

## SUPPLEMENTARY MATERIAL

### MERGING BIO-OPTICAL DATA FROM BIOGEOCHEMICAL-ARGO FLOATS AND MODELS IN MARINE BIOGEOCHEMISTRY

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	<b>LOV NAME</b>	<b>WMO CODE</b>
<b>1</b>	lovbio001i	6901032
<b>2</b>	lovbio015c	6901513
<b>3</b>	lovbio016c	6901510
<b>4</b>	lovbio016d	6902700
<b>5</b>	lovbio017b	6901512
<b>6</b>	lovbio018c	6901528
<b>7</b>	lovbio035b	6901511
<b>8</b>	lovbio039b	6901483
<b>9</b>	lovbio042c	6901490
<b>10</b>	lovbio053b	6901529
<b>11</b>	lovbio058c	6901491
<b>12</b>	lovbio063c	6901653
<b>13</b>	lovbio064b	6901496
<b>14</b>	lovbio064c	6901776
<b>15</b>	lovbio066c	6901605
<b>16</b>	lovbio066d	6901655
<b>17</b>	lovbio067c	6901649
<b>18</b>	lovbio068d	6901648
<b>19</b>	lovbio072c	6901600
<b>20</b>	lovbio083d	6901764
<b>21</b>	lovbio085d	6901766
<b>22</b>	lovbio088d	6901768
<b>23</b>	lovbio089d	6901769
<b>24</b>	lovbio090d	6901770
<b>25</b>	lovbio091d	6901771
<b>26</b>	lovbio093d	6901773
<b>27</b>	ogsbio001b	6901861
<b>28</b>	ogsbio002b	6901864
<b>29</b>	ogsbio003b	6901862
<b>30</b>	ogsbio004b	6901863
<b>31</b>	ogsbio006b	6901865

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#=====
# NAMELISTS
#=====
#=====
# BFM - Biogeochemical Flux Model
#=====
#
# COPYING
#
# Copyright (C) 2015 BFM System Team (bfm_st@lists.cmcc.it)
#
# This program is free software; you can redistribute it and/or modify
# it under the terms of the GNU General Public License as published by
# the Free Software Foundation;
# This program is distributed in the hope that it will be useful,
# but WITHOUT ANY WARRANTY; without even the implied warranty of
# MERCHANTABILITY or FITNESS FOR A PARTICULAR PURPOSE. See the
# GNU General Public License for more details.
#
#-----
!-----!
!NAMELIST bfm_nml
!-----!
! Main initialisation and output specifications
! NAME          KIND   DESCRIPTION
! bio_calc      logical Switch on/off BFM (for coupled configurations)
! bfm_init      integer Initialization state
!                 0. from constant values in bfm_init_nml below
!                 1. from restart
! bfm_RSTCTL   logical Save initial state of bfm in the output file.
! bio_setup     integer BFM configuration:
!                 1. pelagic
!                 2. benthic
!                 3. pelagic and benthic
!                 4. sea ice
!                 5. pelagic and sea ice
! out_fname     string  Name of NetCDF output file
! out_dir       string  Path to the output file
! out_title     string  Name of the experiment in NetCDF file
! out_delta     integer Output is saved every out_delta timesteps
!                 Use -1 to store with real monthly frequency
! parallel_log  logical Set true in parallel jobs for a single log file
!-----!
&bfm_nml
    bio_calc      = .TRUE.
    bfm_init      = 0
    bfm_RSTCTL   = .FALSE.
    bio_setup     = 1
    out_fname     = 'BFM_standalone_pelagic'
    out_dir       = '.'
    out_title     = 'BFM_STANDALONE_PELAGIC'
    in_RST fname = 'in_bfm_restart'
    out_delta     = 300 !every 30 days
    parallel_log  = .FALSE.
        filename_nml_conf = 'BFM_General.nml',
/

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!-----!
! NAMELIST Param_parameters
!-----!
