# MARBL diagnostics for vertical and horizontal nutrient distribution for ocean history files

```
In [1]: import warnings
warnings.filterwarnings('ignore')
%matplotlib inline
import os
from glob import glob
from collections import OrderedDict
import xarray as xr
import numpy as np
import esmlab
import matplotlib.pyplot as plt
import cartopy
import cartopy.crs as ccrs
import seawater as sw
from seawater.library import T90conv
from scipy import stats
```

# Define year range and get the CESM data

```
In [2]: start_yr = 306
num_years = 61
endyr = start_yr + num_years
case = 'g.e22.GOMIPECOIAF_JRA-1p4-2018.TL319_g17.SMYLE.005'
user = 'klindsay'
```

```
In [3]: files = []
                for year in range(start yr,endyr):
                        yr4="{:04d}".format(year)
                        print('doing simulation year', year, '!')
                        for month in range(1, 13):
                                mo2="{:02d}".format(month)
                                files.extend(sorted(glob(f'/glade/scratch/{user}/archive/{case}))
                                #files.extend(sorted(glob(f'/glade/scratch/{user}/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case}/run/{case
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In [4]: %%time
                cesm mon ds=xr.open mfdataset(files,decode times=False,decode coords=Fal
```

```
CPU times: user 6min 28s, sys: 12.9 s, total: 6min 41s
```

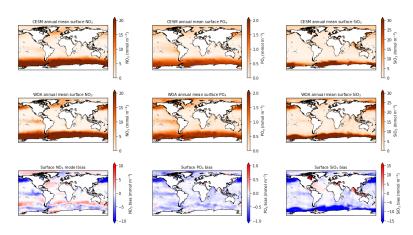
Wall time: 24min 12s

```
In [6]: coords = {'x':'TLONG','y':'TLAT'}
                          keepthese=['z_t','z_t_150m','time_bound','TAREA','PO4','Fe','NO3','SiO3 keep_vars = keepthese +list(coords.values())+['dz','KMT']
                          cesm_mon_ds = cesm_mon_ds.drop([v for v in cesm_mon_ds.variables if v notest
   In [7]: cesm_ann_ds=cesm_mon_ds.mean(dim='time')
   In [3]: file='tmp_mean.nc'
                           #cesm_ann_ds.to_netcdf(file)
                          cesm_ann_ds = xr.load_dataset(file,decode_times=False)
                          World Ocean Atlas 2013
   In [4]: file = '/glade/work/kristenk/WOA_data/regrid_POP/WOA2013_POPgrid.nc'
   In [5]: ds_woa = xr.load_dataset(file, decode_times=False, decode_coords=False)
   In [6]: ds_woa['z_t'] = cesm_ann_ds.z_t
   In [7]: NO3 diff = cesm ann ds.NO3 - ds woa.NO3
                           PO4_diff = cesm_ann_ds.PO4 - ds_woa.PO4
                          SiO3_diff = cesm_ann_ds.SiO3 - ds_woa.SiO3
   In [8]: lons=ds_woa.TLONG
                          lats=ds_woa.TLAT
                          area=ds_woa.TAREA
                          In [14]: rmse_global = xr.Dataset({v: cesm_ann_ds[v] for v in ['z_t']})
                          rmse_global['N03']=esmlab.statistics.weighted_rmsd(cesm_ann_ds.N03, ds_vrmse_global['P04']=esmlab.statistics.weighted_rmsd(cesm_ann_ds.P04, ds_vrmse_global['P04']=esmlab.stat
                           rmse_global['SiO3']=esmlab.statistics.weighted_rmsd(cesm_ann_ds.SiO3, ds
```

