Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032123\
Data File : BM039065.D
Acq On : 21 Mar 2023 12:33
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Mar 22 03:38:21 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM030823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 22 03:36:20 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	113	0.00
2		1,4-Dioxane	40.000	37.514	6.2	103	0.00
3		Pyridine	40.000	37.647	5.9	101	0.00
4		n-Nitrosodimethylamine	40.000	36.861	7.8	101	0.00
5	S	2-Fluorophenol	80.000	77.745	2.8	106	0.00
6		Aniline	40.000	38.758	3.1	106	0.00
7	S	Phenol-d6	80.000	77.009	3.7	106	0.00
8		2-Chlorophenol	40.000	39.683	0.8	109	0.00
9		Benzaldehyde	40.000	35.392	11.5	106	0.00
10	C	Phenol	40.000	38.470	3.8	106	0.00
11		bis(2-Chloroethyl)ether	40.000	38.962	2.6	108	0.00
12		1,3-Dichlorobenzene	40.000	38.799	3.0	106	0.00
13	C	1,4-Dichlorobenzene	40.000	38.827	2.9	107	0.00
14		1,2-Dichlorobenzene	40.000	38.602	3.5	106	0.00
15		Benzyl Alcohol	40.000	39.498	1.3	107	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.709	0.7	111	0.00
17		2-Methylphenol	40.000	39.225	1.9	107	0.00
18		Hexachloroethane	40.000	40.127	-0.3	109	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.463	-1.2	107	0.00
20		3+4-Methylphenols	40.000	39.562	1.1	108	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	114	0.00
22		Acetophenone	40.000	38.272	4.3	105	0.00
23	S	Nitrobenzene-d5	80.000	78.861	1.4	106	0.00
24		Nitrobenzene	40.000	39.065	2.3	106	0.00
25		Isophorone	40.000	39.929	0.2	107	0.00
26	C	2-Nitrophenol	40.000	40.384	-1.0	117	0.00
27		2,4-Dimethylphenol	40.000	39.343	1.6	106	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.174	2.1	107	0.00
29	C	2,4-Dichlorophenol	40.000	39.968	0.1	107	0.00
30		1,2,4-Trichlorobenzene	40.000	38.680	3.3	107	0.00
31		Naphthalene	40.000	38.187	4.5	105	0.00
32		Benzoic acid	40.000	40.519	-1.3	117	0.00
33		4-Chloroaniline	40.000	38.703	3.2	104	0.00
34	C	Hexachlorobutadiene	40.000	39.476	1.3	109	0.00
35		Caprolactam	40.000	38.815	3.0	101	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.890	2.8	104	0.00
37		2-Methylnaphthalene	40.000	38.153	4.6	105	0.00
38		1-Methylnaphthalene	40.000	37.997	5.0	103	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	108	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.128	2.2	103	0.00
41	P	Hexachlorocyclopentadiene	40.000	42.255	-5.6	109	0.00
42	S	2,4,6-Tribromophenol	80.000	78.265	2.2	100	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.980	-2.4	103	0.00
44		2,4,5-Trichlorophenol	40.000	40.342	-0.9	102	0.00
45	S	2-Fluorobiphenyl	80.000	78.131	2.3	102	0.00
46		1,1'-Biphenyl	40.000	38.905	2.7	103	0.00
47		2-Chloronaphthalene	40.000	39.018	2.5	102	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032123\
Data File : BM039065.D
Acq On : 21 Mar 2023 12:33
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Mar 22 03:38:21 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM030823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 22 03:36:20 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	Amount	Calc.	%Dev	Area%	Dev(min)
48		2-Nitroaniline	40.000	39.464	1.3	104	0.00
49		Acenaphthylene	40.000	39.368	1.6	100	0.00
50		Dimethylphthalate	40.000	38.757	3.1	98	0.00
51		2,6-Dinitrotoluene	40.000	38.426	3.9	101	0.00
52	C	Acenaphthene	40.000	38.398	4.0	99	0.00
53		3-Nitroaniline	40.000	38.195	4.5	98	0.00
54	P	2,4-Dinitrophenol	40.000	40.226	-0.6	108	0.00
55		Dibenzofuran	40.000	38.111	4.7	99	0.00
56	P	4-Nitrophenol	40.000	38.813	3.0	95	0.00
57		2,4-Dinitrotoluene	40.000	38.399	4.0	99	0.00
58		Fluorene	40.000	38.211	4.5	97	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.140	2.1	97	0.00
60		Diethylphthalate	40.000	39.788	0.5	98	0.00
61		4-Chlorophenyl-phenylether	40.000	38.698	3.3	99	0.00
62		4-Nitroaniline	40.000	37.708	5.7	95	0.00
63		Azobenzene	40.000	39.624	0.9	97	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
65		4,6-Dinitro-2-methylphenol	40.000	40.156	-0.4	102	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.576	1.1	96	0.00
67		4-Bromophenyl-phenylether	40.000	40.591	-1.5	99	0.00
68		Hexachlorobenzene	40.000	38.739	3.2	96	0.00
69		Atrazine	40.000	42.257	-5.6	97	0.00
70	C	Pentachlorophenol	40.000	39.556	1.1	96	0.00
71		Phenanthrene	40.000	38.449	3.9	94	0.00
72		Anthracene	40.000	39.542	1.1	94	0.00
73		Carbazole	40.000	39.036	2.4	91	0.00
74		Di-n-butylphthalate	40.000	40.469	-1.2	99	0.00
75	C	Fluoranthene	40.000	38.664	3.3	90	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	92	0.00
77		Benzidine	40.000	40.047	-0.1	86	0.00
78		Pyrene	40.000	40.225	-0.6	89	0.00
79	S	Terphenyl-d14	80.000	80.804	-1.0	87	0.00
80		Butylbenzylphthalate	40.000	43.062	-7.7	98	0.00
81		Benzo(a)anthracene	40.000	39.774	0.6	86	0.00
82		3,3'-Dichlorobenzidine	40.000	43.542	-8.9	93	0.00
83		Chrysene	40.000	39.165	2.1	85	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	44.269	-10.7	98	0.00
85	c	Di-n-octyl phthalate	40.000	45.553	-13.9	102	0.00
86	I	Perylene-d12	20.000	20.000	0.0	96	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	39.902	0.2	88	0.00
88		Benzo(b)fluoranthene	40.000	39.238	1.9	87	0.00
89		Benzo(k)fluoranthene	40.000	37.809	5.5	85	0.00
90	C	Benzo(a)pyrene	40.000	39.910	0.2	88	0.00
91		Dibenzo(a,h)anthracene	40.000	39.428	1.4	87	0.00
92		Benzo(g,h,i)perylene	40.000	39.337	1.7	87	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032123\
Data File : BM039065.D
Acq On : 21 Mar 2023 12:33
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Mar 22 03:38:21 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM030823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 22 03:36:20 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	Amount	Calc.	%Dev Area%	Dev(min)

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