! Global Switches : turn on/off or choose model components
! NAME                      KIND      DESCRIPTION
! CalcPelagicFlag           logical   Pelagic System
! CalcBenthicFlag            numeric   Benthic system
!                                         0 = No Benthic System
!                                         The following are Not Yet Activated
!                                         1 = Simple Benthic Return
!                                         2 = Benthic organisms and intermediate
!                                              complexity nutrient regeneration
!                                         3 = Benthic organisms and full nutrient
!                                              regeneration (early diagenesis)
! CalcTransportFlag          logical   Compute Transport Term (when coupled
!                                         with a OGCM)
! CalcConservationFlag       logical   Mass Conservation Check
! CalcPhytoPlankton          logical   Pelagic Phytoplankton (vector)
! CalcPelBacteria             logical   Pelagic Bacteria (vector)
! CalcMesoZooPlankton        logical   Mesozooplankton (vector)
! CalcMicroZooPlankton       logical   Microzooplankton (vector)
! CalcPelChemistry            logical   Pelagic Hydrochemical Processes
! AssignPelBenFluxesInBFMFlag logical   Benthic-pelagic fluxes are added to the
!                                         time integration
! AssignAirPelFluxesInBFMFlag logical   Air-sea fluxes are added to the
!                                         time integration
! ChlDynamicsFlag            numeric   Choose the dynamics of Chl-a
!                                         1 = diagnostic, optimal light property
!                                              in phytoplankton
!                                              (Ebenhoech et al 1995, ERSEM-II)
!                                         2 = state variable, constituent of
!                                              phytoplankton
! check_fixed_quota          numeric   Check whether zooplankton have fixed
quota
!
! Global Parameters : used throughout the model and not related
!                     to a specific component
! NAME      UNIT      DESCRIPTION
! p_small    [-]       Smallest numeric value (the model "zero")
! slp0       [mbar]    Reference sea level pressure
! p_pe_R1c   [-]       Fractional content of C in cytoplasm
! p_pe_R1n   [-]       Fractional content of N in cytoplasm
! p_pe_R1p   [-]       Fractional content of P in cytoplasm
! p_qro      [mmolHS-/mmolO2] Stoichiometric coefficient for
!                               anaerobic reactions
! p_qon_dentri [mmolO2/mmolN] Stoichiometric coefficient for
!                               denitrification
! p_qon_nitri [mmolO2/mmolN] Stoichiometric coefficient for
!                               nitrification
!-----!
&Param_parameters
! Switches :
    CalcPelagicFlag = .TRUE.
    CalcBenthicFlag = 0
    CalcConservationFlag = .FALSE.
    CalcTransportFlag = .FALSE.
    CalcPhytoPlankton(1) = .TRUE.
    CalcPhytoPlankton(2) = .TRUE.
    CalcPhytoPlankton(3) = .TRUE.
    CalcPhytoPlankton(4) = .TRUE.

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        CalcPelBacteria(1) = .TRUE.
CalcMicroZooPlankton(1) = .TRUE.
CalcMicroZooPlankton(2) = .TRUE.
CalcMesoZooPlankton(1) = .TRUE.
CalcMesoZooPlankton(2) = .TRUE.
        CalcPelChemistry = .TRUE.
AssignPelBenFluxesInBFMFlag = .FALSE.
AssignAirPelFluxesInBFMFlag = .TRUE.
        ChlDynamicsFlag = 2
        check_fixed_quota = 0
! Parameters :
        p_small = 1.0e-20
        slp0 = 1013.25E0
        p_pe_Rlc = 0.60
        p_pe_Rln = 0.72
        p_pe_Rlp = 0.832
        p_qro = 0.5
        p_qon_dentri = 1.25
        p_qon_nitri = 2.0
filename_nml_conf = 'BFM_General.nml',
/
!-----
! NAMELIST bfm_init_nml
!-----
! Pelagic initialisation of standard variables
!<variablename>0 = <realvalue>
!-----
&bfm_init_nml
    O2o0 = 300.0,
    N1p0 = 1.0,
    N3n0 = 5.0,
    N4n0 = 1.0,
    N5s0 = 8.0,
    N6r0 = 1.0,
    O3c0 = 27060.00,
    O3h0 = 2660.0,
    O4n0 = 200.0,
    P1c0 = 1.0,
    P2c0 = 1.0,
    P3c0 = 1.0,
    P4c0 = 1.0,
    Z3c0 = 1.0,
    Z4c0 = 1.0,
    Z5c0 = 1.0,
    Z6c0 = 1.0,
    B1c0 = 1.0,
    R1c0 = 1.0,
    R2c0 = 0.1,
    R3c0 = 1.0,
    R6c0 = 1.0,
filename_nml_conf = 'BFM_General.nml',
/
!-----
! NAMELIST bfm_save_nml
!-----
! Stored variables
!ave_save: average values over the output interval
!var_save: instantaneous value at the output interval
!-----
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&bfm_save_nml
  var_save = ''
  ave_save = 'ETW', 'O2o', 'DIC',
    'EIR',
    'xEPS',
    'Chla',
    'N1p', 'N3n', 'N4n', 'N5s',
    'B1c',
    'P1c', 'P2c', 'P3c', 'P4c',
    'P11', 'P21', 'P31', 'P41',
    'P1n', 'P2n', 'P3n', 'P4n',
    'P1p', 'P2p', 'P3p', 'P4p',
    'Z3c', 'Z4c', 'Z5c', 'Z6c',
    'R1c', 'R2c', 'R6c',
    'R1n', 'R6n',
    'R1p', 'R6p',
    'P1s', 'R6s',
    'eiPPY(iiP1)', 'eiPPY(iiP2)', 'eiPPY(iiP3)', 'eiPPY(iiP4)',
    'sunPPY(iiP1)', 'sunPPY(iiP2)', 'sunPPY(iiP3)', 'sunPPY(iiP4)',
    'ruPPYc', 'resPBac', 'resZOOc',
    'ruPPYn', 'ruPPYp', 'ruPPYs', 'exPPYc',
    'ruZOOc', 'remZOOon', 'remZOOp', 'remPBAn', 'remPBAp'

  filename_nml_conf = 'BFM_General.nml',
/

!-----!