Surface nutrients

```
In [65]: fig = plt.figure(figsize=(18,10))
         plt.suptitle('Surface macronutrients', fontsize=14)
         ###############NO3
         #COLUMN 1 - NO3
         #--- CESM panel
         ax = fig.add_subplot(3,3,1, projection=ccrs.PlateCarree())
         ax.set extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set title('CESM annual mean surface NO$ 3$', fontsize=10)
         pc1=ax.pcolormesh(lons, lats,
                           cesm_ann_ds.NO3.isel(z_t=0), vmin=0, vmax=20, cmap='Ora
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc1, ax=ax,extend='max',label='NO$_3$ (mmol m$^{-3}]
         #--- OBS panel
         ax = fig.add_subplot(3,3,4, projection=ccrs.PlateCarree())
ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set_title('WOA annual mean surface NO$_3$', fontsize=10)
         pc2=ax.pcolormesh(lons, lats,
                           ds_woa.NO3.isel(z_t=0), vmin=0, vmax=20, cmap='Oranges
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc2, ax=ax,extend='max',label='NO$_3$ (mmol m$^{-3})
         #---- DIFF panel
         ax = fig.add_subplot(3,3,7, projection=ccrs.PlateCarree())
         ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set title('Surface NO$ 3$ model bias', fontsize=10)
         pc3=ax.pcolormesh(lons, lats,
                           NO3_diff.isel(z_t=0), vmin=-10, vmax=10, cmap='bwr',
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc3, ax=ax,extend='both',label='NO$_3$ bias (mmol r
         ############PO4
         #--- CESM panel
         ax = fig.add_subplot(3,3,2, projection=ccrs.PlateCarree())
         ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set title('CESM annual mean surface PO$ 4$', fontsize=10)
         pcl=ax.pcolormesh(lons, lats,
                           cesm_ann_ds.PO4.isel(z_t=0), vmin=0, vmax=2, cmap='Ora;
                           transform=ccrs.PlateCarree())
         \label{local_color_bar_pcl}  cbarl = fig.colorbar(pcl, ax=ax,extend='max',label='PO\$\_4\$ \ (mmol m\$^{-3}) 
         #---- OBS panel
         ax = fig.add_subplot(3,3,5, projection=ccrs.PlateCarree())
         ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set title('WOA annual mean surface PO$ 4$', fontsize=10)
         pc2=ax.pcolormesh(lons, lats,
                           ds_woa.PO4.isel(z_t=0), vmin=0, vmax=2, cmap='Oranges',
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc2, ax=ax,extend='max',label='PO$_4$ (mmol m$^{-3})
         #--- DIFF panel
         ax = fig.add_subplot(3,3,8, projection=ccrs.PlateCarree())
         ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set title('Surface PO$ 4$ bias', fontsize=10)
         pc3=ax.pcolormesh(lons, lats,
                           PO4 diff.isel(z t=0), vmin=-1, vmax=1, cmap='bwr',
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc3, ax=ax,extend='both',label='PO$ 4$ bias (mmol r
         ################si03
         #---- CESM panel
         ax = fig.add_subplot(3,3,3, projection=ccrs.PlateCarree())
         ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set_title('CESM annual mean surface SiO$_3$', fontsize=10)
         pc1=ax.pcolormesh(lons, lats,
                           cesm_ann_ds.SiO3.isel(z_t=0),
                            vmin=0, vmax=30,
                            cmap='Oranges',
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc1, ax=ax,extend='max',label='SiO$_3$ (mmol m$^{-1}
         #---- OBS panel
         ax = fig.add_subplot(3,3,6, projection=ccrs.PlateCarree())
         ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set_title('WOA annual mean surface SiO$_3$', fontsize=10)
         pc2=ax.pcolormesh(lons, lats,
                           ds_{woa.Si03.isel(z_t=0)},
                            vmin=0, vmax=30,
                            cmap='Oranges'
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc2, ax=ax,extend='max',label='SiO$_3$ (mmol m$^{-1}
```

