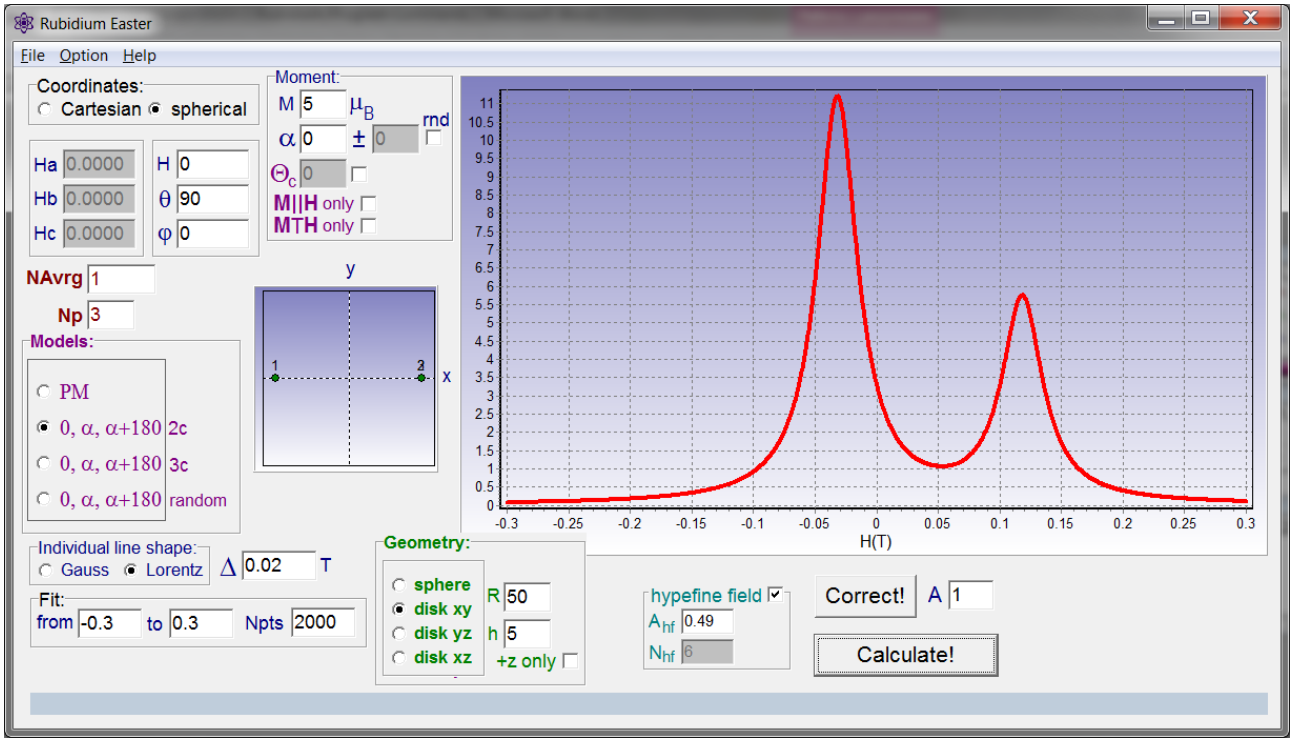
Rubidium Program Comments

The program calculates rubidium NMR spectra in $Rb_{1-x}K_xFe(MoO_4)_2$ for magnetic structures discussed in the paper https://doi.org/10.1103/PhysRevB.99.024419. The following screenshots show parameter sets used in our final calculations (Fig. 6 of the paper). If desired, you can try using other parameter sets to check. The value of the contact field 0.49 arb. units corresponds to 2.5 mT/μB stated in the paper. θ = 90° sets the field $H \perp C^3$, for this field direction the spectral shape does not depend on the value of φ.

Fig.6. (a) 3D *uud* with periodicity 2*c* or 3*c* in the C^3 direction, μ = 5 μB, individual linewidth δ = 20 mT.