!NAMELIST Settling_parameters
!-----!
! BURIAL VELOCITIES into the sediment
! NAME          [UNIT]/KIND      DESCRIPTION
! p_burvel_R6   [m/d]           Bottom Burial Velocity for detritus
! p_burvel_R2   [m/d]           Bottom Burial Velocity for dissolved
! p_burvel_PI   [m/d]           Bottom Burial Velocity for plankton
!-----!
&Settling_parameters
  p_burvel_R6 = 1.5,
  p_burvel_R2 = 0.0,
  p_burvel_PI = 0.0
  filename_nml_conf = 'Benthic_Environment.nml',
/


!-----!
!NAMELIST CO2_parameters
!-----!
! CARBONATE SYSTEM SETTING
! NAME          [UNIT]/KIND      DESCRIPTION
! AtmCO20       [ppmv]          Initial atmospheric concentration of CO2
! calcAtmpCO2   logical         Compute the partial pressure of Atmospheric
CO2
! pCO2Method    integer         pCO2 computation method: 1=MixRatio*slp0,
2=Magnus formula
! phstart       [pH]            Initial pH value
! K1K2          integer         Switch for the acidity constants
parameterization
!                               1 : Roy et al. (1993); DOE (1994); pH on total
scale
!                               2 : Default. OCMIP STANDARD; pH on Sea Water
Scale

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!
!               Mehrbach et al (1973) refit by Dickson &
Millero (1987)
!
!               3 : Mehrbach et al (1973) refit by Lueker et
al. (2000)
!
!               pH on total scale
!
!               4 : Hansson (1973b) data as refitted by
Dickson and
!
!               Millero (1987); pH on Sea Water Scale
MethodCalcCO2 numeric
computation
!
!
!
!               Switch for the choice of [H+] numerical
CalcBioAlkFlag logical
total alkalinity
!
!               ----- Parameters for MethodCalcCO2=2 -----
!
! M2XACC      real          Accuracy of the iterative scheme for OCMIP
(default 1.E-10)
!
! M2PHDELT    [pH]          Delta of pH for the root search (realized
pH+/-DELT)
!
! M2MAXIT     integer       in the OCMIP scheme (default 0.5)
Maximum number of iterations for OCMIP
(default 100 )
!
!               ----- Parameters for calcium and calcite -----
!
! Caconc0     [mol/m3]      Calcium ion concentration
!
! behaviour"                                         ["Seawater : Its composition, properties and
1995]
!
! Seawater concentration = 412 mg / l
!                                     -> atomic weight =
40.078 g / mol
!
! therefore, concentration = 10.279 mmol / l =
10.279 mol / m3
!
! Canorm      logical       Normalize Calcium ion concentration by sea
water salinity
!
!               ----- EXTERNAL DATA INPUT STRUCTURES -----
!
! AtmCO2_N    structure     Read external data for atmospheric CO2 values
!
! AtmSLP_N    structure     Read external data for atmospheric sea level
pressure
!
! AtmTDP_N    structure     Read external data for atmospheric dew-point
temperature
!
! Example of general input structure for the data structure:
!
!           ! Read ! File                                ! NetCDF ! Var      !
!           ! Input ! name                               ! Logical ! name    !
!AtmCO2_N = 0 , 'CMIP5_Historical_GHG_1765_2005.dat' , .FALSE. , 'CO2' ,
!
!           ! RefTime        ! Input      ! Time      !
!           ! yyyyymmdd   ! Frequency  ! interp    !
!           '1764-07-01 00:00' , 'yearly' , .TRUE.
!
!
! Convention for Input reading : 0 = use constant value (default if structure is
not initialized)
!
ratios)                                         2 = read timeseries file ( e.g. CO2 mixing
!
