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ACROPOLIS: A GeneriC FRamework fOr Photodisintegration Of Light ElementS

P. F. Depta, K. Schmidt-Hoberg

Deutsches Elektronen-Synchrotron DESY, Hamburg

M. Hufnagel

Deutsches Elektronen-Synchrotron DESY, Hamburg and

Service de Physique Théorique, Université Libre de Bruxelles, Belgium

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NOTKESTRASSE 85 - 22607 HAMBURG

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Paul Frederik Depta^a, Marco Hufnagel^{a b}, and Kai Schmidt-Hoberg^a

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Abstract:

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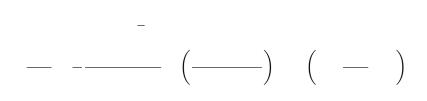
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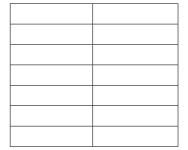
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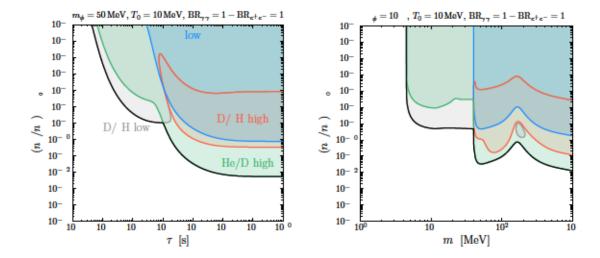
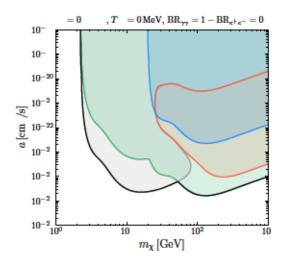


Figure 1. 95% C.L. constraints for decay of a decoupled MeV-scale BSM particle (implemented in DecayModel) into two photons (BR $_{\gamma\gamma}=1$ – BR $_{e^+e^-}=1$) in $\tau_{\phi}-(n_{\phi}/n_{\gamma})|_{T=T_0}$ plane (left) and $m_{\phi}-(n_{\phi}/n_{\gamma})|_{T=T_0}$ plane (right) with $T_0=10\,\mathrm{MeV}$ and $m_{\phi}=50\,\mathrm{MeV}$ (left) as well as $\tau_{\phi}=10^7\,\mathrm{s}$ (right). The limits from individual observables are shown separately: primordial deuterium abundance (orange high, grey low), helium-4 mass fraction \mathcal{Y}_p (blue), and helium-3 abundance normalised by deuterium (green). The overall 95% C.L. BBN limit is given by the black full line as an envelope of individual 95% C.L. constraints neglecting correlations. Using $(n_{\phi}/n_{\gamma})|_{T=T_0}$, i.e. n0a, as a fast parameter on a single computing node with two AMD EPYC 7402 24-Core Processors the scans took $\sim 40\,\mathrm{min}$ (left) and $\sim 2\,\mathrm{h}$ (right) for a 200 \times 200 grid.

deuterium, $m_{\phi} = 2E_{\rm D}^{\rm th} \approx 4.4\,{\rm MeV}$. Apart from some regions with more complex structure due to different disintegration reactions the limits become increasingly strong with larger m_{ϕ} as the energy density injected into the SM becomes larger.

The scans for figure 1 took \sim 40 min (left) and \sim 2 h (right) for a 200 × 200 grid on an AMD EPYC 7402 24-Core Processors, clearly highlighting the performance improvement due to the fast parameter $(n_{\phi}/n_{\gamma})|_{T=T_0}$, i.e. n0a, making the number of points in this direction computationally inexpensive (cf. also appendix B). The runtime is thus determined mostly by the number of points in the direction of τ_{ϕ} or m_{ϕ} (not fast). Note that the longer runtime for the right panel is a result of the database files for the electromagnetic cascade reaction rates having an upper limit on the energy of $m_{\phi}/2 = E_0 = 100\,\text{MeV}$, which often corresponds to the most interesting region in parameter space. For masses above the pion threshold in particular, $m_{\phi} \gtrsim 280\,\text{MeV}$, hadrodisintegration may become relevant if ϕ has non-vanishing couplings to quarks, implying that $\text{BR}_{\gamma\gamma} + \text{BR}_{e^+e^-} < 1$ in general. Also muons are kinematically available in the mass region (which are currently not implemented in ACROPOLIS).

