Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129709.D
Acq On : 11 Aug 2022 10:34
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 11 23:52:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 11 23:50:03 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	56	0.00
2		1,4-Dioxane	40.000	36.344	9.1	51	0.00
3		Pyridine	40.000	35.592	11.0	49	0.00
4		n-Nitrosodimethylamine	40.000	36.168	9.6	50	0.00
5	S	2-Fluorophenol	80.000	78.728	1.6	56	0.00
6		Aniline	40.000	35.627	10.9	50	0.00
7	S	Phenol-d6	80.000	78.086	2.4	56	0.00
8		2-Chlorophenol	40.000	40.492	-1.2	57	0.00
9		Benzaldehyde	40.000	34.160	14.6	51	0.00
10	C	Phenol	40.000	40.180	-0.4	57	0.00
11		bis(2-Chloroethyl)ether	40.000	39.014	2.5	55	0.00
12		1,3-Dichlorobenzene	40.000	40.451	-1.1	57	0.00
13	C	1,4-Dichlorobenzene	40.000	40.328	-0.8	57	0.00
14		1,2-Dichlorobenzene	40.000	40.732	-1.8	57	0.00
15		Benzyl Alcohol	40.000	40.875	-2.2	57	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	34.685	13.3	49	0.00
17		2-Methylphenol	40.000	39.115	2.2	56	0.00
18		Hexachloroethane	40.000	40.695	-1.7	57	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.753	-4.4	60	0.00
20		3+4-Methylphenols	40.000	40.744	-1.9	60	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	56	0.00
22		Acetophenone	40.000	41.714	-4.3	61	0.00
23	S	Nitrobenzene-d5	80.000	81.758	-2.2	58	0.00
24		Nitrobenzene	40.000	39.997	0.0	56	0.00
25		Isophorone	40.000	39.313	1.7	57	0.00
26	C	2-Nitrophenol	40.000	40.894	-2.2	57	0.00
27		2,4-Dimethylphenol	40.000	40.545	-1.4	58	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.468	-1.2	58	0.00
29	C	2,4-Dichlorophenol	40.000	40.803	-2.0	57	0.00
30		1,2,4-Trichlorobenzene	40.000	41.108	-2.8	58	0.00
31		Naphthalene	40.000	41.108	-2.8	59	0.00
32		Benzoic acid	40.000	23.886	40.3#	32	0.00
33		4-Chloroaniline	40.000	39.827	0.4	57	0.00
34	C	Hexachlorobutadiene	40.000	44.275	-10.7	63	0.00
35		Caprolactam	40.000	39.295	1.8	56	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.074	-2.7	58	0.00
37		2-Methylnaphthalene	40.000	41.414	-3.5	59	0.00
38		1-Methylnaphthalene	40.000	41.656	-4.1	61	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	58	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	42.027	-5.1	62	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.440	8.9	49	0.00
42	S	2,4,6-Tribromophenol	80.000	82.174	-2.7	61	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.737	3.2	56	0.00
44		2,4,5-Trichlorophenol	40.000	38.326	4.2	56	0.00
45	S	2-Fluorobiphenyl	80.000	79.140	1.1	65	0.00
46		1,1'-Biphenyl	40.000	41.014	-2.5	61	0.00
47		2-Chloronaphthalene	40.000	40.446	-1.1	60	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129709.D
Acq On : 11 Aug 2022 10:34
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 11 23:52:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 11 23:50:03 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	39.283	1.8	56	0.00
49		Acenaphthylene	40.000	40.447	-1.1	60	0.00
50		Dimethylphthalate	40.000	41.849	-4.6	62	0.00
51		2,6-Dinitrotoluene	40.000	41.044	-2.6	60	0.00
52	C	Acenaphthene	40.000	41.117	-2.8	60	0.00
53		3-Nitroaniline	40.000	37.730	5.7	54	0.00
54	P	2,4-Dinitrophenol	40.000	36.002	10.0	49	0.00
55		Dibenzofuran	40.000	41.019	-2.5	61	0.00
56	P	4-Nitrophenol	40.000	36.699	8.3	52	0.00
57		2,4-Dinitrotoluene	40.000	40.577	-1.4	57	0.00
58		Fluorene	40.000	42.415	-6.0	64	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	37.613	6.0	54	0.00
60		Diethylphthalate	40.000	42.788	-7.0	63	0.00
61		4-Chlorophenyl-phenylether	40.000	43.903	-9.8	66	0.00
62		4-Nitroaniline	40.000	36.369	9.1	53	0.00
63		Azobenzene	40.000	40.112	-0.3	59	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	58	0.00
65		4,6-Dinitro-2-methylphenol	40.000	38.922	2.7	53	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.640	-1.6	60	0.00
67		4-Bromophenyl-phenylether	40.000	43.293	-8.2	64	0.00
68		Hexachlorobenzene	40.000	43.482	-8.7	64	0.00
69		Atrazine	40.000	42.654	-6.6	63	0.00
70	C	Pentachlorophenol	40.000	33.729	15.7	47	0.00
71		Phenanthrene	40.000	41.290	-3.2	62	0.00
72		Anthracene	40.000	41.169	-2.9	62	0.00
73		Carbazole	40.000	39.477	1.3	59	0.00
74		Di-n-butylphthalate	40.000	44.890	-12.2	67	0.00
75	C	Fluoranthene	40.000	41.293	-3.2	62	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	59	0.00
77		Benzidine	40.000	38.975	2.6	55	0.00
78		Pyrene	40.000	41.272	-3.2	61	0.00
79	S	Terphenyl-d14	80.000	90.468	-13.1	69	0.00
80		Butylbenzylphthalate	40.000	45.781	-14.5	66	0.00
81		Benzo(a)anthracene	40.000	40.311	-0.8	61	0.00
82		3,3'-Dichlorobenzidine	40.000	37.567	6.1	55	0.00
83		Chrysene	40.000	39.699	0.8	60	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	46.933	-17.3	70	0.00
85	c	Di-n-octyl phthalate	40.000	48.748	-21.9#	73	0.00
86	I	Perylene-d12	20.000	20.000	0.0	60	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	42.332	-5.8	59	0.00
88		Benzo(b)fluoranthene	40.000	41.385	-3.5	65	0.00
89		Benzo(k)fluoranthene	40.000	39.852	0.4	59	0.00
90	C	Benzo(a)pyrene	40.000	40.149	-0.4	60	0.00
91		Dibenzo(a,h)anthracene	40.000	42.780	-7.0	59	0.00
92		Benzo(g,h,i)perylene	40.000	42.177	-5.4	57	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081122\
Data File : BF129709.D
Acq On : 11 Aug 2022 10:34
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 11 23:52:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 11 23:50:03 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 = 1