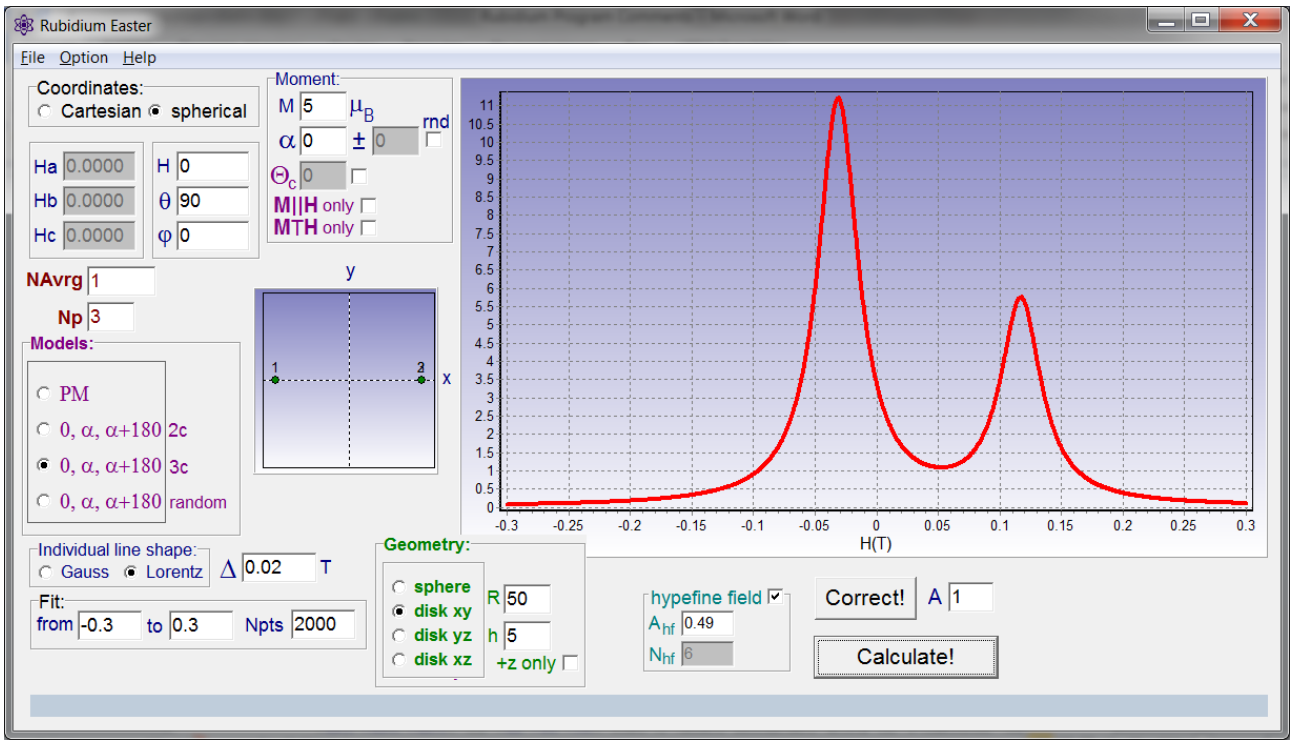
