Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112125\
Data File : BF144306.D
Acq On : 21 Nov 2025 14:03
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Nov 21 14:33:27 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 05 18:37:33 2025
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	81	-0.01
2		1,4-Dioxane	40.000	45.519	-13.8	87	-0.02
3		Pyridine	40.000	45.117	-12.8	88	-0.05
4		n-Nitrosodimethylamine	40.000	41.707	-4.3	83	-0.04
5	S	2-Fluorophenol	80.000	88.539	-10.7	87	-0.01
6		Aniline	40.000	43.733	-9.3	88	0.00
7	S	Phenol-d6	80.000	85.915	-7.4	86	0.00
8		2-Chlorophenol	40.000	44.216	-10.5	87	0.00
9		Benzaldehyde	40.000	51.268	-28.2#	103	-0.01
10	C	Phenol	40.000	43.122	-7.8	83	0.00
11		bis(2-Chloroethyl)ether	40.000	44.005	-10.0	87	0.00
12		1,3-Dichlorobenzene	40.000	43.491	-8.7	88	-0.01
13	C	1,4-Dichlorobenzene	40.000	43.546	-8.9	85	-0.01
14		1,2-Dichlorobenzene	40.000	43.256	-8.1	87	-0.01
15		Benzyl Alcohol	40.000	41.649	-4.1	83	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	44.313	-10.8	88	0.00
17		2-Methylphenol	40.000	42.573	-6.4	83	0.00
18		Hexachloroethane	40.000	43.096	-7.7	84	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	39.932	0.2	81	0.00
20		3+4-Methylphenols	40.000	41.707	-4.3	81	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	81	-0.01
22		Acetophenone	40.000	41.457	-3.6	83	0.00
23	S	Nitrobenzene-d5	80.000	84.777	-6.0	83	0.00
24		Nitrobenzene	40.000	42.961	-7.4	84	0.00
25		Isophorone	40.000	40.968	-2.4	81	0.00
26	C	2-Nitrophenol	40.000	44.784	-12.0	86	-0.01
27		2,4-Dimethylphenol	40.000	40.711	-1.8	77	0.00
28		bis(2-Chloroethoxy)methane	40.000	43.180	-7.9	85	0.00
29	C	2,4-Dichlorophenol	40.000	42.106	-5.3	83	0.00
30		1,2,4-Trichlorobenzene	40.000	41.913	-4.8	83	-0.01
31		Naphthalene	40.000	41.949	-4.9	84	0.00
32		Benzoic acid	40.000	39.480	1.3	91	-0.04
33		4-Chloroaniline	40.000	41.604	-4.0	81	-0.01
34	C	Hexachlorobutadiene	40.000	40.454	-1.1	79	-0.01
35		Caprolactam	40.000	37.384	6.5	72	-0.04
36	C	4-Chloro-3-methylphenol	40.000	40.008	-0.0	79	0.00
37		2-Methylnaphthalene	40.000	41.167	-2.9	82	-0.01
38		1-Methylnaphthalene	40.000	40.040	-0.1	80	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	74	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	42.076	-5.2	76	0.00
41	P	Hexachlorocyclopentadiene	40.000	41.498	-3.7	78	-0.02
42	S	2,4,6-Tribromophenol	80.000	71.948	10.1	63	0.00
43	C	2,4,6-Trichlorophenol	40.000	43.392	-8.5	79	-0.01
44		2,4,5-Trichlorophenol	40.000	41.200	-3.0	75	0.00
45	S	2-Fluorobiphenyl	80.000	85.863	-7.3	80	0.00
46		1,1'-Biphenyl	40.000	42.964	-7.4	79	-0.01
47		2-Chloronaphthalene	40.000	42.681	-6.7	78	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112125\
Data File : BF144306.D
Acq On : 21 Nov 2025 14:03
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 21 14:33:27 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 05 18:37:33 2025
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	41.685	-4.2	73	-0.01
49		Acenaphthylene	40.000	41.807	-4.5	76	-0.01
50		Dimethylphthalate	40.000	41.037	-2.6	75	-0.01
51		2,6-Dinitrotoluene	40.000	41.885	-4.7	75	0.00
52	C	Acenaphthene	40.000	39.578	1.1	73	0.00
53		3-Nitroaniline	40.000	40.044	-0.1	69	-0.01
54	P	2,4-Dinitrophenol	40.000	53.969	-34.9#	115	-0.02
55		Dibenzofuran	40.000	40.746	-1.9	74	0.00
56	P	4-Nitrophenol	40.000	37.887	5.3	73	-0.01
57		2,4-Dinitrotoluene	40.000	39.869	0.3	69	0.00
58		Fluorene	40.000	39.818	0.5	72	-0.01
59		2,3,4,6-Tetrachlorophenol	40.000	39.072	2.3	69	-0.01
60		Diethylphthalate	40.000	39.554	1.1	71	0.00
61		4-Chlorophenyl-phenylether	40.000	40.129	-0.3	71	-0.01
62		4-Nitroaniline	40.000	37.845	5.4	68	0.00
63		Azobenzene	40.000	39.882	0.3	72	-0.01
64	I	Phenanthrene-d10	20.000	20.000	0.0	63	-0.01
65		4,6-Dinitro-2-methylphenol	40.000	40.152	-0.4	66	-0.02
66	c	n-Nitrosodiphenylamine	40.000	45.791	-14.5	72	0.00
67		4-Bromophenyl-phenylether	40.000	44.183	-10.5	67	-0.01
68		Hexachlorobenzene	40.000	41.697	-4.2	66	0.00
69		Atrazine	40.000	39.642	0.9	62	-0.01
70	C	Pentachlorophenol	40.000	47.905	-19.8	72	-0.02
71		Phenanthrene	40.000	43.612	-9.0	68	-0.01
72		Anthracene	40.000	42.459	-6.1	66	-0.01
73		Carbazole	40.000	41.911	-4.8	64	-0.01
74		Di-n-butylphthalate	40.000	41.930	-4.8	64	-0.01
75	C	Fluoranthene	40.000	42.597	-6.5	65	-0.01
76	I	Chrysene-d12	20.000	20.000	0.0	82	0.00
77		Benzidine	40.000	39.483	1.3	77	-0.02
78		Pyrene	40.000	34.294	14.3	66	-0.01
79	S	Terphenyl-d14	80.000	66.870	16.4	65	-0.01
80		Butylbenzylphthalate	40.000	39.116	2.2	76	-0.01
81		Benzo(a)anthracene	40.000	42.837	-7.1	85	-0.01
82		3,3'-Dichlorobenzidine	40.000	42.084	-5.2	85	-0.02
83		Chrysene	40.000	41.375	-3.4	77	-0.01
84		Bis(2-ethylhexyl)phthalate	40.000	40.779	-1.9	78	-0.02
85	c	Di-n-octyl phthalate	40.000	45.431	-13.6	87	-0.02
86	I	Perylene-d12	20.000	20.000	0.0	78	-0.02
87		Indeno(1,2,3-cd)pyrene	40.000	33.989	15.0	64	-0.01
88		Benzo(b)fluoranthene	40.000	46.781	-17.0	84	0.00
89		Benzo(k)fluoranthene	40.000	44.155	-10.4	87	-0.01
90	C	Benzo(a)pyrene	40.000	43.616	-9.0	82	-0.01
91		Dibenzo(a,h)anthracene	40.000	34.802	13.0	65	-0.02
92		Benzo(g,h,i)perylene	40.000	31.827	20.4	59	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112125\
Data File : BF144306.D
Acq On : 21 Nov 2025 14:03
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Nov 21 14:33:27 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 05 18:37:33 2025
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