Surface macronutrients



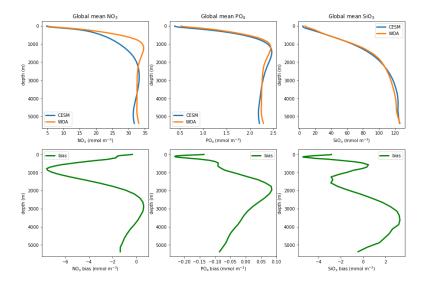
# Global nutrient profiles

```
In [19]: ds_glb = xr.Dataset({v: cesm_ann_ds[v] for v in ['z_t']})
    ds_glb['N03']= esmlab.weighted_mean(cesm_ann_ds['N03'], weights=area, d:
    ds_glb['P04']= esmlab.weighted_mean(cesm_ann_ds['P04'], weights=area, d:
    ds_glb['Si03']= esmlab.weighted_mean(cesm_ann_ds['Si03'], weights=area,
```

```
In [66]: ds_glb_woa = xr.Dataset({v: cesm_ann_ds[v] for v in ['z_t']})
    ds_glb_woa['N03'] = esmlab.weighted_mean(ds_woa['N03'], weights=area, d:
    ds_glb_woa['P04'] = esmlab.weighted_mean(ds_woa['P04'], weights=area, d:
    ds_glb_woa['Si03'] = esmlab.weighted_mean(ds_woa['Si03'], weights=area,
```

```
In [67]: fig = plt.figure(figsize=(15,10))
         plt.suptitle('Global mean macronutrient profiles', fontsize=14)
         #COLUMN 1 - NO3
         ax = fig.add subplot(2,3,1)
         ax.set_title('Global mean NO$_3$')
         ax.plot(ds_glb['NO3'].values, depths, label='CESM', linewidth=3)
         ax.plot(ds_glb_woa['NO3'].values, depths, label='WOA', linewidth=3)
         ax.legend()
         ax.set(ylabel='depth (m)',xlabel='NO\$_3\$ (mmol m\$^{-3}\$)')
         plt.gca().invert_yaxis()
         #COLUMN 2 - PO4
         ax = fig.add_subplot(2,3,2)
         ax.set_title('Global mean PO$_4$')
         ax.plot(ds_glb['P04'].values, depths, label='CESM', linewidth=3)
         ax.plot(ds_glb_woa['PO4'].values, depths, label='WOA', linewidth=3)
         ax.set(ylabel='depth (m)',xlabel='PO$_4$ (mmol <math>m$^{-3}$)')
         plt.gca().invert_yaxis()
         #COLUMN 3 - SiO3
         ax = fig.add_subplot(2,3,3)
         ax.set_title('Global mean SiO$_3$')
         ax.plot(ds glb['SiO3'].values, depths, label='CESM', linewidth=3)
         ax.plot(ds_glb_woa['SiO3'].values, depths, label='WOA', linewidth=3)
         ax.legend()
         ax.set(ylabel='depth \ (m)',xlabel='SiO\$_3\$ \ (mmol \ m\$^{-3}\$)')
         plt.gca().invert_yaxis()
         #COLUMN 1 - NO3 diff
         ax = fig.add_subplot(2,3,4)
         ax.plot(ds_glb['N03'].values - ds_glb_woa['N03'].values, depths, label=
         ax.legend()
         ax.set(ylabel='depth (m)',xlabel='NO$_3$ bias (mmol m$^{-3}$)')
         plt.gca().invert_yaxis()
         #COLUMN 2 - PO4 diff
         ax = fig.add_subplot(2,3,5)
         ax.plot(ds_glb['P04'].values - ds_glb_woa['P04'].values, depths, label=
         ax.set(ylabel='depth (m)',xlabel='PO$_4$ bias (mmol m$^{-3}$)')
         plt.gca().invert_yaxis()
         #COLUMN 3 - SiO3 diff
         ax = fig.add_subplot(2,3,6)
         ax.plot(ds_glb['Si03'].values - ds_glb_woa['Si03'].values, depths, label
         ax.legend()
         ax.set(ylabel='depth (m)',xlabel='SiO$_3$ bias (mmol m$^{-3}$)')
         plt.gca().invert_yaxis()
```

