Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120823\
Data File : BF136464.D
Acq On : 08 Dec 2023 13:26
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Dec 08 13:54:10 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120523.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 06 09:40:27 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	91	0.00
2		1,4-Dioxane	40.000	37.846	5.4	82	-0.01
3		Pyridine	40.000	42.654	-6.6	90	-0.01
4		n-Nitrosodimethylamine	40.000	39.598	1.0	84	-0.02
5	S	2-Fluorophenol	80.000	75.002	6.2	82	0.00
6		Aniline	40.000	47.031	-17.6	101	0.00
7	S	Phenol-d6	80.000	75.111	6.1	82	0.00
8		2-Chlorophenol	40.000	37.868	5.3	83	0.00
9		Benzaldehyde	40.000	41.203	-3.0	87	0.00
10	C	Phenol	40.000	37.696	5.8	82	0.00
11		bis(2-Chloroethyl)ether	40.000	38.175	4.6	84	0.00
12		1,3-Dichlorobenzene	40.000	35.272	11.8	77	0.00
13	C	1,4-Dichlorobenzene	40.000	34.927	12.7	76	0.00
14		1,2-Dichlorobenzene	40.000	35.227	11.9	77	0.00
15		Benzyl Alcohol	40.000	40.279	-0.7	86	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.409	6.5	82	-0.01
17		2-Methylphenol	40.000	40.230	-0.6	87	0.00
18		Hexachloroethane	40.000	35.610	11.0	77	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.268	4.3	83	-0.01
20		3+4-Methylphenols	40.000	39.282	1.8	86	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	90	0.00
22		Acetophenone	40.000	37.924	5.2	82	-0.01
23	S	Nitrobenzene-d5	80.000	77.141	3.6	82	0.00
24		Nitrobenzene	40.000	37.200	7.0	80	0.00
25		Isophorone	40.000	38.025	4.9	81	0.00
26	C	2-Nitrophenol	40.000	38.973	2.6	81	0.00
27		2,4-Dimethylphenol	40.000	51.113	-27.8#	109	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.285	4.3	83	0.00
29	C	2,4-Dichlorophenol	40.000	38.807	3.0	82	0.00
30		1,2,4-Trichlorobenzene	40.000	36.281	9.3	78	0.00
31		Naphthalene	40.000	36.970	7.6	80	0.00
32		Benzoic acid	40.000	37.318	6.7	76	-0.01
33		4-Chloroaniline	40.000	42.743	-6.9	91	-0.01
34	C	Hexachlorobutadiene	40.000	35.731	10.7	77	0.00
35		Caprolactam	40.000	37.718	5.7	80	-0.01
36	C	4-Chloro-3-methylphenol	40.000	37.452	6.4	80	0.00
37		2-Methylnaphthalene	40.000	39.099	2.3	85	0.00
38		1-Methylnaphthalene	40.000	36.483	8.8	79	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	87	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.454	-1.1	85	0.00
41	P	Hexachlorocyclopentadiene	40.000	50.270	-25.7#	101	0.00
42	S	2,4,6-Tribromophenol	80.000	74.226	7.2	76	-0.01
43	C	2,4,6-Trichlorophenol	40.000	37.583	6.0	78	0.00
44		2,4,5-Trichlorophenol	40.000	39.847	0.4	82	0.00
45	S	2-Fluorobiphenyl	80.000	77.751	2.8	82	-0.01
46		1,1'-Biphenyl	40.000	39.210	2.0	82	-0.01
47		2-Chloronaphthalene	40.000	36.550	8.6	77	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120823\
Data File : BF136464.D
Acq On : 08 Dec 2023 13:26
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Dec 08 13:54:10 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120523.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 06 09:40:27 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	39.720	0.7	81	0.00
49		Acenaphthylene	40.000	39.740	0.6	83	0.00
50		Dimethylphthalate	40.000	39.136	2.2	82	-0.01
51		2,6-Dinitrotoluene	40.000	37.824	5.4	78	0.00
52	C	Acenaphthene	40.000	38.891	2.8	81	0.00
53		3-Nitroaniline	40.000	39.441	1.4	81	0.00
54	P	2,4-Dinitrophenol	40.000	32.856	17.9	63	0.00
55		Dibenzofuran	40.000	39.232	1.9	82	0.00
56	P	4-Nitrophenol	40.000	35.126	12.2	69	0.00
57		2,4-Dinitrotoluene	40.000	36.828	7.9	75	-0.01
58		Fluorene	40.000	37.215	7.0	79	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	37.635	5.9	78	0.00
60		Diethylphthalate	40.000	37.563	6.1	78	-0.01
61		4-Chlorophenyl-phenylether	40.000	38.738	3.2	83	0.00
62		4-Nitroaniline	40.000	37.371	6.6	75	-0.01
63		Azobenzene	40.000	38.225	4.4	80	-0.01
64	I	Phenanthrene-d10	20.000	20.000	0.0	79	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.180	2.1	68	0.00
66	c	n-Nitrosodiphenylamine	40.000	42.028	-5.1	82	0.00
67		4-Bromophenyl-phenylether	40.000	41.595	-4.0	79	0.00
68		Hexachlorobenzene	40.000	37.026	7.4	71	0.00
69		Atrazine	40.000	38.297	4.3	79	0.00
70	C	Pentachlorophenol	40.000	34.826	12.9	64	0.00
71		Phenanthrene	40.000	39.872	0.3	77	0.00
72		Anthracene	40.000	40.538	-1.3	78	0.00
73		Carbazole	40.000	36.437	8.9	70	0.00
74		Di-n-butylphthalate	40.000	37.975	5.1	72	0.00
75	C	Fluoranthene	40.000	34.784	13.0	67	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	66	0.00
77		Benzidine	40.000	30.772	23.1	46	0.00
78		Pyrene	40.000	40.927	-2.3	65	0.00
79	S	Terphenyl-d14	80.000	81.623	-2.0	65	0.00
80		Butylbenzylphthalate	40.000	39.904	0.2	61	0.00
81		Benzo(a)anthracene	40.000	39.565	1.1	64	0.00
82		3,3'-Dichlorobenzidine	40.000	43.574	-8.9	71	0.00
83		Chrysene	40.000	39.491	1.3	63	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	40.501	-1.3	62	0.00
85	c	Di-n-octyl phthalate	40.000	42.668	-6.7	66	0.00
86	I	Perylene-d12	20.000	20.000	0.0	86	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	44.203	-10.5	87	0.00
88		Benzo(b)fluoranthene	40.000	33.930	15.2	73	0.00
89		Benzo(k)fluoranthene	40.000	35.205	12.0	72	0.00
90	C	Benzo(a)pyrene	40.000	40.427	-1.1	84	0.00
91		Dibenzo(a,h)anthracene	40.000	44.844	-12.1	88	0.00
92		Benzo(g,h,i)perylene	40.000	43.650	-9.1	85	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120823\
Data File : BF136464.D
Acq On : 08 Dec 2023 13:26
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Dec 08 13:54:10 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120523.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 06 09:40:27 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