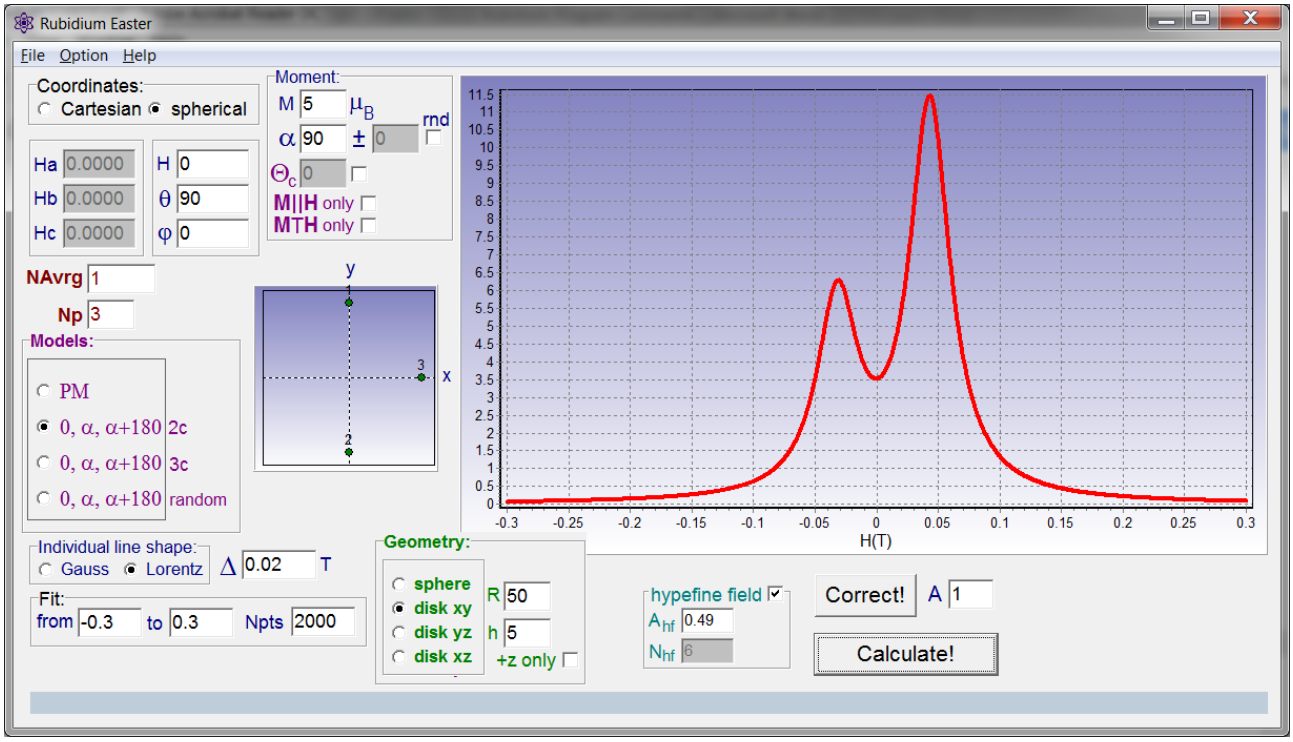