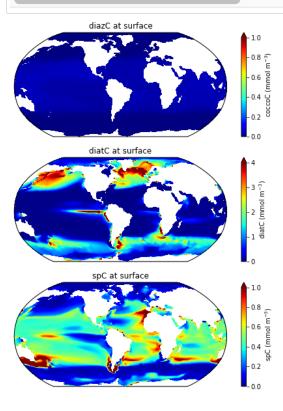




```
In [37]: #most limiting nutrient - concatenate the limitation terms so that nutr.
           \# 0 = PO4
           #1 = Fe
           #2 = NO3 (only for sp and diat)
           # 3 = Si (only for diat)
           limarray_sp=xr.concat((cesm_ann_ds.sp_P_lim_surf, cesm_ann_ds.sp_Fe_lim_
           limarray diat=xr.concat((cesm ann ds.diat P lim surf, cesm ann ds.diat I
           limarray_diaz=xr.concat((cesm_ann_ds.diaz_P_lim_surf, cesm_ann_ds.diaz_I
In [38]: most_lim_sp=limarray_sp.argmin(dim='nutrient', skipna=False).squeeze()
           most_lim_diat=limarray_diat.argmin(dim='nutrient', skipna=False).squeeze
most_lim_diaz=limarray_diaz.argmin(dim='nutrient', skipna=False).squeeze
           mask = np.isnan(cesm_ann_ds.sp_N_lim_surf.squeeze())
In [39]: fig = plt.figure(figsize=(5,9))
           ax = fig.add_subplot(3,1,1, projection=ccrs.Robinson(central_longitude=0)
           ax.set_title('Diat nut lim', fontsize=12)
pc=ax.pcolormesh(lons, lats, most_lim_diat.where(~mask), cmap=plt.cm.get
           colorbar_specs = {'ticks' : np.arange(0,4,1)}
           ax = fig.add_subplot(3,1,2, projection=ccrs.Robinson(central_longitude=0)
           ax.set_title('SP nut lim', fontsize=12)
           pc=ax.pcolormesh(lons, lats, most_lim_sp.where(~mask), cmap=plt.cm.get_c
           colorbar_specs = {'ticks' : np.arange(0,4,1)}
           ax = fig.add_subplot(3,1,3, projection=ccrs.Robinson(central_longitude=:
ax.set_title('Diaz nut lim', fontsize=12)
pc=ax.pcolormesh(lons, lats, most_lim_diaz.where(~mask), cmap=plt.cm.get
           colorbar_specs = {'ticks' : np.arange(0,4,1)}
           fig.subplots_adjust(right=0.8)
           cbar_ax = fig.add_axes([0.85, 0.15, 0.05, 0.7])
           cbar = fig.colorbar(pc, cax=cbar_ax,**colorbar_specs)
           cbar.ax.set_yticklabels(['P lim', 'Fe lim', 'N lim', 'SiO3/C lim']);
                        Diat nut lim
                                                    SiO3/C lim
                                                    N lim
                                                    Fe lim
                                                    - P lim
```

Look at phyto carbon pools

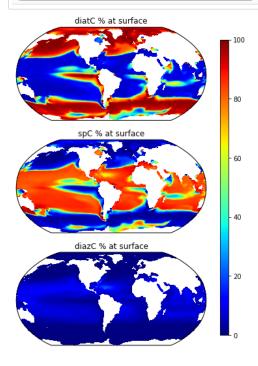
```
In [27]: fig = plt.figure(figsize=(8,10))
    ax = fig.add_subplot(3,1,1, projection=ccrs.Robinson(central_longitude=: ax.set_title('diazC at surface', fontsize=12)
    pc=ax.pcolormesh(lons, lats, cesm_ann_ds.diazC.isel(z_t_150m=0), cmap=':
    cbar1 = fig.colorbar(pc, ax=ax,extend='max',label='coccoC (mmol m$^{-3}{\}$
    ax = fig.add_subplot(3,1,2, projection=ccrs.Robinson(central_longitude=: ax.set_title('diatC at surface', fontsize=12)
    pc=ax.pcolormesh(lons, lats, cesm_ann_ds.diatC.isel(z_t_150m=0), cmap=':
    cbar1 = fig.colorbar(pc, ax=ax,extend='max',label='diatC (mmol m$^{-3}{\}$)
    ax = fig.add_subplot(3,1,3, projection=ccrs.Robinson(central_longitude=: ax.set_title('spC at surface', fontsize=12)
    pc=ax.pcolormesh(lons, lats, cesm_ann_ds.spC.isel(z_t_150m=0), cmap='jet cbar1 = fig.colorbar(pc, ax=ax,extend='max',label='spC (mmol m$^{-3}{\}$)'
```