SLP from OGCM)                                 4 = field from a coupled model (e.g. atmospheric
!
! NOTE: The file "CMIP5_Historical_GHG_1765_2005.dat" is located in
"$BFMDIR/tools" folder
!
-----!
&CO2_parameters

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AtmCO20      = 370.0E0
calcAtmpCO2  = .FALSE.
pCO2Method   = 1
AtmCO2_N     = 0 , 'CMIP5_Historical_GHG_1765_2005.dat' , .FALSE. , 'CO2'
, '1764-07-01 00:00' , 'yearly' , .TRUE.
AtmSLP_N     = 0 , 'AtmSLP.nc' , .TRUE. , 'AtmSLP' , '1764-07-01 00:00' ,
'dummy' , .TRUE.
AtmTDP_N     = 0 , 'AtmTDP.nc' , .TRUE. , 'AtmTDP' , '1764-07-01 00:00' ,
'dummy' , .TRUE.
phstart      = 8.10E0
K1K2         = 2
MethodCalcco2 = 2
CalcBioAlkFlag = .FALSE.
M2XACC       = 1.0E-10
M2PHDELT    = 0.3
M2MAXIT     = 100
Caconc0     = 10.279E0
Canorm      = .TRUE.
filename_nml_conf = 'Carbonate_Dynamics.nml',
/
!-----!
!NAMELIST PelBac_parameters
!-----!
! PELAGIC BACTERIA
!
! NAME      [UNIT]/KIND          DESCRIPTION
! p_version integer            Switch for bacteria parameterization
!                               1 : Baretta-Bekker et al. 1995;
!                               Vichi et al., 2007
!                               2 : Vichi et al., 2004
!                               3 : Polimene et al., 2006
! p_q10     [-]                Q10-value (temperature dependency)
! p_chdo    [mmol/m3]          Half-saturation constant for O2 limitation
! p_sd      [1/d]              Specific mortality rate
! p_sd2     [1/d]              Density dependent specific mortality rate
! p_suhR1   [1/d]              Specific potential uptake for nutrient-rich DOM
! p_sulR1   [1/d]              Specific potential uptake for nutrient-poor DOM
! p_sur2    [1/d]              Specific potential uptake for semi-labile DOC
! p_sur3    [1/d]              Specific potential uptake for semi-refractory DOC
! p_sur6    [1/d]              Specific potential uptake for POM (1/d)
! p_sum     [1/d]              Potential specific growth rate
! p_pu_ra   [-]               Activity respiration fraction
! p_pu_ra_o [-]              Additional respiration fraction at low O2 conc
! p_srs     [1/d]              Specific rest respiration
! p_qncPBA  [mmolN/mgC]        Optimal N/C ratio
! p_qpcPBA  [mmolP/mgC]        Optimal P/C ratio
! p_qlnc    [mmolN/mgC]        Minimal N/C ratio
! p qlpc   [mmolP/mgC]        Minimal P/C ratio
! p_qun    [mmolN/mgC/day]    Membrane affinity for N
! p_qup    [mmolP/mgC/day]    Membrane affinity for N
! p_chn    [mmolN/m3]          Half saturation ammonium conc. for uptake
! p_chp    [mmolP/m3]          Half saturation phosphate conc. for uptake
! p_ruen   [1/d]              Relaxation timescale for N uptake/remin.
! p_ruep   [1/d]              Relaxation timescale for P uptake/remin.
! p_rec    [1/d]              Relaxation timescale for semi-labile excretion
! p_pu_ea_R3 [-]             Excretion of semi-refractory DOC
!-----!
&PelBacteria_parameters
!
!           B1
p_version = 2

```



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!
!           2. Combined
!           ---- N limitation control ----
!   p_qun      [m3/mgC/d]      Membrane affinity for N
!   p_ln4      [mmolN/m3]      Half saturation constant for NH4 uptake preference
over NO3
!   p_qnlc     [mmolN/mgC]    Minimum quotum Si:C
!   p_qncPPY   [mmolN/mgC]    Reference quotum Si:C
!   p_xqn      [-]           Multiplication factor for luxury storage
!
!           ---- P limitation control ----
!   p_qup      [m3/mgC/d]      Membrane affinity for P
!   p_qplc     [mmolP/mgC]    Minimum quotum Si:C
!   p_qpcPPY   [mmolP/mgC]    Reference quotum Si:C
!   p_xqp      [-]           Multiplication factor for luxury storage
!
!           ---- Si limitation control ----
!   p_switchSi [1-2]         Switch for Silica limitation
!                           1. Si limitation is controlled by external Si
conc.
