Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011822\
Data File : BM033774.D
Acq On : 18 Jan 2022 21:20
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Jan 19 04:22:47 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM011722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 17 16:30:56 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	118	0.00
2		1,4-Dioxane	40.000	38.023	4.9	125	0.00
3		Pyridine	40.000	38.169	4.6	120	0.00
4		n-Nitrosodimethylamine	40.000	37.775	5.6	118	0.00
5	S	2-Fluorophenol	80.000	75.242	5.9	120	0.00
6		Aniline	40.000	37.568	6.1	118	0.00
7	S	Phenol-d6	80.000	75.700	5.4	118	0.00
8		2-Chlorophenol	40.000	37.039	7.4	117	0.00
9		Benzaldehyde	40.000	41.393	-3.5	121	0.00
10	C	Phenol	40.000	38.259	4.4	119	0.00
11		bis(2-Chloroethyl)ether	40.000	36.880	7.8	118	0.00
12		1,3-Dichlorobenzene	40.000	36.455	8.9	117	0.00
13	C	1,4-Dichlorobenzene	40.000	36.531	8.7	117	0.00
14		1,2-Dichlorobenzene	40.000	36.484	8.8	117	0.00
15		Benzyl Alcohol	40.000	37.004	7.5	116	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.933	2.7	124	-0.01
17		2-Methylphenol	40.000	37.297	6.8	116	0.00
18		Hexachloroethane	40.000	36.712	8.2	119	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.902	7.7	117	0.00
20		3+4-Methylphenols	40.000	37.656	5.9	117	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	116	0.00
22		Acetophenone	40.000	36.181	9.5	116	0.00
23	S	Nitrobenzene-d5	80.000	72.474	9.4	116	0.00
24		Nitrobenzene	40.000	36.488	8.8	117	0.00
25		Isophorone	40.000	36.146	9.6	116	0.00
26	C	2-Nitrophenol	40.000	36.516	8.7	114	0.00
27		2,4-Dimethylphenol	40.000	35.908	10.2	114	0.00
28		bis(2-Chloroethoxy)methane	40.000	36.166	9.6	118	0.00
29	C	2,4-Dichlorophenol	40.000	35.759	10.6	114	0.00
30		1,2,4-Trichlorobenzene	40.000	35.365	11.6	114	0.00
31		Naphthalene	40.000	35.750	10.6	115	0.00
32		Benzoic acid	40.000	39.146	2.1	117	0.00
33		4-Chloroaniline	40.000	36.266	9.3	115	0.00
34	C	Hexachlorobutadiene	40.000	34.682	13.3	112	0.00
35		Caprolactam	40.000	36.815	8.0	117	0.00
36	C	4-Chloro-3-methylphenol	40.000	36.253	9.4	116	0.00
37		2-Methylnaphthalene	40.000	35.082	12.3	114	0.00
38		1-Methylnaphthalene	40.000	35.215	12.0	114	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	115	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	34.775	13.1	112	0.00
41	P	Hexachlorocyclopentadiene	40.000	34.334	14.2	107	0.00
42	S	2,4,6-Tribromophenol	80.000	68.639	14.2	109	0.00
43	C	2,4,6-Trichlorophenol	40.000	35.854	10.4	114	0.00
44		2,4,5-Trichlorophenol	40.000	36.138	9.7	115	0.00
45	S	2-Fluorobiphenyl	80.000	69.616	13.0	113	0.00
46		1,1'-Biphenyl	40.000	35.381	11.5	114	0.00
47		2-Chloronaphthalene	40.000	35.679	10.8	115	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011822\
Data File : BM033774.D
Acq On : 18 Jan 2022 21:20
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Jan 19 04:22:47 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM011722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 17 16:30:56 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	38.069	4.8	117	0.00
49		Acenaphthylene	40.000	35.813	10.5	115	0.00
50		Dimethylphthalate	40.000	35.415	11.5	115	0.00
51		2,6-Dinitrotoluene	40.000	36.526	8.7	116	0.00
52	C	Acenaphthene	40.000	36.083	9.8	115	0.00
53		3-Nitroaniline	40.000	36.997	7.5	116	0.00
54	P	2,4-Dinitrophenol	40.000	37.288	6.8	112	0.00
55		Dibenzofuran	40.000	35.257	11.9	115	0.00
56	P	4-Nitrophenol	40.000	38.807	3.0	120	-0.01
57		2,4-Dinitrotoluene	40.000	37.661	5.8	118	0.00
58		Fluorene	40.000	34.831	12.9	113	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	35.568	11.1	114	0.00
60		Diethylphthalate	40.000	35.811	10.5	115	0.00
61		4-Chlorophenyl-phenylether	40.000	34.909	12.7	114	0.00
62		4-Nitroaniline	40.000	37.992	5.0	118	0.00
63		Azobenzene	40.000	36.958	7.6	118	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	117	0.00
65		4,6-Dinitro-2-methylphenol	40.000	36.962	7.6	113	0.00
66	c	n-Nitrosodiphenylamine	40.000	35.217	12.0	115	0.00
67		4-Bromophenyl-phenylether	40.000	33.671	15.8	111	0.00
68		Hexachlorobenzene	40.000	33.925	15.2	113	0.00
69		Atrazine	40.000	35.037	12.4	114	-0.01
70	C	Pentachlorophenol	40.000	36.185	9.5	114	-0.01
71		Phenanthrene	40.000	35.077	12.3	116	0.00
72		Anthracene	40.000	35.480	11.3	116	0.00
73		Carbazole	40.000	36.198	9.5	118	0.00
74		Di-n-butylphthalate	40.000	36.345	9.1	117	-0.01
75	C	Fluoranthene	40.000	36.382	9.0	119	-0.01
76	I	Chrysene-d12	20.000	20.000	0.0	132	0.00
77		Benzidine	40.000	37.576	6.1	141	0.00
78		Pyrene	40.000	32.922	17.7	119	0.00
79	S	Terphenyl-d14	80.000	65.509	18.1	119	0.00
80		Butylbenzylphthalate	40.000	35.602	11.0	127	-0.01
81		Benzo(a)anthracene	40.000	36.117	9.7	131	-0.01
82		3,3'-Dichlorobenzidine	40.000	35.957	10.1	132	0.00
83		Chrysene	40.000	36.066	9.8	133	-0.01
84		Bis(2-ethylhexyl)phthalate	40.000	35.860	10.4	130	-0.01
85	c	Di-n-octyl phthalate	40.000	38.701	3.2	142	-0.01
86	I	Perylene-d12	20.000	20.000	0.0	148	-0.01
87		Indeno(1,2,3-cd)pyrene	40.000	34.440	13.9	143	-0.02
88		Benzo(b)fluoranthene	40.000	36.592	8.5	148	-0.01
89		Benzo(k)fluoranthene	40.000	36.419	9.0	147	-0.01
90	C	Benzo(a)pyrene	40.000	36.442	8.9	147	-0.01
91		Dibenzo(a,h)anthracene	40.000	34.577	13.6	142	-0.02
92		Benzo(g,h,i)perylene	40.000	33.921	15.2	140	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011822\
Data File : BM033774.D
Acq On : 18 Jan 2022 21:20
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Jan 19 04:22:47 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM011722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 17 16:30:56 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