## Look at percent phytoC

```
In [30]: phytoC = cesm_ann_ds.spC + cesm_ann_ds.diatC + cesm_ann_ds.diazC
    perc_sp = cesm_ann_ds.spC / (phytoC) * 100.
    perc_diat = cesm_ann_ds.diatC / (phytoC) * 100.
    perc_diaz = cesm_ann_ds.diazC / (phytoC) * 100.
```

```
In [62]: fig = plt.figure(figsize=(6,9))
    ax = fig.add_subplot(3,1,1, projection=ccrs.Robinson(central_longitude=: ax.set_title('diatC % at surface', fontsize=12)
    pc=ax.pcolormesh(lons, lats, perc_diat.isel(z_t_150m=0), cmap='jet',vmin
    ax = fig.add_subplot(3,1,2, projection=ccrs.Robinson(central_longitude=: ax.set_title('spC % at surface', fontsize=12)
    pc=ax.pcolormesh(lons, lats, perc_sp.isel(z_t_150m=0), cmap='jet',vmin=(
    ax = fig.add_subplot(3,1,3, projection=ccrs.Robinson(central_longitude=: ax.set_title('diazC % at surface', fontsize=12)
    pc=ax.pcolormesh(lons, lats, perc_diaz.isel(z_t_150m=0), cmap='jet',vmin
    fig.subplots_adjust(right=0.8)
    cbar_ax = fig.add_axes([0.85, 0.15, 0.03, 0.7])
    fig.colorbar(pc, cax=cbar_ax);
```



# **NPP**

```
In [31]: variables = [f'photoC_{phyto}_zint' for phyto in ['diat', 'sp', 'diaz',
                                                                   ds_glb = xr.Dataset()
for v in variables:
                                                                                                \label{eq:cosm_ann_ds} ds\_glb[v] = esmlab.statistics.weighted\_sum(cesm\_ann\_ds[v], weights=\epsilon to the control of the control of the cost of
                                                                                                ds_glb[v].attrs = cesm_ann_ds[v].attrs
                                                                                                 #if ds1 annmean[v].units == 'mmo1/m^3 cm/s':
                                                                                              ds_glb[v] = ds_glb[v] * nmols_to_PgCyr
ds_glb[v].attrs['units'] = 'Pg C yr$^{-1}$'
                                                                    ds_glb = ds_glb.compute()
                                                                    ds_glb
```

Out[31]: xarray.Dataset

- ▶ Dimensions:
- ► Coordinates: (0)
- ▼ Data variables:

```
photoC_diat_zint () float64 21.37
                                                               photoC_sp_zint () float64 22.43
                                                               photoC_diaz_zint () float64 1.607
                                                               photoC_TOT_zint () float64 45.41
```

► Attributes: (0)

Globally integrated NPP should be between 50 and 60 Pg C  ${\rm yr}^{-1}$  according to satellite derived NPP algorithms

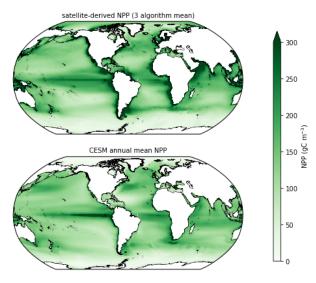
### Compare to satellite-derived NPP

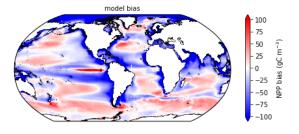
```
In [9]: file='/glade/work/kristenk/satellite data/POP regrid/NPP 3methodmean mea
        ds_npp=xr.open_dataset(file, decode_times=False)
```

In [10]: npp = cesm\_ann\_ds.photoC\_TOT\_zint \* 0.01 \* 86400 \* 365 \* 0.001 \* 12. #cc

```
In [33]: #plot depth integrated NPP
         fig = plt.figure(figsize=(8,7))
         plt.suptitle('depth integrated NPP', fontsize=14)
         ###########NO3
         #COLUMN 1 - NO3
         #---- CESM panel
         ax = fig.add_subplot(2,1,2, projection=ccrs.Robinson(central_longitude=
         ax.coastlines('10m',linewidth=0.5)
ax.set_title('CESM annual mean NPP', fontsize=10)
         pc1=ax.pcolormesh(lons, lats,
                           npp, cmap='Greens',
                            vmin=0, vmax=300,
                           transform=ccrs.PlateCarree())
         #--- OBS panel
         ax = fig.add_subplot(2,1,1, projection=ccrs.Robinson(central_longitude=
         ax.coastlines('10m',linewidth=0.5)
         ax.set_title('satellite-derived NPP (3 algorithm mean)', fontsize=10)
         pc2=ax.pcolormesh(lons, lats,
                           ds_npp.NPP, cmap='Greens',
                           vmin=0, vmax=300,
                           transform=ccrs.PlateCarree())
         fig.subplots_adjust(right=0.8)
         cbar_ax = fig.add_axes([0.85, 0.15, 0.02, 0.7])
         fig.colorbar(pc2, cax=cbar_ax,extend='max',label='NPP (gC m$^{-3}$)');
```