In figure 2 we show the constraints for residual annihilations of DM into two photons $(BR_{\gamma\gamma} = 1 - BR_{e^+e^-} = 1)$ for purely s-wave annihilations (left, b = 0) and purely p-wave annihilations (right, a = 0, $T_{\rm kd} = 1 \, {\rm MeV}$) as implemented in AnnihilationModel. These limits start at the disintegration threshold of deuterium, $m_{\chi} = E_{\rm D}^{\rm th} \approx 2.2 \, {\rm MeV}$, and closely



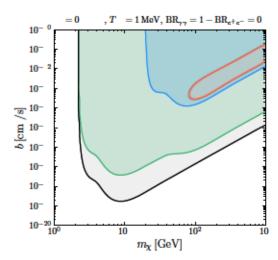


Figure 2. 95% C.L. constraints for residual annihilations of DM (implemented in AnnihilationModel) into two photons (BR $_{\gamma\gamma}=1-\mathrm{BR}_{e^+e^-}=1$) for purely s-wave annihilations (left, b=0) and purely p-wave annihilations (right, a=0, $T_{\mathrm{kd}}=1\,\mathrm{MeV}$). For the explanation of the colour-coding see figure 1. Using $(n_\phi/n_\gamma)|_{T=T_0}$, i.e. n0a, as a fast parameter on a single computing node with two AMD EPYC 7402 24-Core Processors the scans took ~ 10 h each for a 200×200 grid.

resemble those presented in [26], albeit with updated observationally inferred primordial abundances. We therefore refer to [26] for a detailed discussion. The scans took ~ 10 h each for a 200×200 grid on the aforementioned computing node.

6 Implementing your own models

6.1 The model framework acropolis.models

While the provided example models should suffice to tackle most problems of interest, it may sometimes still happen that a scenario cannot directly be mapped to the standard implementation in ACROPOLIS. For such cases, ACROPOLIS provides further tools that allow for an easy implementation of additional models. The most important class in this context is acropolis.models.AbstractModel, which is an abstract base class containing most of the low-level implementation needed to run its method run_disintegration(). In fact, using this class as a base, any new model can be implemented in only two steps:

- create a new class, say NewModel, that uses AbstractModel as a base class, and
- (ii) implement all abstract methods that are provided by AbstractModel, i.e.
 - AbstractModel._temperature_range()

⁷By default, the function AbstractModel._source_positron() simply returns the output of AbstractModel._source_electron(), which is justified for most scenarios. However, if your specific scenario predicts different source terms for electrons and positrons, it is always possible to simply overwrite the former function.

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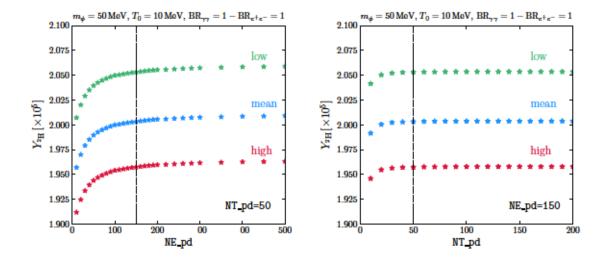


Figure 3. Convergence of the abundance of deuterium as a function of the grid points NE_pd and NT_pd. The dashed line indicates the default value in ACROPOLIS.

Acknowledgments

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A Rates for the cascade processes

In this appendix, we collect for completeness all relevant total and differential interaction rates $\Gamma_x(E)$ and $K_{x'\to x}(E,E')$ for the cascade processes of high-energetic photons, electrons, and positrons on the background photons, electrons, and nuclei (see eqs. (2.4) and (2.7)). Large parts are directly taken from [24].

Target densities

The thermal photon spectrum differential in energy $f_{\gamma}(\bar{\epsilon})$ is given by

$$f_{\gamma}(\bar{\epsilon}) = \frac{\bar{\epsilon}^2}{\pi^2} \times \frac{1}{\exp(\bar{\epsilon}/T) - 1}$$
, (A.1)

while the total baryon number density can be calculated from the baryon-to-photon ratio η and the number density of photons $n_{\gamma}(T)$,

$$n_b(T) = \eta \times n_\gamma(T) = \eta \times \frac{2\zeta(3)}{\pi^2} T^3$$
 (A.2)

Via charge neutrality we obtain for the number density of background electrons

$$n_e(T) = \sum_N Z_N n_N \simeq [Y_p(T) + 2Y_{4He}(T)] \times n_b(T), \quad Y_N(T) = \frac{n_N(T)}{n_b(T)}.$$
 (A.3)

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