Implementation of non-trivial species in CLASS

decaying dark matter, scalar fields, ...

Expanding CLASS

- Adding exotic new physics to CLASS is easy*
- New features will be merged into the main code and will be available in all future versions
- This lecture is essentially a complete step by step tutorial for adding decaying dark matter at the level of the code.

*given sufficient understanding of the underlying physics and provided that the rules of CLASS are respected...

A word on quintessence

- Implemented in a CLASS branch by Miguel Zumalacárregui (ITP - Uni. of Heidelberg)
- Merged into the development version of CLASS this week, available in v2.4

$$\frac{d\phi^2}{dt^2} = -3H\frac{d\phi}{dt} - \frac{dV}{d\phi},$$

$$V(\phi) = M^4[(\phi - B)^{\alpha} + A]e^{-\lambda\phi/M}$$

Changes similar to DCDM case

input.c

Do we have to deal with boundary conditions?
 (See Julien's lecture 1)

```
int input init(...){
 /** These two arrays must contain the strings of names to be searched
     for and the coresponding new parameter */
  char * const target namestrings[] =
    {"100*theta_s", "Omega_dcdmdr", "omega_dcdmdr", "Omega_scf"};
  char * const unknown namestrings[] =
    {"h","Omega_ini_dcdm","Omega_ini_dcdm","scf_lambda"};
  enum computation_stage target_cs[] =
    {cs_thermodynamics, cs_background, cs_background};
int input try unknown parameters()
int input get guess()
```

Setting flags, defining _bg_ and _bi_ indices:

```
int background_indices(...){
    if (pba->Omega0_dcdmdr != 0.){
        pba->has_dcdm = _TRUE_;
        pba->has_dr = _TRUE_;
}
    ...
    /* - index for dcdm */
    class_define_index(pba->index_bg_rho_dcdm,pba->has_dcdm,index_bg,1);
    ...
    /* -> energy density in DCDM */
    class_define_index(pba->index_bi_rho_dcdm,pba->has_dcdm,index_bi,1);
    ...
}
```

Compute {A} variables from {B} variables:

```
int background_functions(...){
    ...
    /* dcdm */
    if (pba->has_dcdm == _TRUE_) {
        /* Pass value of rho_dcdm to output */
        pvecback[pba->index_bg_rho_dcdm] = pvecback_B[pba->index_bi_rho_dcdm];
        rho_tot += pvecback[pba->index_bg_rho_dcdm];
        p_tot += 0.;
        rho_m += pvecback[pba->index_bg_rho_dcdm];
    }
    ...
}
```

Set initial conditions for {B} quantities:

Write e.o.m for {B} quantities:

```
int background_derivs(...){
    ...
    if (pba->has_dcdm == _TRUE_){
        /** compute dcdm density rho' = -3aH rho - a Gamma rho*/
        dy[pba->index_bi_rho_dcdm] = -3.*y[pba->index_bi_a]*
            pvecback[pba->index_bg_H]*y[pba->index_bi_rho_dcdm]-
            y[pba->index_bi_a]*pba->Gamma_dcdm*y[pba->index_bi_rho_dcdm];
    }
    ...
}
```

Initialise perturbation vector:

Set initial conditions for perturbations.

```
int perturb_initial_conditions(...){
  if (pba->has_dcdm == _TRUE_)
    rho_m += ppw->pvecback[pba->index_bg_rho_dcdm];
}
```

Differential equation for perturbations:

```
int perturb_derivs(...){
  if ((pba->has_dcdm == _TRUE_)&&(pba->has_dr == _TRUE_)) {
    /** -> dcdm */
    dy[pv->index_pt_delta_dcdm] =
        -(y[pv->index_pt_theta_dcdm]+metric_continuity);

    dy[pv->index_pt_theta_dcdm] =
        - a_prime_over_a*y[pv->index_pt_theta_dcdm] + metric_euler;
    }
}
```