#### depth integrated NPP





#### Calcification

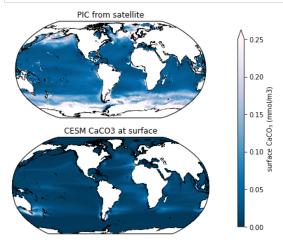
In [22]: ## GLOBALLY integrated calcification; observation-based estimates range
 ## (Feely et al., 2004, Maranon et al., 2016, Smith et al., 2016, Balch
 ds\_glb['CaCO3\_PROD\_zint'] = esmlab.statistics.weighted\_sum(cesm\_ann\_ds[
 ds\_glb['CaCO3\_PROD\_zint'] = ds\_glb['CaCO3\_PROD\_zint'] \* nmols\_to\_PgCyr
 ds\_glb['CaCO3\_PROD\_zint'].values
 print('Globally integrated calcification is',ds\_glb['CaCO3\_PROD\_zint'].v
 print('(should be between 0.6 to 2.4 Pg C per year)')

Globally integrated calcification is 0.5967147978656123 Pg C per year (should be between 0.6 to 2.4 Pg C per year)

# Compare satellite-derived Particulate Inorganic Carbon (PIC; i.e., ${\sf CaCO}_3$ ) to ${\sf spCaCO}_3$ in the top 10m

In [23]: #read in mean PIC
 pic\_file = '/glade/work/kristenk/satellite\_data/processed/PIC\_mean2003tc
 ds\_pic=xr.open\_dataset(pic\_file, decode\_times=False)
 #convert to mmol/m3 from mg/m3
 ds\_pic['PIC']=ds\_pic.PIC / 12.011

In [35]: fig = plt.figure(figsize=(7,6))
 ax = fig.add\_subplot(2,1,1, projection=ccrs.Robinson(central\_longitude=: ax.set\_title('PIC from satellite', fontsize=12)
 ax.coastlines('10m',linewidth=0.5)
 pc=ax.pcolormesh(ds\_pic.lon.values, ds\_pic.lat.values, ds\_pic.PIC, cmap=
 ax = fig.add\_subplot(2,1,2, projection=ccrs.Robinson(central\_longitude=: ax.set\_title('CESM CaCO3 at surface', fontsize=12)
 ax.coastlines('10m',linewidth=0.5)
 pc=ax.pcolormesh(lons, lats, cesm\_ann\_ds.spCaCO3.isel(z\_t\_150m=0), cmap=
 fig.subplots\_adjust(right=0.8)
 cbar\_ax = fig.add\_axes([0.85, 0.15, 0.02, 0.7])
 fig.colorbar(pc, cax=cbar\_ax,extend='max',label='surface CaCO\$\_3\$ (mmol, other cax is a contracted by the color cax is a contracted by the cax is a contracted by the color cax is a contracted by the color cax is a contracted by the cax is a contracted by the color cax is a contracted by the c

