Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012720\
Data File : BM024625.D
Acq On : 27 Jan 2020 13:20
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jan 28 05:38:06 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 23 14:57:07 2020
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	115	0.00
2		1,4-Dioxane	0.416	0.447	-7.5	122	0.00
3		Pyridine	1.206	1.181	2.1	113	0.00
4		n-Nitrosodimethylamine	0.535	0.479	10.5	100	0.00
5	S	2-Fluorophenol	1.052	1.116	-6.1	119	0.00
6		Aniline	1.764	1.787	-1.3	113	0.00
7	S	Phenol-d6	1.449	1.467	-1.2	113	0.00
8		2-Chlorophenol	1.197	1.280	-6.9	119	0.00
9		Benzaldehyde	0.856	0.764	10.7	103	0.00
10	C	Phenol	1.444	1.447	-0.2	111	0.00
11		bis(2-Chloroethyl)ether	1.156	1.169	-1.1	111	0.00
12		1,3-Dichlorobenzene	1.471	1.651	-12.2	125	0.00
13	C	1,4-Dichlorobenzene	1.492	1.657	-11.1	124	0.00
14		1,2-Dichlorobenzene	1.440	1.581	-9.8	123	0.00
15		Benzyl Alcohol	1.194	1.151	3.6	108	0.00
16		2,2'-oxybis(1-Chloropropane	1.058	0.931	12.0	96	0.00
17		2-Methylphenol	1.018	1.050	-3.1	114	0.00
18		Hexachloroethane	0.576	0.603	-4.7	117	0.00
19	P	n-Nitroso-di-n-propylamine	1.088	0.995	8.5	102	0.00
20		3+4-Methylphenols	1.409	1.436	-1.9	112	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.516	0.536	-3.9	110	0.00
23	S	Nitrobenzene-d5	0.408	0.404	1.0	104	0.00
24		Nitrobenzene	0.391	0.387	1.0	105	0.00
25		Isophorone	0.693	0.689	0.6	105	0.00
26	C	2-Nitrophenol	0.181	0.204	-12.7	118	0.00
27		2,4-Dimethylphenol	0.246	0.268	-8.9	116	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.443	-4.7	110	0.00
29	C	2,4-Dichlorophenol	0.332	0.389	-17.2	123	0.00
30		1,2,4-Trichlorobenzene	0.401	0.474	-18.2	127	0.00
31		Naphthalene	1.010	1.110	-9.9	117	0.00
32		Benzoic acid	0.225	0.254	-12.9	116	0.00
33		4-Chloroaniline	0.443	0.491	-10.8	116	0.00
34	C	Hexachlorobutadiene	0.277	0.328	-18.4	125	0.00
35		Caprolactam	0.107	0.117	-9.3	115	0.00
36	C	4-Chloro-3-methylphenol	0.353	0.372	-5.4	111	0.00
37		2-Methylnaphthalene	0.780	0.871	-11.7	119	0.00
38		1-Methylnaphthalene	0.742	0.828	-11.6	119	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	113	0.00
40		1,2,4,5-Tetrachlorobenzene	0.714	0.803	-12.5	125	0.00
41	P	Hexachlorocyclopentadiene	0.400	0.466	-16.5	127	0.00
42	S	2,4,6-Tribromophenol	0.325	0.385	-18.5	133	0.00
43	C	2,4,6-Trichlorophenol	0.454	0.512	-12.8	123	0.00
44		2,4,5-Trichlorophenol	0.464	0.526	-13.4	125	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012720\
Data File : BM024625.D
Acq On : 27 Jan 2020 13:20
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jan 28 05:38:06 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 23 14:57:07 2020
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.471	1.619	-10.1	121	0.00
46		1,1'-Biphenyl	1.514	1.670	-10.3	122	0.00
47		2-Chloronaphthalene	1.222	1.348	-10.3	122	0.00
48		2-Nitroaniline	0.373	0.346	7.2	101	0.00
49		Acenaphthylene	1.829	2.009	-9.8	121	0.00
50		Dimethylphthalate	1.617	1.786	-10.5	122	0.00
51		2,6-Dinitrotoluene	0.344	0.389	-13.1	124	0.00
52	C	Acenaphthene	1.135	1.241	-9.3	121	0.00
53		3-Nitroaniline	0.357	0.386	-8.1	117	0.00
54	P	2,4-Dinitrophenol	0.206	0.236	-14.6	126	0.00
55		Dibenzofuran	1.851	2.059	-11.2	124	0.00
56	P	4-Nitrophenol	0.279	0.304	-9.0	118	0.00
57		2,4-Dinitrotoluene	0.493	0.554	-12.4	123	0.00
58		Fluorene	1.556	1.696	-9.0	120	0.00
59		2,3,4,6-Tetrachlorophenol	0.470	0.533	-13.4	126	0.00
60		Diethylphthalate	1.642	1.760	-7.2	118	0.00
61		4-Chlorophenyl-phenylether	0.897	0.975	-8.7	121	0.00
62		4-Nitroaniline	0.394	0.423	-7.4	116	0.00
63		Azobenzene	1.401	1.327	5.3	104	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	113	0.00
65		4,6-Dinitro-2-methylphenol	0.118	0.135	-14.4	126	0.00
66	c	n-Nitrosodiphenylamine	0.574	0.636	-10.8	124	0.00
67		4-Bromophenyl-phenylether	0.249	0.285	-14.5	128	0.00
68		Hexachlorobenzene	0.285	0.337	-18.2	133	0.00
69		Atrazine	0.221	0.224	-1.4	111	0.00
70	C	Pentachlorophenol	0.174	0.201	-15.5	128	0.00
71		Phenanthrene	1.083	1.201	-10.9	124	0.00
72		Anthracene	1.060	1.182	-11.5	125	0.00
73		Carbazole	0.965	1.071	-11.0	123	0.00
74		Di-n-butylphthalate	1.233	1.308	-6.1	117	0.00
75	C	Fluoranthene	1.352	1.505	-11.3	125	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	111	0.00
77		Benzidine	0.466	0.618	-32.6#	143	0.00
78		Pyrene	1.318	1.523	-15.6	125	0.00
79	S	Terphenyl-d14	1.071	1.263	-17.9	119	0.00
80		Butylbenzylphthalate	0.553	0.592	-7.1	114	0.00
81		Benzo(a)anthracene	1.388	1.530	-10.2	119	0.00
82		3,3'-Dichlorobenzidine	0.499	0.571	-14.4	119	0.00
83		Chrysene	1.325	1.458	-10.0	119	0.00
84		Bis(2-ethylhexyl)phthalate	0.816	0.852	-4.4	112	0.00
85	c	Di-n-octyl phthalate	1.349	1.366	-1.3	107	0.00
86		Indeno(1,2,3-cd)pyrene	1.444	1.401	3.0	101	0.00
87	I	Perylene-d12	1.000	1.000	0.0	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012720\
Data File : BM024625.D
Acq On : 27 Jan 2020 13:20
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Jan 28 05:38:06 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM012320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 23 14:57:07 2020
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.297	1.479	-14.0	112	0.00
89	Benzo(k)fluoranthene	1.266	1.443	-14.0	114	0.00
90 C	Benzo(a)pyrene	1.179	1.326	-12.5	111	0.00
91	Dibenzo(a,h)anthracene	1.119	1.164	-4.0	101	0.00
92	Benzo(g,h,i)perylene	1.016	1.090	-7.3	103	0.00

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