Fig.6. (b) 3D *fan* with periodicity 2*c* or 3*c* in the C^3 direction, μ = 5 μB, δ = 20 mT.



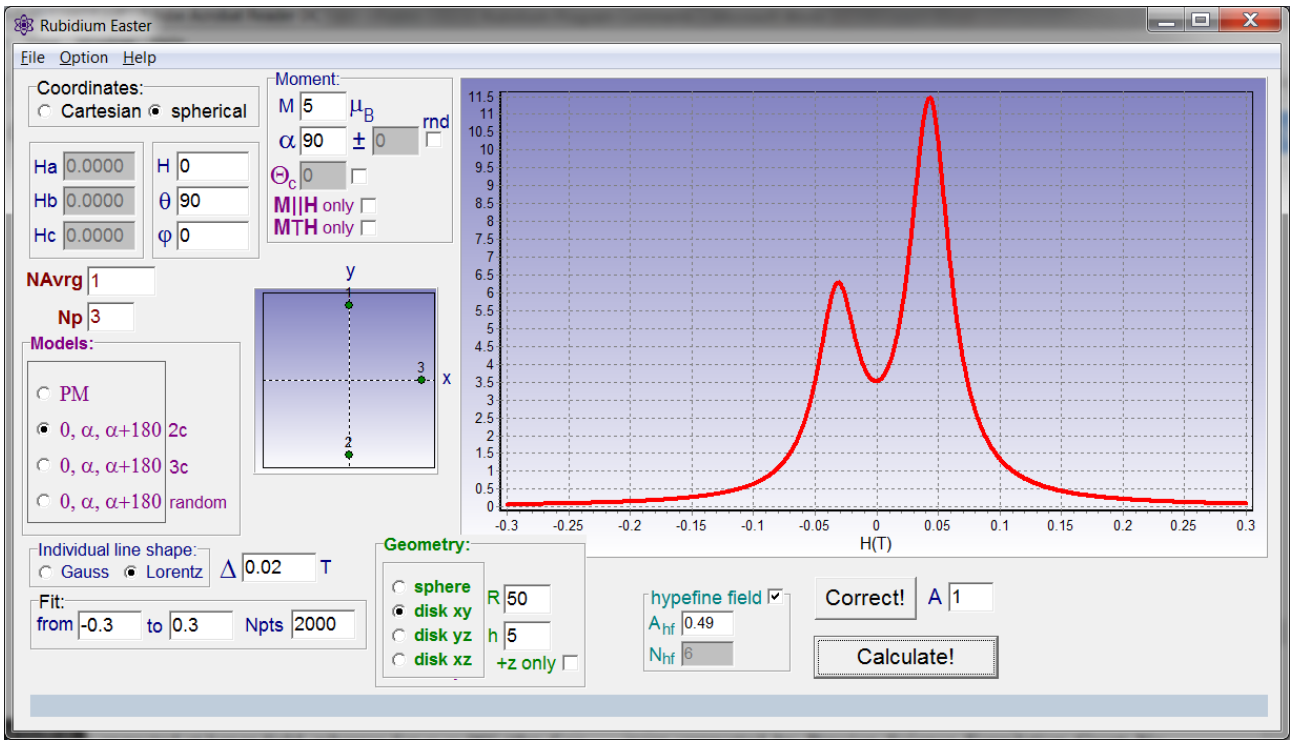


Fig.6. (c) 2D *uud*, μ = 5 μB, δ = 20 mT.

