Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091522\
Data File : BP011719.D
Acq On : 15 Sep 2022 23:22
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 16 03:01:48 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 14 19:12:50 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	112	0.00
2		1,4-Dioxane	40.000	39.043	2.4	114	0.00
3		Pyridine	40.000	41.467	-3.7	118	0.00
4		n-Nitrosodimethylamine	40.000	42.521	-6.3	121	0.00
5	S	2-Fluorophenol	80.000	81.428	-1.8	116	0.00
6		Aniline	40.000	41.860	-4.6	117	0.00
7	S	Phenol-d6	80.000	84.605	-5.8	118	0.00
8		2-Chlorophenol	40.000	41.022	-2.6	116	0.00
9		Benzaldehyde	40.000	41.658	-4.1	123	0.00
10	C	Phenol	40.000	41.917	-4.8	118	0.00
11		bis(2-Chloroethyl)ether	40.000	40.921	-2.3	116	0.00
12		1,3-Dichlorobenzene	40.000	39.691	0.8	113	-0.01
13	C	1,4-Dichlorobenzene	40.000	39.708	0.7	113	0.00
14		1,2-Dichlorobenzene	40.000	39.973	0.1	114	0.00
15		Benzyl Alcohol	40.000	42.602	-6.5	119	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	42.861	-7.2	121	-0.01
17		2-Methylphenol	40.000	42.102	-5.3	118	0.00
18		Hexachloroethane	40.000	41.050	-2.6	117	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.836	-7.1	118	0.00
20		3+4-Methylphenols	40.000	42.048	-5.1	117	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	116	-0.01
22		Acetophenone	40.000	40.668	-1.7	118	0.00
23	S	Nitrobenzene-d5	80.000	83.229	-4.0	119	0.00
24		Nitrobenzene	40.000	41.398	-3.5	119	0.00
25		Isophorone	40.000	41.362	-3.4	117	0.00
26	C	2-Nitrophenol	40.000	36.398	9.0	111	0.00
27		2,4-Dimethylphenol	40.000	41.074	-2.7	116	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.624	-1.6	117	0.00
29	C	2,4-Dichlorophenol	40.000	40.533	-1.3	115	0.00
30		1,2,4-Trichlorobenzene	40.000	38.842	2.9	113	0.00
31		Naphthalene	40.000	39.836	0.4	116	0.00
32		Benzoic acid	40.000	36.057	9.9	110	0.00
33		4-Chloroaniline	40.000	40.640	-1.6	116	0.00
34	C	Hexachlorobutadiene	40.000	37.873	5.3	111	0.00
35		Caprolactam	40.000	39.330	1.7	112	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.920	-2.3	116	0.00
37		2-Methylnaphthalene	40.000	39.766	0.6	115	0.00
38		1-Methylnaphthalene	40.000	39.876	0.3	116	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	113	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.026	2.4	111	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.127	2.2	109	0.00
42	S	2,4,6-Tribromophenol	80.000	75.742	5.3	104	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.450	-1.1	112	0.00
44		2,4,5-Trichlorophenol	40.000	40.317	-0.8	113	0.00
45	S	2-Fluorobiphenyl	80.000	80.639	-0.8	114	0.00
46		1,1'-Biphenyl	40.000	40.169	-0.4	115	0.00
47		2-Chloronaphthalene	40.000	39.916	0.2	114	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091522\
Data File : BP011719.D
Acq On : 15 Sep 2022 23:22
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 16 03:01:48 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 14 19:12:50 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	43.842	-9.6	117	0.00
49		Acenaphthylene	40.000	40.903	-2.3	113	0.00
50		Dimethylphthalate	40.000	39.602	1.0	112	0.00
51		2,6-Dinitrotoluene	40.000	41.349	-3.4	112	0.00
52	C	Acenaphthene	40.000	40.255	-0.6	113	-0.01
53		3-Nitroaniline	40.000	38.737	3.2	113	0.00
54	P	2,4-Dinitrophenol	40.000	34.260	14.4	102	0.00
55		Dibenzofuran	40.000	39.765	0.6	113	0.00
56	P	4-Nitrophenol	40.000	39.800	0.5	111	0.00
57		2,4-Dinitrotoluene	40.000	37.804	5.5	111	0.00
58		Fluorene	40.000	40.567	-1.4	114	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	38.770	3.1	108	0.00
60		Diethylphthalate	40.000	40.383	-1.0	113	-0.01
61		4-Chlorophenyl-phenylether	40.000	39.451	1.4	112	0.00
62		4-Nitroaniline	40.000	39.280	1.8	114	0.00
63		Azobenzene	40.000	42.856	-7.1	117	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	110	0.00
65		4,6-Dinitro-2-methylphenol	40.000	35.686	10.8	103	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.360	-3.4	113	0.00
67		4-Bromophenyl-phenylether	40.000	39.064	2.3	107	0.00
68		Hexachlorobenzene	40.000	38.409	4.0	106	0.00
69		Atrazine	40.000	39.832	0.4	107	0.00
70	C	Pentachlorophenol	40.000	38.231	4.4	105	0.00
71		Phenanthrene	40.000	40.182	-0.5	111	0.00
72		Anthracene	40.000	41.269	-3.2	111	0.00
73		Carbazole	40.000	41.124	-2.8	111	0.00
74		Di-n-butylphthalate	40.000	41.612	-4.0	110	-0.01
75	C	Fluoranthene	40.000	39.993	0.0	109	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	106	0.00
77		Benzidine	40.000	38.750	3.1	89	0.00
78		Pyrene	40.000	41.444	-3.6	110	0.00
79	S	Terphenyl-d14	80.000	81.772	-2.2	108	0.00
80		Butylbenzylphthalate	40.000	42.541	-6.4	108	0.00
81		Benzo(a)anthracene	40.000	40.486	-1.2	108	0.00
82		3,3'-Dichlorobenzidine	40.000	41.965	-4.9	106	-0.01
83		Chrysene	40.000	40.225	-0.6	107	-0.01
84		Bis(2-ethylhexyl)phthalate	40.000	43.509	-8.8	107	0.00
85	c	Di-n-octyl phthalate	40.000	41.034	-2.6	106	-0.01
86	I	Perylene-d12	20.000	20.000	0.0	105	-0.01
87		Indeno(1,2,3-cd)pyrene	40.000	41.221	-3.1	106	-0.01
88		Benzo(b)fluoranthene	40.000	40.436	-1.1	104	-0.01
89		Benzo(k)fluoranthene	40.000	41.744	-4.4	108	0.00
90	C	Benzo(a)pyrene	40.000	41.296	-3.2	106	-0.01
91		Dibenzo(a,h)anthracene	40.000	41.299	-3.2	107	-0.02
92		Benzo(g,h,i)perylene	40.000	40.397	-1.0	105	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091522\
Data File : BP011719.D
Acq On : 15 Sep 2022 23:22
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 16 03:01:48 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 14 19:12:50 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