Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090222\
Data File : BF130111.D
Acq On : 02 Sep 2022 15:08
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Sep 03 05:44:28 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 31 18:15:32 2022
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	87	0.00
2		1,4-Dioxane	40.000	38.005	5.0	87	-0.04
3		Pyridine	40.000	38.812	3.0	88	-0.05
4		n-Nitrosodimethylamine	40.000	37.496	6.3	83	-0.04
5	S	2-Fluorophenol	80.000	77.670	2.9	90	-0.01
6		Aniline	40.000	38.261	4.3	85	0.00
7	S	Phenol-d6	80.000	77.167	3.5	89	0.00
8		2-Chlorophenol	40.000	39.051	2.4	88	0.00
9		Benzaldehyde	40.000	36.859	7.9	92	0.00
10	C	Phenol	40.000	38.847	2.9	89	0.00
11		bis(2-Chloroethyl)ether	40.000	38.029	4.9	87	0.00
12		1,3-Dichlorobenzene	40.000	37.579	6.1	86	0.00
13	C	1,4-Dichlorobenzene	40.000	38.219	4.5	88	0.00
14		1,2-Dichlorobenzene	40.000	38.030	4.9	89	0.00
15		Benzyl Alcohol	40.000	35.753	10.6	81	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	36.037	9.9	81	0.00
17		2-Methylphenol	40.000	37.735	5.7	86	0.00
18		Hexachloroethane	40.000	39.232	1.9	87	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.541	8.6	84	-0.01
20		3+4-Methylphenols	40.000	38.063	4.8	90	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	86	0.00
22		Acetophenone	40.000	36.859	7.9	86	0.00
23	S	Nitrobenzene-d5	80.000	84.276	-5.3	91	0.00
24		Nitrobenzene	40.000	40.907	-2.3	90	0.00
25		Isophorone	40.000	37.145	7.1	84	0.00
26	C	2-Nitrophenol	40.000	44.261	-10.7	101	0.00
27		2,4-Dimethylphenol	40.000	39.604	1.0	90	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.429	6.4	85	0.00
29	C	2,4-Dichlorophenol	40.000	39.169	2.1	88	0.00
30		1,2,4-Trichlorobenzene	40.000	38.249	4.4	88	0.00
31		Naphthalene	40.000	37.619	6.0	87	0.00
32		Benzoic acid	40.000	41.388	-3.5	94	-0.02
33		4-Chloroaniline	40.000	38.208	4.5	86	0.00
34	C	Hexachlorobutadiene	40.000	37.947	5.1	86	0.00
35		Caprolactam	40.000	38.132	4.7	85	-0.01
36	C	4-Chloro-3-methylphenol	40.000	38.619	3.5	87	0.00
37		2-Methylnaphthalene	40.000	37.706	5.7	88	0.00
38		1-Methylnaphthalene	40.000	37.227	6.9	87	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	85	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.272	1.8	89	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.116	-0.3	84	0.00
42	S	2,4,6-Tribromophenol	80.000	79.429	0.7	86	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.703	0.7	86	0.00
44		2,4,5-Trichlorophenol	40.000	39.136	2.2	85	0.00
45	S	2-Fluorobiphenyl	80.000	73.598	8.0	88	0.00
46		1,1'-Biphenyl	40.000	38.673	3.3	89	0.00
47		2-Chloronaphthalene	40.000	38.688	3.3	88	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090222\
Data File : BF130111.D
Acq On : 02 Sep 2022 15:08
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 03 05:44:28 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 31 18:15:32 2022
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	46.060	-15.2	92	0.00
49		Acenaphthylene	40.000	38.359	4.1	87	0.00
50		Dimethylphthalate	40.000	37.749	5.6	86	-0.01
51		2,6-Dinitrotoluene	40.000	43.487	-8.7	90	0.00
52	C	Acenaphthene	40.000	39.040	2.4	88	0.00
53		3-Nitroaniline	40.000	43.037	-7.6	91	0.00
54	P	2,4-Dinitrophenol	40.000	46.514	-16.3	111	0.00
55		Dibenzofuran	40.000	38.441	3.9	87	0.00
56	P	4-Nitrophenol	40.000	42.344	-5.9	88	-0.01
57		2,4-Dinitrotoluene	40.000	46.984	-17.5	93	0.00
58		Fluorene	40.000	36.928	7.7	87	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.371	1.6	86	0.00
60		Diethylphthalate	40.000	37.270	6.8	83	0.00
61		4-Chlorophenyl-phenylether	40.000	37.216	7.0	88	0.00
62		4-Nitroaniline	40.000	41.720	-4.3	87	0.00
63		Azobenzene	40.000	36.967	7.6	83	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	81	0.00
65		4,6-Dinitro-2-methylphenol	40.000	46.337	-15.8	101	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.190	2.0	85	0.00
67		4-Bromophenyl-phenylether	40.000	39.144	2.1	84	0.00
68		Hexachlorobenzene	40.000	38.239	4.4	81	0.00
69		Atrazine	40.000	37.298	6.8	78	0.00
70	C	Pentachlorophenol	40.000	41.005	-2.5	81	0.00
71		Phenanthrene	40.000	37.201	7.0	83	0.00
72		Anthracene	40.000	37.678	5.8	83	0.00
73		Carbazole	40.000	36.424	8.9	80	0.00
74		Di-n-butylphthalate	40.000	36.047	9.9	77	0.00
75	C	Fluoranthene	40.000	34.622	13.4	77	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	70	0.00
77		Benzidine	40.000	43.286	-8.2	75	0.00
78		Pyrene	40.000	42.373	-5.9	76	0.00
79	S	Terphenyl-d14	80.000	81.400	-1.8	74	0.00
80		Butylbenzylphthalate	40.000	39.123	2.2	65	0.00
81		Benzo(a)anthracene	40.000	38.902	2.7	72	0.00
82		3,3'-Dichlorobenzidine	40.000	41.764	-4.4	74	-0.01
83		Chrysene	40.000	39.341	1.6	73	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	37.205	7.0	64	0.00
85	c	Di-n-octyl phthalate	40.000	39.569	1.1	66	0.00
86	I	Perylene-d12	20.000	20.000	0.0	88	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	47.961	-19.9	103	0.00
88		Benzo(b)fluoranthene	40.000	34.270	14.3	82	0.00
89		Benzo(k)fluoranthene	40.000	37.756	5.6	84	0.00
90	C	Benzo(a)pyrene	40.000	38.664	3.3	88	0.00
91		Dibenzo(a,h)anthracene	40.000	48.448	-21.1	103	0.00
92		Benzo(g,h,i)perylene	40.000	49.031	-22.6	104	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090222\
Data File : BF130111.D
Acq On : 02 Sep 2022 15:08
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Sep 03 05:44:28 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 31 18:15:32 2022
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