!
!                           2. Si limitation is controlled by internal quota
!   p_chPs      [mmolSi/m3]   Half saturation conc. for dissolved Si limitation
!   p_Contois   [>=0]        If >0, use Contois formulation
!   p_quS       [m3/mgC/d]    Membrane affinity for Si
!   p_qsIc      [mmolSi/mgC]  Minimum quotum Si:C
!   p_qsCPPY   [mmolSi/mgC]  Reference quotum Si:C
!                           Brzezinski (1985) 0.13 mol/mol
!
!           ---- nutrient stressed sinking ----
!   p_esNI      [-]          Nutrient stress threshold for sinking
!   p_res       [m/d]        Maximum Sinking vel city (m/d)
!
!           ----- Chlorophyll parameters -----
!   p_switchChl [1-4]        Switch for Chla-a synthesis
!   p_sdchl     [1/d]        Specific turnover rate for Chla
!   p_alpha_chl [mgC s m2/   Initial slope of the P-E curve
!                 mgChl/uE]
!   p_qlcPPY    [mgChla/mgC] Reference quotum Chla:C
!   p_epsChla   [m2/mgChla]  Chla-specific extinction coefficient
!   p_tochl_relt [1/d]       Relaxation rate towards maximum Chla:C
!   p_EpEk_or   [-]          Optimal value of E_PAR/E_K
!
!           ----- Iron parameters -----
!   p_quf       [m3/mgC/d]   Membrane affinity for Fe
!   p_qflc      [umolFe/mgC] Minimum quotum Fe:C derived from 3 umol Fe/mol C
!                           Sunda & Huntsman (1997), Nature, 390, p 389-392
!   p_qfcPPY   [umolFe/mgC] Reference quotum Fe:C
!   p_xqf       [-]          Multiplication factor for luxury storage
!
!           ----- Light parameters ERSEM-II -----
!   p_iswLtyp   [0-6]        Shape of the productivity function
!   p_chELIPPY  [W/m2]       Maximum Iopt
!   p_cLELiPPY  [W/m2]       Minimum Iopt
!   p_ruELIPPY  [1/d]        Maximum daily shift in Iopt (1/d)
!   p_addepth   [m]          Adaptation depth. Meaningless with high-res models
!
!           ----- Sinking parameters -----
!   p_rPIm     [m/d]        Phytoplankton background sinking rate
!-----!
&Phyto_parameters
!
!           P1      P2      P3      P4
p_q10 = 2.0,    2.0,    2.0,    2.0
p_temp = 0.0,    0.0,    0.75,    0.0
p_sum = 2.5,    3.0,    3.5,    1.5
p_srs = 0.1,    0.05,   0.1,    0.1
p_sdmo = 0.0,    0.0,    0.0,    0.0
p_thdo = 0.0,    0.0,    0.0,    0.0
p_seo = 0.0,    0.0,    0.0,    0.0
p_sheo = 0.0,    0.0,    0.0,    100.0

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    p_pu_ea = 0.05,    0.1,      0.1,      0.15
    p_pu_ra = 0.1,     0.1,      0.2,      0.1
p_switchDOC = 2,      2,       2,       2
p_netgrowth = .TRUE.,.TRUE.,.TRUE.,.TRUE.
    p_limnut = 1,      1,       1,       1
    p_qun = 0.025,   0.025,   0.25,   0.025
    p_lN4 = 1.0,     0.5,     0.1,     1.0
    p_qnlc = 0.00687, 0.00687, 0.00687, 0.00687
p_qncPPY = 1.26e-2,1.26e-2,1.26e-2,1.26e-2
    p_xqn = 2.0,     2.0,     2.0,     2.0
    p_qup = 0.0025,  0.0025,  0.0025,  0.0025
    p_qplc = 0.0004288,0.0004288,0.0004288,0.0004288
p_qpcPPY = 7.86e-4,7.86e-4,7.86e-4,7.86e-4
    p_xqp = 2.0,     2.0,     2.0,     2.0
p_switchSi = 2,      0,       0,       0
    p_chPs = 0.3,    0.0,     0.0,     0.0
p_Contois = 0.0,    0.0,     0.0,     0.0
    p_qua = 0.0025,  0.0,     0.0,     0.0
    p_qslc = 0.007,   0.0,     0.0,     0.0
p_qscPPY = 0.01,    0.0,     0.0,     0.0
    p_esNI = 0.7,    0.75,   0.75,   0.75
    p_res = 5.0,    0.0,     0.0,     2.5
p_switchChl = 2,      2,       2,       2
    p_sdchl = 0.2,   0.2,     0.2,     0.2
p_alpha_chl = 2.50e-5, 1.70e-5, 2.75e-5, 0.68e-5
    p qlcPPY = 0.02,  0.02,   0.02,   0.02
    p_epsChla = 0.03, 0.03,   0.03,   0.03
    p_EpEk_or = 0.0,  0.0,     0.0,     0.0
p_tochl_relt = 0.0,  0.0,     0.0,     0.0
    p_iswLtyp = 5,    5,       5,       5
    p_addepth = 50.0, 50.0,   50.0,   50.0
p_cheliPPY = 100.0, 100.0, 100.0, 100.0
p_cleliPPY = 8.0,   10.0,   6.0,    12.0
p_rueLiPPY = 0.2,   0.25,  0.3,    0.15
    p_rPIm = 0.0,   0.0,     0.0,     0.0
    filename_nml_conf = 'Pelagic_Ecology.nml',
/
&Phyto_parameters_iron
    filename_nml_conf = 'Pelagic_Ecology.nml',
/
!-----!
