Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051122\
Data File : BF128191.D
Acq On : 11 May 2022 13:23
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: May 12 05:59:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF050922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 09 15:30:11 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	98	0.00
2		1,4-Dioxane	40.000	41.444	-3.6	99	0.01
3		Pyridine	40.000	41.228	-3.1	101	0.01
4		n-Nitrosodimethylamine	40.000	42.338	-5.8	100	0.01
5	S	2-Fluorophenol	80.000	80.398	-0.5	99	0.00
6		Aniline	40.000	41.010	-2.5	99	0.00
7	S	Phenol-d6	80.000	79.782	0.3	98	0.00
8		2-Chlorophenol	40.000	39.718	0.7	97	0.00
9		Benzaldehyde	40.000	45.145	-12.9	107	0.00
10	C	Phenol	40.000	40.110	-0.3	96	0.00
11		bis(2-Chloroethyl)ether	40.000	39.725	0.7	97	0.00
12		1,3-Dichlorobenzene	40.000	40.187	-0.5	98	0.00
13	C	1,4-Dichlorobenzene	40.000	40.276	-0.7	98	0.00
14		1,2-Dichlorobenzene	40.000	39.022	2.4	96	0.00
15		Benzyl Alcohol	40.000	39.942	0.1	97	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.904	0.2	96	0.00
17		2-Methylphenol	40.000	40.523	-1.3	98	0.00
18		Hexachloroethane	40.000	41.168	-2.9	100	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.125	2.2	96	0.00
20		3+4-Methylphenols	40.000	39.839	0.4	98	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	96	0.00
22		Acetophenone	40.000	40.016	-0.0	98	0.00
23	S	Nitrobenzene-d5	80.000	81.650	-2.1	98	0.00
24		Nitrobenzene	40.000	40.977	-2.4	97	0.00
25		Isophorone	40.000	40.381	-1.0	96	0.00
26	C	2-Nitrophenol	40.000	41.729	-4.3	98	0.00
27		2,4-Dimethylphenol	40.000	40.641	-1.6	97	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.542	-1.4	97	0.00
29	C	2,4-Dichlorophenol	40.000	41.283	-3.2	98	0.00
30		1,2,4-Trichlorobenzene	40.000	40.697	-1.7	98	0.00
31		Naphthalene	40.000	40.153	-0.4	96	0.00
32		Benzoic acid	40.000	40.967	-2.4	91	0.00
33		4-Chloroaniline	40.000	39.264	1.8	95	0.00
34	C	Hexachlorobutadiene	40.000	40.408	-1.0	96	0.00
35		Caprolactam	40.000	38.280	4.3	90	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.208	-0.5	95	0.00
37		2-Methylnaphthalene	40.000	39.943	0.1	96	0.00
38		1-Methylnaphthalene	40.000	39.722	0.7	97	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	95	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.193	-3.0	97	0.00
41	P	Hexachlorocyclopentadiene	40.000	37.109	7.2	85	0.00
42	S	2,4,6-Tribromophenol	80.000	78.766	1.5	92	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.903	-2.3	94	0.00
44		2,4,5-Trichlorophenol	40.000	41.239	-3.1	96	0.00
45	S	2-Fluorobiphenyl	80.000	79.708	0.4	96	0.00
46		1,1'-Biphenyl	40.000	40.470	-1.2	95	0.00
47		2-Chloronaphthalene	40.000	40.282	-0.7	95	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051122\
Data File : BF128191.D
Acq On : 11 May 2022 13:23
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: May 12 05:59:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF050922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 09 15:30:11 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	41.119	-2.8	93	0.00
49		Acenaphthylene	40.000	40.145	-0.4	94	0.00
50		Dimethylphthalate	40.000	40.119	-0.3	95	0.00
51		2,6-Dinitrotoluene	40.000	39.726	0.7	92	0.00
52	C	Acenaphthene	40.000	40.156	-0.4	94	0.00
53		3-Nitroaniline	40.000	39.072	2.3	91	0.00
54	P	2,4-Dinitrophenol	40.000	38.399	4.0	86	0.00
55		Dibenzofuran	40.000	39.716	0.7	95	0.00
56	P	4-Nitrophenol	40.000	38.196	4.5	87	0.00
57		2,4-Dinitrotoluene	40.000	40.136	-0.3	92	0.00
58		Fluorene	40.000	39.242	1.9	93	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.057	-0.1	92	0.00
60		Diethylphthalate	40.000	39.789	0.5	94	0.00
61		4-Chlorophenyl-phenylether	40.000	39.718	0.7	94	0.00
62		4-Nitroaniline	40.000	38.711	3.2	90	0.00
63		Azobenzene	40.000	39.958	0.1	94	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	91	0.00
65		4,6-Dinitro-2-methylphenol	40.000	41.804	-4.5	89	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.500	-1.3	92	0.00
67		4-Bromophenyl-phenylether	40.000	41.074	-2.7	94	0.00
68		Hexachlorobenzene	40.000	40.788	-2.0	93	0.00
69		Atrazine	40.000	39.756	0.6	91	0.00
70	C	Pentachlorophenol	40.000	40.575	-1.4	83	0.00
71		Phenanthrene	40.000	39.763	0.6	92	0.00
72		Anthracene	40.000	39.913	0.2	92	0.00
73		Carbazole	40.000	39.265	1.8	89	0.00
74		Di-n-butylphthalate	40.000	39.046	2.4	90	0.00
75	C	Fluoranthene	40.000	38.237	4.4	89	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	82	0.00
77		Benzidine	40.000	46.331	-15.8	94	0.00
78		Pyrene	40.000	43.286	-8.2	88	0.00
79	S	Terphenyl-d14	80.000	84.645	-5.8	88	0.00
80		Butylbenzylphthalate	40.000	42.390	-6.0	85	0.00
81		Benzo(a)anthracene	40.000	40.751	-1.9	83	0.00
82		3,3'-Dichlorobenzidine	40.000	41.822	-4.6	88	0.00
83		Chrysene	40.000	39.984	0.0	83	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.186	-3.0	83	0.00
85	c	Di-n-octyl phthalate	40.000	38.956	2.6	78	0.00
86	I	Perylene-d12	20.000	20.000	0.0	81	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	43.935	-9.8	91	0.00
88		Benzo(b)fluoranthene	40.000	42.173	-5.4	85	0.00
89		Benzo(k)fluoranthene	40.000	38.418	4.0	76	0.00
90	C	Benzo(a)pyrene	40.000	40.801	-2.0	82	0.00
91		Dibenzo(a,h)anthracene	40.000	44.234	-10.6	91	0.00
92		Benzo(g,h,i)perylene	40.000	43.484	-8.7	92	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051122\
Data File : BF128191.D
Acq On : 11 May 2022 13:23
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: May 12 05:59:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF050922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 09 15:30:11 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