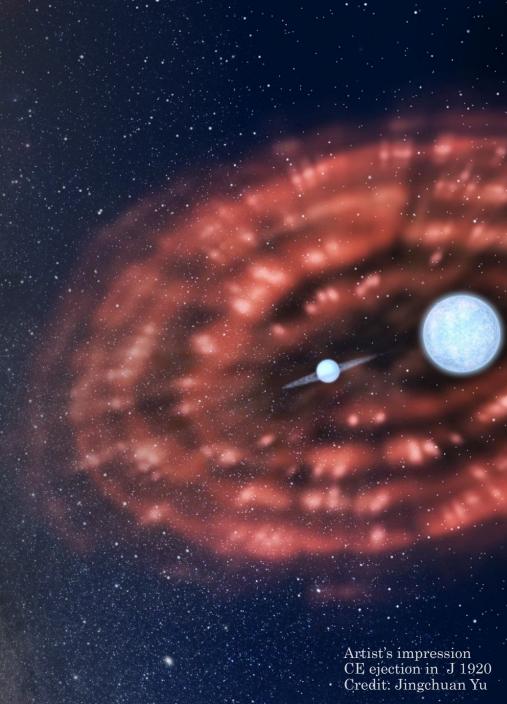
Close white dwarfs with unevolved companions: constraints on the common envelope phase

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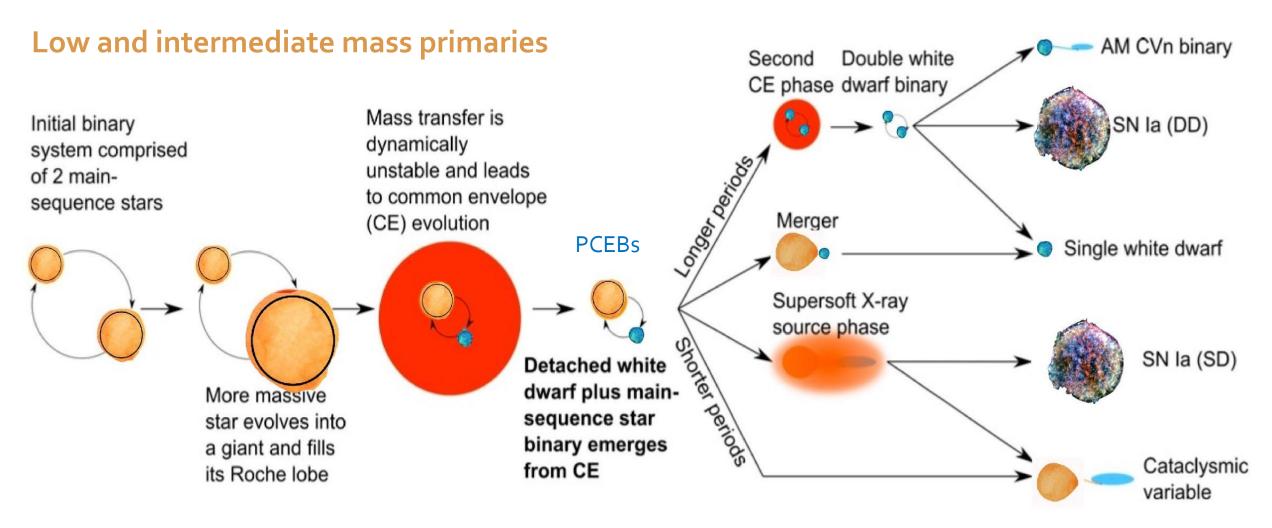






#### **CE** evolution

Paczynski 1976

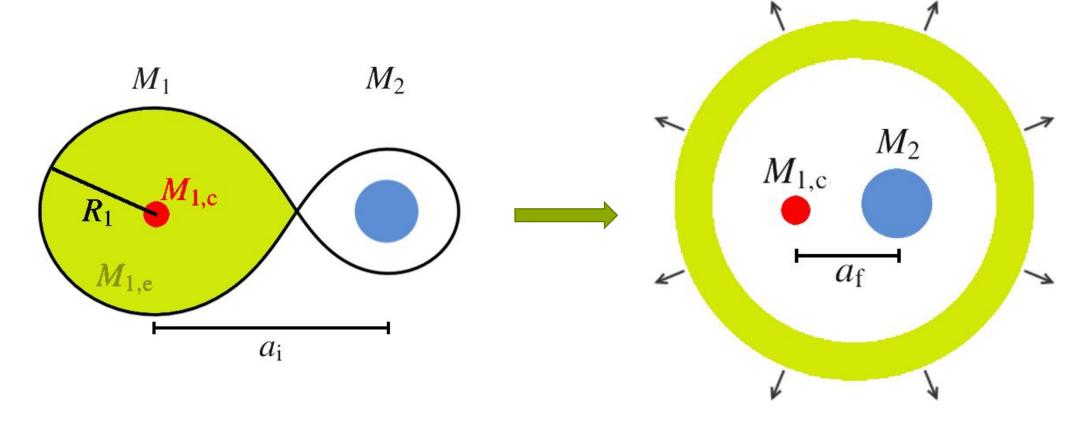


Adapted by M. S. Hernandez from Kulkarni et al. (2021)