!NAMELIST PAR_parameters
!-----!
! LightPeriodFlag      numeric Choose the light averaging period
!                           1 = Instantaneous irradiance
!                           2 = Daily average
!                           3 = Daylight average with explicit
!                                photoperiod
! LightLocationFlag    numeric Choose the parameterization of light
!                           location in the discrete grid
!                           1 = Light at the top of the cell
!                           2 = Light in the middle of the cell
!                           3 = Average Light in the cell
! ChlAttenFlag         numeric Choose the PAR attenuation due to Chl
!                           1 = broadband linear attenuation
!                           2 = 3-band tabulated attenuation coefficients
!                                (Morel, 1988; Lengaigne et al, 2007)
! p_PAR                [-]      Fraction of Photosynthetically Available Radiation
! p_eps0               [1/m]    Background extinction coefficient

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! p_epsIR      [1/m]          Infrared extinction coefficient
!                                         (to use with ChlAttenFlag=2)
! p_epsESS     [m2/g]          Specific attenuation coefficient of
!                                         suspended sediments
! p_epsR6      [m2/mgC]        Specific attenuation coefficient of particulate
!                                         detritus
!-----
!&PAR_parameters
LightPeriodFlag    = 1
LightLocationFlag  = 2
ChlAttenFlag       = 1
p_PAR              = 0.40
p_eps0             = 0.0435
p_epsIR            = 2.857
p_epsESS            = 0.04d-3
p_epsR6            = 0.1d-3
filename_nml_conf = 'Pelagic_Environment.nml',
/
!-----
!NAMELIST MicroZoo_parameters
!-----
! MICRO-ZOOPLANKTON
!
! NAME          [UNIT]/KIND           DESCRIPTION
! p_q10          [-]                Q10 value for physiological rates
! p_srs          [1/d]              Respiration rate at 10 degrees Celsius
! p_sum          [1/d]              Potential growth rate
! p_sdo          [1/d]              Mortality rate due to oxygen limitation
! p_sd           [1/d]              Temperature independent mortality rate
! p_pu           [-]                Assimilation efficiency
! p_pu_ea         [-]                Fraction of activity excretion
! p_chro          [mmolO2/m3]       Half-saturation oxygen concentration
! p_chuc          [mgC/m3]          Half-saturation Food concentration for Type II
! p_minfood       [mgC/m3]          Half-saturation food concentration for
!                                         preference factor
! p_qncMIZ        [mmolN/mgC]      Maximum quotum P:C
! p_qpcMIZ        [mmolN/mgC]      Maximum quotum N:C
! p_paPBA(z,b)   [-]              Availability of pelagic Bacteria group b
!                                         to Zooplankton group z
! p_paPPY(z,p)   [-]              Availability of PhytoPlankton group p
!                                         to Zooplankton group z
! p_paMIZ(z,m)   [-]              Availability of MicroZooplankton group m
!                                         to Zooplankton group z
!-----
!&MicroZoo_parameters
!
!          Z5      Z6
p_q10 = 2.0,    2.0
p_srs = 0.02,   0.02
p_sum = 2.0,    5.0
p_sdo = 0.05,   0.05
p_sd = 0.0,    0.0
p_pu = 0.5,    0.3
p_pu_ea = 0.5, 0.5
p_chro = 8,     8
p_chuc = 30.0, 100.0
p_minfood = 50.0, 50.0
p_qpcMIZ = 1.85d-3, 1.85d-3
p_qncMIZ = 1.67d-2, 1.67d-2
! Food matrix parameters: take care of the notation
!
!          Z5      Z6

```

```

    p_paPBA = 0.1,      1.0
!           P1        P2        P3        P4
!
!           Z5
p_paPPY(1,:) = 0.7,      1.0,      0.1,      0.1
!
!           Z6
p_paPPY(2,:) = 0.0,      0.2,      1.0,      0.0
!
!           Z5
p_paMIZ(1,:) = 1.0,      1.0
!
!           Z6
p_paMIZ(2,:) = 0.0,      0.2
    filename_nml_conf = 'Pelagic_Ecology.nml',
/
!

!-----!
!NAMELIST MesoZoo_parameters
!-----!
! MESO-ZOOPLANKTON
! NAME          [UNIT]/KIND      DESCRIPTION
! p_q10         [-]            Q10 value for physiological rates
! p_srs         [1/d]          Respiration rate at 10 degrees C
! p_sum         [1/d]          Maximal productivity at 10 degrees C
! p_sd          [1/d]          Background natural mortality
! p_vum         [m3/mgC/d]     Specific search volume
! p_pui         [-]            Assimilation efficiency
! p_peI          [-]            Fraction of Faeces production
! p_sdo         [m3/mgC/d]     Specific density-dependent mortality
! p_sds         [-]            Exponent of density-dependent mortality
! p_qpcMEZ      [mmolP/mgC]    Maximum quotum P:C
! p_qncMEZ      [mmolN/mgC]    Maximum quotum N:C
! p_clo2o       [mmolO2/m3]    Half-saturation oxygen concentration
! p_paPPY(z,p) [-]          Availability of PhytoPlankton group p
!                           to Zooplankton group z
! p_paMIZ(z,m)  [-]          Availability of MicroZooplankton group m
!                           to Zooplankton group z
! p_paMEZ(z,m)  [-]          Availability of MesoZooplankton group m
!                           to Zooplankton group z
!
!-----!
&MesoZoo_parameters
!
!           Z3        Z4
p_q10 = 2.0,      2.0
p_srs = 0.01,      0.02
p_sum = 2.0,      2.0
p_vum = 0.008,      0.02
p_pui = 0.6,      0.6
p_peI = 0.3,      0.35
p_sdo = 0.01,      0.01
p_sd = 0.02,      0.02
p_sds = 2.0,      2.0
p_qpcMEZ = 1.67d-3, 1.67d-3
p_qncMEZ = 0.015, 0.015
p_clo2o = 30.0, 30.0
!
! Food matrix parameters: take care of the notation
!           P1        P2        P3        P4
!
!           Z3
p_paPPY(1,:) = 0.0,      0.0,      0.0,      1.0
!
!           Z4
p_paPPY(2,:) = 1.0,      0.75,     0.0,      1.0
!
!           Z3
p_paMIZ(1,:) = 0.0,      0.0

```

```

!      Z4
p_paMIZ(2,:) = 1.0,    0.0,
!           Z3      Z4
!
!      Z3
p_paMEZ(1,:) = 1.0,    1.0
!           Z4
p_paMEZ(2,:) = 0.0,    1.0
      filename_nml_conf = 'Pelagic_Ecology.nml',
/
!-----!
!NAMELIST PelChem_parameters, PelChem_parameters_iron
!-----!
! Pelagic Chemistry parameters
! NAME      [UNIT]/KIND      DESCRIPTION
! p_q10N4N3  [-]            Q10 factor for nitrification/denit
! p_sN4N3    [1/d]          Specific nitrification rate at 10 degC
! p_cLO2o    [mmolO2/m3]    Half-saturation O2 concentration for
!                           nitrification and reoxidation
!
! p_rOS      [1/d]          Specific reoxidation rate of reduction
!                           equivalents
!
! p_sN3O4n   [1/d]          Specific denitrification rate
! p_cLN6r    [mmolHS/m3]    Half-saturation concentration of
!                           reduction equivalents for denitrification
!
! p_rPAo     [mmolO2/m3/d]  Reference anoxic mineralization rate
! p_q10R6N5   [-]            Q10 factor for biogenic silica
! p_sR6N5    [1/d]          Specific remineralization rate of
!                           biogenic silica
!
!           ----- Iron parameters -----
! p_q10R6N7   [-]            Q10 temperature dependence
! p_sR6N7    [1/d]          Specific remineralization rate of particulate
! p_sR1N7    [1/d]          Specific remineralization rate of dissolved
! p_scavN7f   [1/d]          Specific scavenging rate
! p_N7fsol   [umolFe/m3]   Solubility concentration
!
!-----!
&PelChem_parameters
  p_q10N4N3 = 2.367
  p_sN4N3 = 0.01
  p_cLO2o = 10.0
  p_rOS = 0.05
  p_sN3O4n = 0.35
  p_cLN6r = 1.0
  p_rPAo = 1.0
  p_q10R6N5 = 1.49
  p_sR6N5 = 0.1
      filename_nml_conf = 'Pelagic_Environment.nml',
/
&PelChem_parameters_iron
      filename_nml_conf = 'Pelagic_Environment.nml',
/
!-----!
!NAMELIST PelGlobal_parameters
!-----!
! Sinking rates of Pelagic Variables
! : for mem_PelGlobal filled by InitPelGlobal
! NAME      UNIT      DESCRIPTION
! p_rR6m    [m/d]    detritus sinking rate
! KSINK_rPPY [m]      prescribe sinking rate for phytoplankton below this
!                           depth threshold to p_rR6m value. Use 0.0 to disable.
!
```

```

! AggregateSink logic use aggregation = true to enhance the sink rate
!                                         and bypass the prescribed sinking
! depth_factor [m]      depth factor for aggregation method
!-----
!&PelGlobal_parameters
  p_rR6m      = 5.0
  KSINK_rPPY   = 150.0
  AggregateSink = .FALSE.
  depth_factor = 2000.0
    filename_nml_conf = 'Pelagic_Environment.nml',
/
!-----
!NAMELIST standalone_nml
!-----
!NAME      KIND DESCRIPTION
!nboxes    integer   Number of water volumes (boxes)
!indepth   real     Depth of each box (m)
!latitude  real     Latitude of each box
!longitude real     Longitude of each box
!maxdelt   real     Maximum timestep duration (s)
!mindelt   real     Minimum timestep duration (s)
!method    integer  Integration method
!
!           1. Euler forward
!           2. Runge-Kutta 2nd order
!           3. Leap-frog
!-----
&standalone_nml
  nboxes      = 1,
  indepth     = 5.0,
  latitude    = 45.0,
  longitude   = 13.5,
  maxdelt     = 8640.0,
  mindelt     = 1.0,
  method       = 2,
    filename_nml_conf = 'Standalone.nml',
/
!-----
!NAMELIST time_nml
!-----
!Specify time related formats and variables here.
!timefmt    integer   implicitly uses timestep=maxdelt
!
!           1- MaxN only - fake start time used.
!           2- start and stop - MaxN calculated.
!           3- start and MaxN - stop calculated.
!           4- simdays - fake start time used
!                     and MaxN calculated.
!MaxN      integer   do loop from n=1,MaxN
!start      string    Initial time: YYYY/MM/DD HH:MM:SS
!stop       string    Final   time: YYYY/MM/DD HH:MM:SS
!-----
!&time_nml
  timefmt= 2,
  MaxN=    144,
  simdays= 5760,
  start=   "2000-01-01 00:00:00",
  stop=    "2010-01-01 00:00:00",
    filename_nml_conf = 'Standalone.nml',
/

```

```

!-----!
!NAMELIST forcings_nml
!-----!
!forcing_method integer      Choice of the external forcing functions
!          1 = analytical forcings
!          2 = from file
!          3 = interactive fluxes (not yet implemented)
!
!Method 1: Analytical forcing functions
!lw       real Sinusoidal light intensity (winter) W m^-2
!ls       real Sinusoidal light intensity (summer) W m^-2
!sw       real Sinusoidal salinity (winter)
!ss       real Sinusoidal salinity (summer)
!tw       real Sinusoidal temperature (winter) degC
!ts       real Sinusoidal temperature (summer) degC
!tde      real Sinusoidal temperature daily excursion degC
!ww       real Sinusoidal wind (winter) m/s
!ws       real Sinusoidal wind (summer) m/s
!CO2inc    real     Linear increase in CO2 air partial pressure [% per year]
!
!Method 2: data file
!forcing_file   char  Filename for external forcings
!
!use_external_data logical Read external data (user defined)
!data_file      char      Filename for external data
!
!-----!
&forcings_nml
  forcing_method      = 1,
  ltype               = 1,
  lw                  = 20.0,
  ls                  = 300.0,
  sw                  = 37.0,
  ss                  = 34.0,
  tw                  = 8.0,
  ts                  = 28.0,
  tde                 = 1.0,
  ww                  = 20.0,
  ws                  = 10.0,
  forcing_file        = '',
  use_external_data   = .false.,
  data_file           = '',
  filename_nml_conf  = 'Standalone.nml',
/
!=====
! MODEL  BFM - Biogeochemical Flux Model
!=====
```