

# Traceurs et isotopes dans LDMZ6

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# Nouveau fichier **tracer.def** - objectifs

- Plus de souplesse tout en gardant une bonne lisibilité  $\Rightarrow$  clefs, factorisation, val. par défaut, commentaires

```
&version=1.0
&default
# Note: currently, the 2 recognized types are "tracer" and "tag"
  parent=air hadv=10 vadv=10 phases=g type=tracer
&lmdz
  blue,red          parent=H[2]HO,H2O      type=tag          # tagging tracer of gen 1&2 tracers
  H2O               phases=g              hadv=14   vadv=14   # water vapour
  water,H2[18]O,H[2]HO,H[3]HO parent=H2O          # H2O childs
  yellow           parent=H[3]HO          type=tag          # tagging tracer for single isotope
  H2O              phases=ls              # H2O in liquid and solid phases
```

- Gestion de plusieurs listes de traceurs (fusion/cumul)  $\Rightarrow$  **type\_trac=t1,t2 (tracers\_t[12].def)**
- Stockage synthétique des grandeurs utiles  $\Rightarrow$  base de données + valeurs par défaut  
 $\Rightarrow$  accès simple: **getKey(tra/itr,keyn,val)**
- Fichier de paramètres non modifiables a priori: **isotopes\_params.def**

```
&H2O
  H2[16]O  tnat=1.0000   alpha=1.000
  H2[17]O  tnat=40.000e-6 alpha=1.003
  H2[18]O  tnat=2005.2e-6 alpha=1.006
  H[2]HO   tnat=155.78e-6 alpha=1.010
  H[3]HO   tnat=0.0000   alpha=1.000
```

# Nouveau fichier **tracer.def** - résultat

```
infotrac_init: Information stored in infotrac :
infotrac_init:  iq | tname | ttext | iadv | niadv | ipar
infotrac_init: -----
infotrac_init:  1 | H2O-g | H2O-gVLH | 14 | 1 | 0
infotrac_init:  2 | H2O-l | H2O-lVL1 | 10 | 2 | 0
infotrac_init:  3 | H2O-s | H2O-sVL1 | 10 | 3 | 0
infotrac_init:  4 | H2O-g_blue | H2O-g_blueVL1 | 10 | 4 | 1
infotrac_init:  5 | H2O-l_blue | H2O-l_blueVL1 | 10 | 5 | 2
infotrac_init:  6 | H2O-s_blue | H2O-s_blueVL1 | 10 | 6 | 3
infotrac_init:  7 | H2O-g_red | H2O-g_redVL1 | 10 | 7 | 1
infotrac_init:  8 | H2O-l_red | H2O-l_redVL1 | 10 | 8 | 2
infotrac_init:  9 | H2O-s_red | H2O-s_redVL1 | 10 | 9 | 3
infotrac_init: 10 | water-g | water-gVL1 | 10 | 10 | 1
infotrac_init: 11 | water-l | water-lVL1 | 10 | 11 | 2
infotrac_init: 12 | water-s | water-sVL1 | 10 | 12 | 3
infotrac_init: 13 | H2[18]O-g | H2[18]O-gVL1 | 10 | 13 | 1
infotrac_init: 14 | H2[18]O-l | H2[18]O-lVL1 | 10 | 14 | 2
infotrac_init: 15 | H2[18]O-s | H2[18]O-sVL1 | 10 | 15 | 3
infotrac_init: 16 | H[2]HO-g | H[2]HO-gVL1 | 10 | 16 | 1
infotrac_init: 17 | H[2]HO-l | H[2]HO-lVL1 | 10 | 17 | 2
infotrac_init: 18 | H[2]HO-s | H[2]HO-sVL1 | 10 | 18 | 3
infotrac_init: 19 | H[3]HO-g | H[3]HO-gVL1 | 10 | 19 | 1
infotrac_init: 20 | H[3]HO-l | H[3]HO-lVL1 | 10 | 20 | 2
infotrac_init: 21 | H[3]HO-s | H[3]HO-sVL1 | 10 | 21 | 3
infotrac_init: 22 | H[2]HO-g_blue | H[2]HO-g_blueVL1 | 10 | 22 | 16
infotrac_init: 23 | H[2]HO-l_blue | H[2]HO-l_blueVL1 | 10 | 23 | 17
infotrac_init: 24 | H[2]HO-s_blue | H[2]HO-s_blueVL1 | 10 | 24 | 18
infotrac_init: 25 | H[2]HO-g_red | H[2]HO-g_redVL1 | 10 | 25 | 16
infotrac_init: 26 | H[2]HO-l_red | H[2]HO-l_redVL1 | 10 | 26 | 17
infotrac_init: 27 | H[2]HO-s_red | H[2]HO-s_redVL1 | 10 | 27 | 18
infotrac_init: 28 | H[3]HO-g_yellow | H[3]HO-g_yellowVL1 | 10 | 28 | 19
infotrac_init: 29 | H[3]HO-l_yellow | H[3]HO-l_yellowVL1 | 10 | 29 | 20
infotrac_init: 30 | H[3]HO-s_yellow | H[3]HO-s_yellowVL1 | 10 | 30 | 21
infotrac_init: fin
```

# Isotopes – quelques outils

- Routines de vérification plus générales (xD) et génériques ; quelques dizaines de routines remplacées par:
  - NaNs: LOGICAL FUNCTION `checkNaN(x, err_msg[, name])`
  - Positivité: LOGICAL FUNCTION `checkPos(x, err_msg[, threshold])`
  - Égalité: LOGICAL FUNCTION `checkEqual(a, b, err_msg[, abs_err][, rel_err][, name])`
  - Valeurs aberrantes: LOGICAL FUNCTION `checkAbsurdDelta(iso_idx, q, err_msg[, qmin][, deltaMax])`
- Encapsulations supplémentaires possibles.
- Routines d'affichage: SUBROUTINE `dispOutliers(mask, a[, b], p, subname, err_msg[, nmax][, name])`
- $q(klon, klev), xt(niso, klon, klev) \Rightarrow q(klon, klev, niso+1)$   
 $\Rightarrow q\%q(klon, klev), q\%xt(klon, klev, niso)$
- Appels isotopiques "légers":

```
IF (turb_fcg_gcssold) THEN
  DO k = 1, klev
    zzdt(1:klon, k ) = hthturb_gcssold(k)*dtime_frcg
    zzdq(1:klon, k, 1:nisoH2O+1) = hqturb_gcssold(k)*dtime_frcg
  END DO
  ...
END IF
```