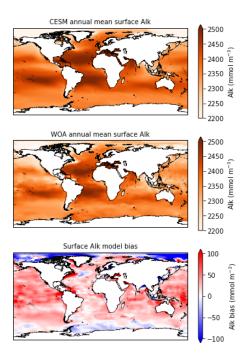


# **Alkalinity**

In [42]: #just getting this for the coords (that don't have nans)
 file = '/glade/work/kristenk/GLODAPv2\_regridded/glodap\_pop\_grid/GLODAP\_1
 ds\_glodap = xr.load\_dataset(file, decode\_times=False, decode\_coords=False, ds\_glodap['z\_t']=cesm\_ann\_ds.z\_t
 ALK\_diff = cesm\_ann\_ds.ALK - ds\_glodap.ALK
 #rmse\_global['ALK']=esmlab.statistics.weighted\_rmsd(cesm\_ann\_ds.ALK, ds\_mrmse\_global.ALK.values

```
In [63]: #plot surface alkalinity
         fig = plt.figure(figsize=(6,9))
         plt.suptitle('Surface alkalinity', fontsize=14)
         ###########NO3
         #COLUMN 1 - NO3
         #---- CESM panel
         ax = fig.add_subplot(3,1,1, projection=ccrs.PlateCarree())
         ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set_title('CESM annual mean surface Alk', fontsize=10)
         pc1=ax.pcolormesh(lons, lats,
                           cesm_ann_ds.ALK.isel(z_t=0), cmap='Oranges',
                            vmin=2200, vmax=2500,
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc1, ax=ax,extend='max',label='Alk (mmol m$^{-3}$)
         #--- OBS panel
         ax = fig.add_subplot(3,1,2, projection=ccrs.PlateCarree())
         ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set_title('WOA annual mean surface Alk', fontsize=10)
         pc2=ax.pcolormesh(lons, lats,
                           ds_glodap.ALK.isel(z_t=0), cmap='Oranges',
                            vmin=2200, vmax=2500,
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc2, ax=ax,extend='max',label='Alk (mmol m$^{-3}$)
         #--- DIFF panel
         ax = fig.add_subplot(3,1,3, projection=ccrs.PlateCarree())
ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
         ax.coastlines('10m',linewidth=0.5)
         ax.set_title('Surface Alk model bias', fontsize=10)
         pc3=ax.pcolormesh(lons, lats,
                           ALK_diff.isel(z_t=0), vmin=-100, vmax=100, cmap='bwr',
                           transform=ccrs.PlateCarree())
         cbar1 = fig.colorbar(pc3, ax=ax,extend='both',label='Alk bias (mmol m$^-
```

#### Surface alkalinity



```
In [70]: # integrate SiO2 production over top 150 m
           SiO2_prod_int = cesm_ann_ds['SiO2_PROD'].isel(z_t=slice(0,15)) * 10. #ea
           Si02_prod_int = Si02_prod_int.sum(dim='z_t'), #units go from mmolSi/m3/s
Si02_prod_int = Si02_prod_int * 86400. * 365. * 0.001 #convert to molSi,
In [71]: fig = plt.figure(figsize=(11,5))
           #---- CESM panel
           ax = fig.add_subplot(1,1,1, projection=ccrs.PlateCarree())
ax.set_extent([-180, 180, -90, 90], ccrs.PlateCarree())
           ax.coastlines('10m',linewidth=0.5)
           ax.set_title('CESM 150m-integrated SiO$_2$ production', fontsize=10)
           pc1=ax.pcolormesh(lons, lats,
                                SiO2_prod_int,
                                 vmin=0, vmax=1,
                                 cmap='BuGn',
                                transform=ccrs.PlateCarree())
           cbar1 = fig.colorbar(pc1, ax=ax,extend='max',label='SiO$_2$ Production
                                  CESM 150m-integrated SiO<sub>2</sub> production
                                                                                              m^{-2} y^{-1}
                                                                                              SiO
                                                                                          0.2 🖔
In [73]: ### GLOBALLY integrated biogenic SiO2 production is between 166 and 280
           gc_si_prod = SiO2_prod_int * area_m2 #molSi/gc/yr
           ds_glb['SiO2_PROD'] = gc_si_prod.sum(dim='nlon').sum(dim='nlat') * 1.e-
           ds_glb['SiO2_PROD'].values
Out[73]: array(111.15986714)
           GLOBALLY integrated biogenic SiO2 production is between 166 and 280 mol Si per year, but
           half is dissolved in the upper 100m (Nelson et al., 1995; Holzer et al., 2014)
           POC flux at 100m
In [61]: ds_glb['POC_FLUX_100m'] = esmlab.statistics.weighted_sum(cesm_ann_ds['POC_GLUX_100m'] = ds_glb['POC_FLUX_100m'] * nmols_to_PgCyr
           ds_glb['POC_FLUX_100m'].values
           print('Globally integrated POC flux is',ds_glb['POC_FLUX_100m'].values,
           Globally integrated POC flux is 7.05632704198471 Pg C per year
 In [ ]:
 In [ ]:
```

In [ ]:

In [ ]:

In [ ]: