolactam	40.000	40.136	-0.3	98	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.896	-2.2	98	0.00
37		2-Methylnaphthalene	40.000	41.100	-2.8	99	0.00
38		1-Methylnaphthalene	40.000	40.464	-1.2	98	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	95	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.056	-2.6	97	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.042	-0.1	101	0.00
42	S	2,4,6-Tribromophenol	80.000	84.248	-5.3	101	0.00
43	C	2,4,6-Trichlorophenol	40.000	41.232	-3.1	96	0.00
44		2,4,5-Trichlorophenol	40.000	41.366	-3.4	98	0.00
45	S	2-Fluorobiphenyl	80.000	81.103	-1.4	98	0.00
46		1,1'-Biphenyl	40.000	40.884	-2.2	99	0.00
47		2-Chloronaphthalene	40.000	41.563	-3.9	99	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131963.D
Acq On : 09 Jan 2023 15:05
Operator : CG\JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF010923

Quant Time: Jan 10 00:35:05 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	41.238	-3.1	98	0.00
49		Acenaphthylene	40.000	40.805	-2.0	99	0.00
50		Dimethylphthalate	40.000	41.442	-3.6	100	0.00
51		2,6-Dinitrotoluene	40.000	41.061	-2.7	97	0.00
52	C	Acenaphthene	40.000	41.126	-2.8	99	0.00
53		3-Nitroaniline	40.000	40.557	-1.4	98	0.00
54	P	2,4-Dinitrophenol	40.000	43.440	-8.6	100	0.00
55		Dibenzofuran	40.000	40.650	-1.6	98	0.00
56	P	4-Nitrophenol	40.000	43.454	-8.6	101	0.00
57		2,4-Dinitrotoluene	40.000	41.587	-4.0	101	0.00
58		Fluorene	40.000	40.616	-1.5	100	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	42.795	-7.0	100	0.00
60		Diethylphthalate	40.000	41.099	-2.7	100	0.00
61		4-Chlorophenyl-phenylether	40.000	40.852	-2.1	99	0.00
62		4-Nitroaniline	40.000	42.674	-6.7	102	0.00
63		Azobenzene	40.000	40.935	-2.3	99	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	40.000	42.957	-7.4	102	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.882	0.3	100	0.00
67		4-Bromophenyl-phenylether	40.000	39.966	0.1	101	0.00
68		Hexachlorobenzene	40.000	39.532	1.2	100	0.00
69		Atrazine	40.000	40.366	-0.9	104	0.00
70	C	Pentachlorophenol	40.000	39.288	1.8	101	0.00
71		Phenanthrene	40.000	39.532	1.2	100	0.00
72		Anthracene	40.000	40.142	-0.4	102	0.00
73		Carbazole	40.000	40.035	-0.1	102	0.00
74		Di-n-butylphthalate	40.000	40.833	-2.1	104	0.00
75	C	Fluoranthene	40.000	39.653	0.9	101	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
77		Benzidine	40.000	37.306	6.7	103	0.00
78		Pyrene	40.000	41.574	-3.9	103	0.00
79	S	Terphenyl-d14	80.000	82.677	-3.3	103	0.00
80		Butylbenzylphthalate	40.000	42.069	-5.2	104	0.00
81		Benzo(a)anthracene	40.000	40.308	-0.8	102	0.00
82		3,3'-Dichlorobenzidine	40.000	38.735	3.2	98	0.00
83		Chrysene	40.000	40.344	-0.9	102	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.265	-3.2	102	0.00
85	c	Di-n-octyl phthalate	40.000	41.570	-3.9	105	0.00
86	I	Perylene-d12	20.000	20.000	0.0	106	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	41.787	-4.5	102	0.00
88		Benzo(b)fluoranthene	40.000	40.196	-0.5	104	0.00
89		Benzo(k)fluoranthene	40.000	41.438	-3.6	108	0.00
90	C	Benzo(a)pyrene	40.000	40.672	-1.7	106	0.00
91		Dibenzo(a,h)anthracene	40.000	42.020	-5.1	104	0.00
92		Benzo(g,h,i)perylene	40.000	38.390	4.0	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
Data File : BF131963.D
Acq On : 09 Jan 2023 15:05
Operator : CG\JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF010923

Quant Time: Jan 10 00:35:05 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