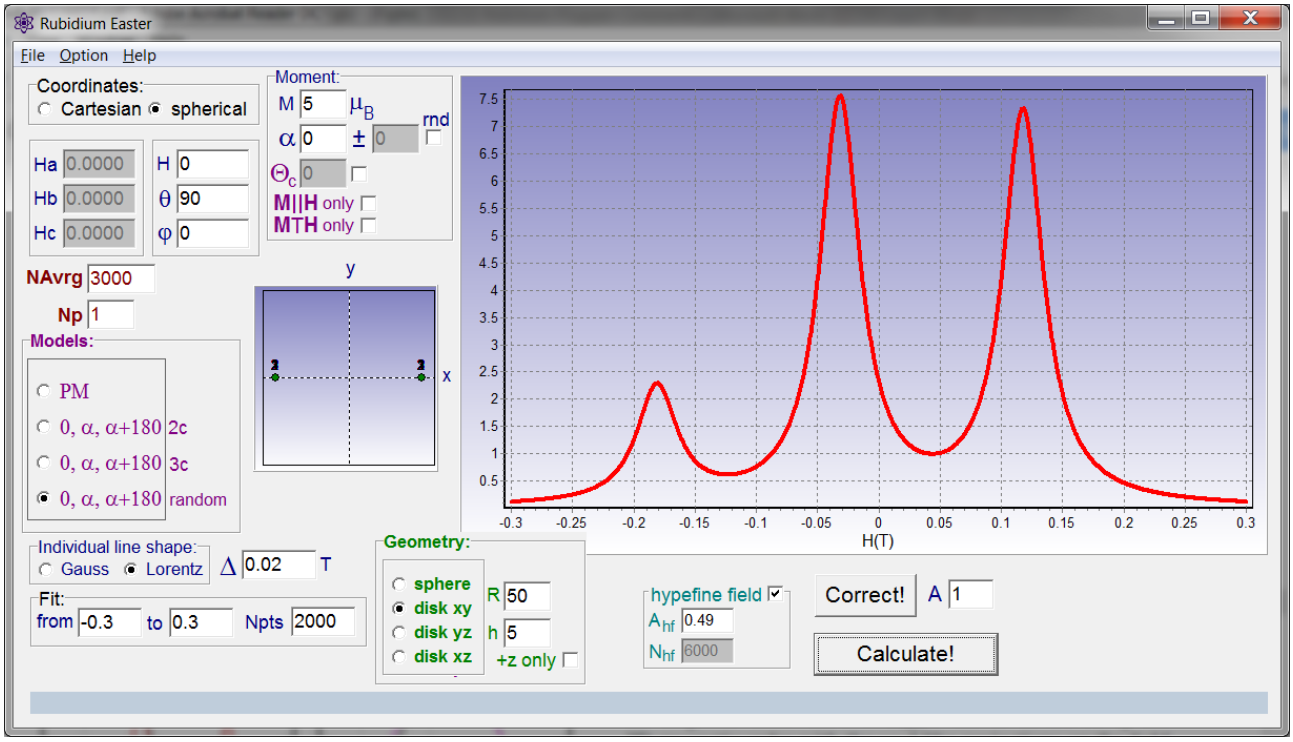
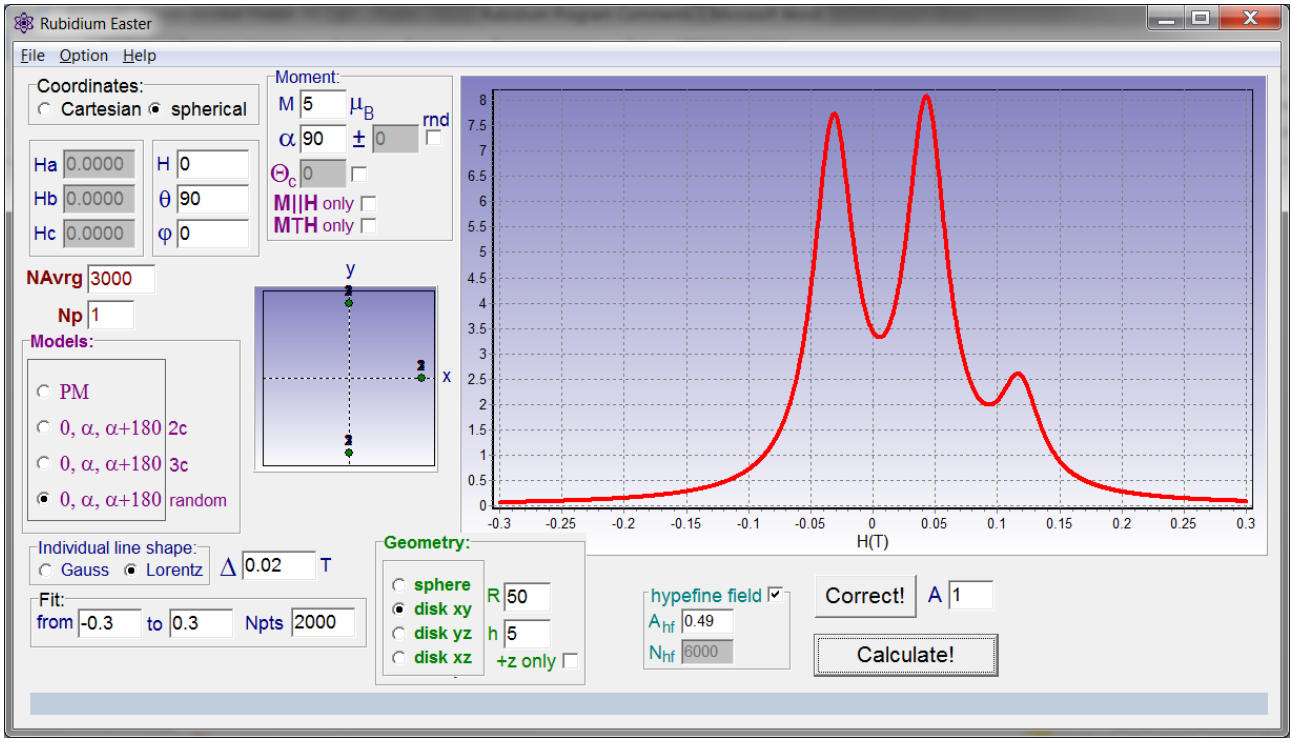
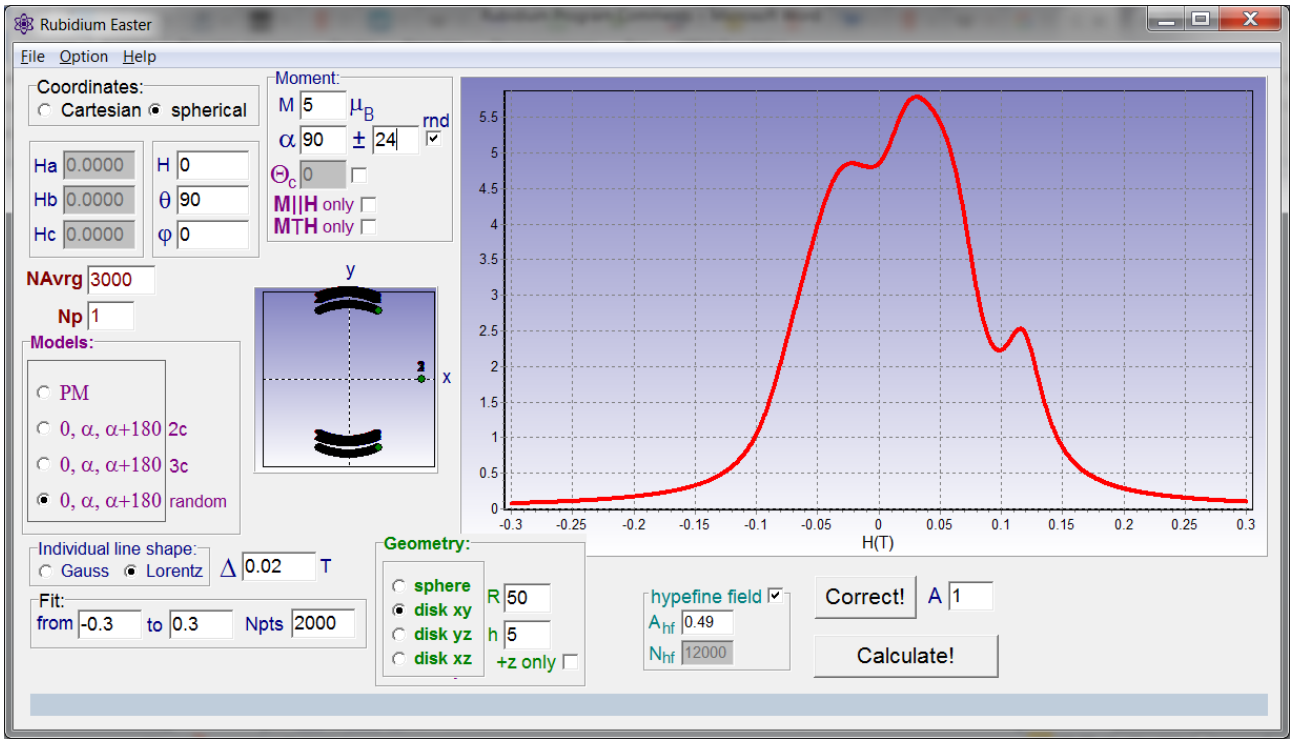


Fig. 6. (d) 2D *fan*, μ = 5 μB, α = 90°, δ = 20 mT,



2D *fan*, μ = 5 μB, α = 90°±24°, δ = 20 mT.



Questions, suggestions and bug reports are welcomed!

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