Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090523\
Data File : BF135104.D
Acq On : 05 Sep 2023 20:46
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Sep 06 01:56:51 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 31 03:48:08 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	67	0.00
2		1,4-Dioxane	0.548	0.503	8.2	66	-0.01
3		Pyridine	1.479	1.345	9.1	65	0.00
4		n-Nitrosodimethylamine	0.723	0.691	4.4	69	-0.01
5	S	2-Fluorophenol	1.295	1.221	5.7	71	0.00
6		Aniline	2.001	1.844	7.8	68	0.00
7	S	Phenol-d6	1.614	1.521	5.8	70	0.00
8		2-Chlorophenol	1.315	1.249	5.0	70	0.00
9		Benzaldehyde	0.822	0.837	-1.8	79	0.00
10	C	Phenol	1.694	1.578	6.8	69	0.00
11		bis(2-Chloroethyl)ether	1.276	1.156	9.4	67	0.00
12		1,3-Dichlorobenzene	1.355	1.265	6.6	69	0.00
13	C	1,4-Dichlorobenzene	1.356	1.265	6.7	68	0.00
14		1,2-Dichlorobenzene	1.266	1.195	5.6	70	0.00
15		Benzyl Alcohol	1.106	1.089	1.5	72	0.00
16		2,2'-oxybis(1-Chloropropane	2.137	1.852	13.3	64	0.00
17		2-Methylphenol	1.146	1.077	6.0	69	0.00
18		Hexachloroethane	0.502	0.467	7.0	68	0.00
19	P	n-Nitroso-di-n-propylamine	0.959	0.864	9.9	70	0.00
20		3+4-Methylphenols	1.352	1.308	3.3	73	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	68	0.00
22		Acetophenone	0.445	0.421	5.4	72	0.00
23	S	Nitrobenzene-d5	0.356	0.335	5.9	71	0.00
24		Nitrobenzene	0.365	0.346	5.2	70	0.00
25		Isophorone	0.675	0.612	9.3	68	0.00
26	C	2-Nitrophenol	0.180	0.173	3.9	70	0.00
27		2,4-Dimethylphenol	0.290	0.267	7.9	68	0.00
28		bis(2-Chloroethoxy)methane	0.407	0.362	11.1	67	0.00
29	C	2,4-Dichlorophenol	0.276	0.259	6.2	70	0.00
30		1,2,4-Trichlorobenzene	0.298	0.278	6.7	69	0.00
31		Naphthalene	0.977	0.902	7.7	70	0.00
32		Benzoic acid	0.232	0.212	8.6	66	0.00
33		4-Chloroaniline	0.427	0.395	7.5	69	0.00
34	C	Hexachlorobutadiene	0.169	0.161	4.7	70	0.00
35		Caprolactam	0.090	0.087	3.3	70	0.00
36	C	4-Chloro-3-methylphenol	0.300	0.290	3.3	72	0.00
37		2-Methylnaphthalene	0.652	0.602	7.7	69	0.00
38		1-Methylnaphthalene	0.610	0.563	7.7	70	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	68	0.00
40		1,2,4,5-Tetrachlorobenzene	0.560	0.523	6.6	71	0.00
41	P	Hexachlorocyclopentadiene	0.238	0.239	-0.4	72	0.00
42	S	2,4,6-Tribromophenol	0.212	0.207	2.4	74	0.00
43	C	2,4,6-Trichlorophenol	0.376	0.349	7.2	70	0.00
44		2,4,5-Trichlorophenol	0.435	0.403	7.4	70	0.00
45	S	2-Fluorobiphenyl	1.228	1.156	5.9	73	0.00
46		1,1'-Biphenyl	1.528	1.401	8.3	70	0.00
47		2-Chloronaphthalene	1.149	1.063	7.5	70	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090523\
Data File : BF135104.D
Acq On : 05 Sep 2023 20:46
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Sep 06 01:56:51 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 31 03:48:08 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.393	0.386	1.8	72	0.00
49		Acenaphthylene	1.737	1.620	6.7	71	0.00
50		Dimethylphthalate	1.400	1.285	8.2	70	0.00
51		2,6-Dinitrotoluene	0.307	0.296	3.6	71	0.00
52	C	Acenaphthene	1.107	1.038	6.2	72	0.00
53		3-Nitroaniline	0.358	0.330	7.8	69	0.00
54	P	2,4-Dinitrophenol	0.147	0.140	4.8	65	0.00
55		Dibenzofuran	1.597	1.484	7.1	71	0.00
56	P	4-Nitrophenol	0.252	0.228	9.5	66	0.00
57		2,4-Dinitrotoluene	0.383	0.382	0.3	73	0.00
58		Fluorene	1.217	1.179	3.1	76	0.00
59		2,3,4,6-Tetrachlorophenol	0.335	0.306	8.7	68	0.00
60		Diethylphthalate	1.318	1.236	6.2	72	0.00
61		4-Chlorophenyl-phenylether	0.602	0.559	7.1	72	0.00
62		4-Nitroaniline	0.350	0.337	3.7	71	0.00
63		Azobenzene	1.327	1.228	7.5	70	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	72	0.00
65		4,6-Dinitro-2-methylphenol	0.121	0.115	5.0	69	0.00
66	c	n-Nitrosodiphenylamine	0.622	0.551	11.4	70	0.00
67		4-Bromophenyl-phenylether	0.212	0.189	10.8	71	0.00
68		Hexachlorobenzene	0.219	0.199	9.1	72	0.00
69		Atrazine	0.161	0.135	16.1	68	0.00
70	C	Pentachlorophenol	0.133	0.116	12.8	67	0.00
71		Phenanthrene	1.032	0.934	9.5	73	0.00
72		Anthracene	1.054	0.952	9.7	72	0.00
73		Carbazole	0.909	0.837	7.9	73	0.00
74		Di-n-butylphthalate	1.096	0.977	10.9	71	0.00
75	C	Fluoranthene	1.043	0.951	8.8	73	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	71	0.00
77		Benzidine	0.361	0.328	9.1	66	0.00
78		Pyrene	1.899	1.750	7.8	72	0.00
79	S	Terphenyl-d14	1.226	1.132	7.7	73	0.00
80		Butylbenzylphthalate	0.686	0.627	8.6	69	0.00
81		Benzo(a)anthracene	1.364	1.250	8.4	73	0.00
82		3,3'-Dichlorobenzidine	0.412	0.402	2.4	78	0.00
83		Chrysene	1.341	1.207	10.0	72	0.00
84		Bis(2-ethylhexyl)phthalate	0.710	0.668	5.9	72	0.00
85	c	Di-n-octyl phthalate	1.084	1.015	6.4	73	0.00
86	I	Perylene-d12	1.000	1.000	0.0	77	0.00
87		Indeno(1,2,3-cd)pyrene	1.347	1.315	2.4	78	0.00
88		Benzo(b)fluoranthene	1.244	1.113	10.5	79	0.00
89		Benzo(k)fluoranthene	1.202	1.029	14.4	71	0.00
90	C	Benzo(a)pyrene	1.156	1.039	10.1	76	0.00
91		Dibenzo(a,h)anthracene	1.091	1.075	1.5	78	0.00
92		Benzo(g,h,i)perylene	1.135	1.133	0.2	80	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090523\
Data File : BF135104.D
Acq On : 05 Sep 2023 20:46
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Sep 06 01:56:51 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 31 03:48:08 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	AvgRF	CCRF	%Dev	Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 0