• Add contribution to $\delta T^{\mu,\nu}$:

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```
int perturb_total_stress_energy(...){
  /* dcdm contribution */
  if (pba->has dcdm == TRUE ) {
    ppw->delta rho += ppw->pvecback[pba->index bg rho dcdm]*
                        y[ppw->pv->index pt delta dcdm];
    ppw->rho plus p theta += ppw->pvecback[pba->index bg rho dcdm]*
                                y[ppw->pv->index pt theta dcdm];
  if (ppt->has_source_delta_m == _TRUE_) {
    /* include decaying cold dark matter */
    if (pba->has_dcdm == _TRUE_) {
      delta_rho_m += ppw->pvecback[pba->index_bg_rho_dcdm]*
        y[ppw->pv->index pt delta dcdm];
      rho_m += ppw->pvecback[pba->index_bg_rho_dcdm];
Priday, 16 May 2014
```

 Print perturbations? (Note: will be moved to output.c in an upcoming release.)

• Store source functions $S_i(k,\tau)$ for later use?

```
int perturb_indices_of_perturbs(...){
    ppt->has_source_delta_dcdm = _FALSE_;
    ppt->has_source_theta_dcdm = _FALSE_;
    if (ppt->has_density_transfers == _TRUE_) {
        if (pba->has_dcdm == _TRUE_)
            ppt->has_source_delta_dcdm = _TRUE_;
    }
    if (ppt->has_velocity_transfers == _TRUE_) {
        if (pba->has_dcdm == _TRUE_)
            ppt->has_source_theta_dcdm = _TRUE_;
    }
    class_define_index(ppt->index_tp_delta_dcdm, ppt->has_source_delta_dcdm,...);
    class_define_index(ppt->index_tp_theta_dcdm, ppt->has_source_theta_dcdm,...);
}
```

• Store the sources $S_i(\tau, k)$:

```
int perturb_sources(...){
    /* delta_dcdm */
    if (ppt->has_source_delta_dcdm == _TRUE_) {
        _set_source_(ppt->index_tp_delta_dcdm) = y[ppw->pv->index_pt_delta_dcdm];
    }
    /* theta_dcdm */
    if (ppt->has_source_theta_dcdm == _TRUE_) {
        _set_source_(ppt->index_tp_theta_dcdm) = y[ppw->pv->index_pt_theta_dcdm];
    }
}
```

spectra.c

• Define indices for transfer functions $T_i(k,z)$:

```
int spectra_indices(...){
    /* indices for species ass. with a matter trsf. function in Fourier space*/
    class_define_index(psp->index_tr_delta_dcdm,ppt->has_source_delta_dcdm,...);
    /* indices for species ass. with a vel. trsf. function in Fourier space */
    class_define_index(psp->index_tr_theta_dcdm,ppt->has_source_theta_dcdm,...);
}
```

spectra.c

• Compute matter transfer functions $T_i(k, z)$:

```
int spectra_matter_transfers(...){
  /* T dcdm(k,tau) */
  if (pba->has_dcdm == _TRUE_) {
    rho_i = pvecback_sp_long[pba->index_bg_rho_dcdm];
      if (ppt->has source delta dcdm == TRUE ) {
        delta_i = ppt->sources[][(...)+ppt->index_tp_delta_dcdm][(...)+index_k];
        psp->matter_transfer[(...) + psp->index_tr_delta_dcdm] = delta_i;
        delta rho tot += rho i * delta i;
        rho tot += rho i;
      if (ppt->has source theta dcdm == TRUE ) {
        theta_i = ppt->sources[][(...)+ppt->index_tp_theta_dcdm][(...)+index_k];
        psp->matter transfer[(...) + psp->index tr theta dcdm] = theta i;
        rho_plus_p_theta_tot += rho_i * theta_i;
        rho plus p tot += rho i;
```

output.c

Write transfer functions if requested: