Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081721\
Data File : BM031776.D
Acq On : 17 Aug 2021 10:31
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 17 12:17:43 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 16 16:42:38 2021
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	114	0.00
2		1,4-Dioxane	40.000	40.610	-1.5	116	0.00
3		Pyridine	40.000	40.116	-0.3	111	0.00
4		n-Nitrosodimethylamine	40.000	41.616	-4.0	115	0.00
5	S	2-Fluorophenol	80.000	84.278	-5.3	117	0.00
6		Aniline	40.000	40.448	-1.1	111	0.00
7	S	Phenol-d6	80.000	84.009	-5.0	113	0.00
8		2-Chlorophenol	40.000	42.156	-5.4	115	0.00
9		Benzaldehyde	40.000	41.418	-3.5	118	0.00
10	C	Phenol	40.000	42.041	-5.1	114	0.00
11		bis(2-Chloroethyl)ether	40.000	40.895	-2.2	113	0.00
12		1,3-Dichlorobenzene	40.000	41.686	-4.2	114	0.00
13	C	1,4-Dichlorobenzene	40.000	41.375	-3.4	115	0.00
14		1,2-Dichlorobenzene	40.000	41.001	-2.5	116	0.00
15		Benzyl Alcohol	40.000	41.676	-4.2	112	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.883	0.3	110	0.00
17		2-Methylphenol	40.000	41.948	-4.9	112	0.00
18		Hexachloroethane	40.000	42.045	-5.1	114	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.020	-2.6	111	0.00
20		3+4-Methylphenols	40.000	41.799	-4.5	115	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	112	0.00
22		Acetophenone	40.000	42.056	-5.1	114	0.00
23	S	Nitrobenzene-d5	80.000	83.267	-4.1	112	0.00
24		Nitrobenzene	40.000	40.993	-2.5	111	0.00
25		Isophorone	40.000	41.029	-2.6	110	0.00
26	C	2-Nitrophenol	40.000	44.130	-10.3	116	0.00
27		2,4-Dimethylphenol	40.000	41.990	-5.0	113	0.00
28		bis(2-Chloroethoxy)methane	40.000	41.720	-4.3	112	0.00
29	C	2,4-Dichlorophenol	40.000	43.482	-8.7	114	0.00
30		1,2,4-Trichlorobenzene	40.000	43.123	-7.8	117	0.00
31		Naphthalene	40.000	42.307	-5.8	114	0.00
32		Benzoic acid	40.000	37.264	6.8	98	0.00
33		4-Chloroaniline	40.000	42.469	-6.2	113	-0.01
34	C	Hexachlorobutadiene	40.000	43.313	-8.3	116	0.00
35		Caprolactam	40.000	42.764	-6.9	110	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.098	-5.2	110	0.00
37		2-Methylnaphthalene	40.000	42.302	-5.8	113	0.00
38		1-Methylnaphthalene	40.000	42.202	-5.5	113	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	116	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.929	-4.8	115	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.369	1.6	103	0.00
42	S	2,4,6-Tribromophenol	80.000	85.194	-6.5	115	0.00
43	C	2,4,6-Trichlorophenol	40.000	42.472	-6.2	115	0.00
44		2,4,5-Trichlorophenol	40.000	41.751	-4.4	113	0.00
45	S	2-Fluorobiphenyl	80.000	83.223	-4.0	115	0.00
46		1,1'-Biphenyl	40.000	41.306	-3.3	115	0.00
47		2-Chloronaphthalene	40.000	41.501	-3.8	114	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081721\
Data File : BM031776.D
Acq On : 17 Aug 2021 10:31
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Aug 17 12:17:43 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 16 16:42:38 2021
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	41.825	-4.6	112	0.00
49		Acenaphthylene	40.000	41.590	-4.0	114	0.00
50		Dimethylphthalate	40.000	40.221	-0.6	111	0.00
51		2,6-Dinitrotoluene	40.000	41.025	-2.6	112	0.00
52	C	Acenaphthene	40.000	41.338	-3.3	114	0.00
53		3-Nitroaniline	40.000	41.540	-3.8	112	0.00
54	P	2,4-Dinitrophenol	40.000	36.355	9.1	97	0.00
55		Dibenzofuran	40.000	41.452	-3.6	115	0.00
56	P	4-Nitrophenol	40.000	39.101	2.2	103	0.00
57		2,4-Dinitrotoluene	40.000	42.003	-5.0	112	0.00
58		Fluorene	40.000	42.257	-5.6	116	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.469	-3.7	114	0.00
60		Diethylphthalate	40.000	41.322	-3.3	113	0.00
61		4-Chlorophenyl-phenylether	40.000	41.983	-5.0	116	0.00
62		4-Nitroaniline	40.000	41.613	-4.0	111	0.00
63		Azobenzene	40.000	41.422	-3.6	112	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	112	0.00
65		4,6-Dinitro-2-methylphenol	40.000	37.231	6.9	95	0.00
66	c	n-Nitrosodiphenylamine	40.000	42.631	-6.6	113	0.00
67		4-Bromophenyl-phenylether	40.000	42.713	-6.8	113	0.00
68		Hexachlorobenzene	40.000	42.871	-7.2	114	0.00
69		Atrazine	40.000	42.947	-7.4	114	0.00
70	C	Pentachlorophenol	40.000	39.952	0.1	103	0.00
71		Phenanthrene	40.000	41.800	-4.5	112	0.00
72		Anthracene	40.000	42.377	-5.9	113	0.00
73		Carbazole	40.000	41.172	-2.9	108	0.00
74		Di-n-butylphthalate	40.000	42.124	-5.3	111	0.00
75	C	Fluoranthene	40.000	40.515	-1.3	108	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	93	0.00
77		Benzidine	40.000	48.772	-21.9	113	0.00
78		Pyrene	40.000	48.573	-21.4	106	0.00
79	S	Terphenyl-d14	80.000	99.102	-23.9	107	0.00
80		Butylbenzylphthalate	40.000	46.028	-15.1	99	0.00
81		Benzo(a)anthracene	40.000	42.004	-5.0	92	0.00
82		3,3'-Dichlorobenzidine	40.000	40.252	-0.6	86	0.00
83		Chrysene	40.000	41.495	-3.7	92	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	42.231	-5.6	90	0.00
85	c	Di-n-octyl phthalate	40.000	38.060	4.8	83	0.00
86	I	Perylene-d12	20.000	20.000	0.0	80	-0.01
87		Indeno(1,2,3-cd)pyrene	40.000	42.542	-6.4	81	0.00
88		Benzo(b)fluoranthene	40.000	43.898	-9.7	82	-0.01
89		Benzo(k)fluoranthene	40.000	41.855	-4.6	80	0.00
90	C	Benzo(a)pyrene	40.000	42.256	-5.6	79	-0.01
91		Dibenzo(a,h)anthracene	40.000	42.187	-5.5	81	-0.01
92		Benzo(g,h,i)perylene	40.000	43.485	-8.7	84	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081721\
Data File : BM031776.D
Acq On : 17 Aug 2021 10:31
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Aug 17 12:17:43 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 16 16:42:38 2021
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