## **Energy formalism**

Webbink 1984, Livio and Soker 1988

$$E_{\rm bind} = \alpha_{\rm CE} \Delta E_{\rm orb}$$



$$E_{\text{orb,i}} = \frac{1}{2} \frac{GM_{1,c}M_2}{a_{i}}$$
;  $E_{\text{bind}} = -\frac{GM_1M_{1,e}}{\lambda R_1}$ 

$$E_{\text{orb,f}} = \frac{1}{2} \frac{GM_{1,c}M_2}{a_{\text{f}}}$$

## **Energy formalism**

Webbink 1984, Livio and Soker 1988

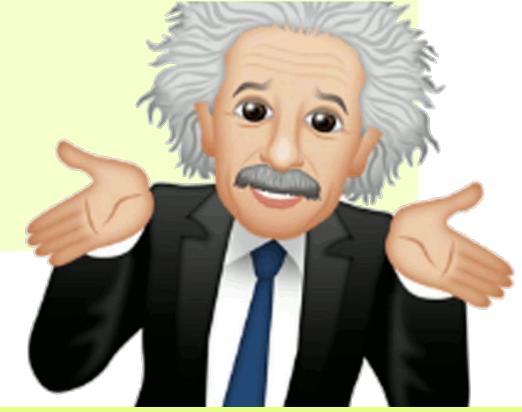


# Is $\alpha_{CE}$ a universal constant?

What is  $\lambda$ ?

Careful with E orb,i (some authors use M<sub>1</sub> here)

$$E_{\text{orb,i}} = \frac{1}{2} \frac{GM_{1,c}M_2}{a_{\text{i}}}; E_{\text{bind}} = -\frac{GM_1M_{1,e}}{\lambda R_1}$$



## The λ parameter

Accounts for the "real" binding energy of the envelope  $E_{\rm bind} = -\frac{GM_1M_{1,\rm e}}{\lambda~R}$ 

$$E_{\rm bind} = -\frac{GM_1M_{1,\epsilon}}{\lambda R_1}$$

Different studies assumed:

- Fixed  $\lambda$  value (typically 0.5 or 1.0)
- λ calculated from pure gravitational energy of the envelope

$$-\frac{GM_1M_{1,e}}{\lambda R_1} = -\int_{M_{1,e}}^{M_1} \frac{Gm}{r(m)} dm$$

• λ calculated from gravitational + a fraction of "internal" energy:

$$-\frac{GM_1M_{1,e}}{\lambda R_1} = \int_{M_{1,e}}^{M_1} \left( -\frac{Gm}{r(m)} + \alpha_{\text{int}}U_{\text{int}}(m) \right) dm$$

## The λ parameter

Internal energy can include:

- Thermal energy
- Recombination
- Radiative losses
- Enthalpy
- Convective transport

• ...

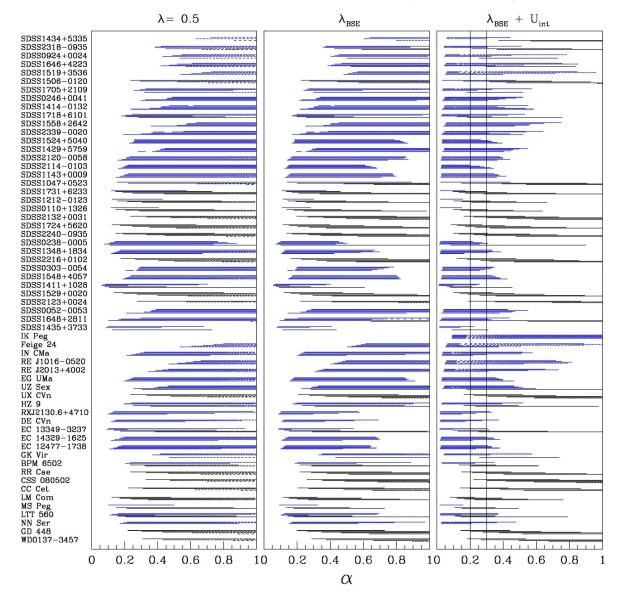
(e.g. Dewi & Tauris 2000; Ivanova et al 2013; Wilson & Nordhaus 2022)

all with different efficiencies!



$$E_{\text{bind}} = -\frac{GM_1M_{1,e}}{\lambda R_1} = \int_{M_{1,e}}^{M_1} \left( -\frac{Gm}{r(m)} dm + \alpha_{\text{th}} U_{\text{th}} + \alpha_{\text{rec}} U_{\text{rec}} + \dots \right) dm$$

## Zorotovic et al. (2010)



• Left:  $\lambda = 0.5$ 

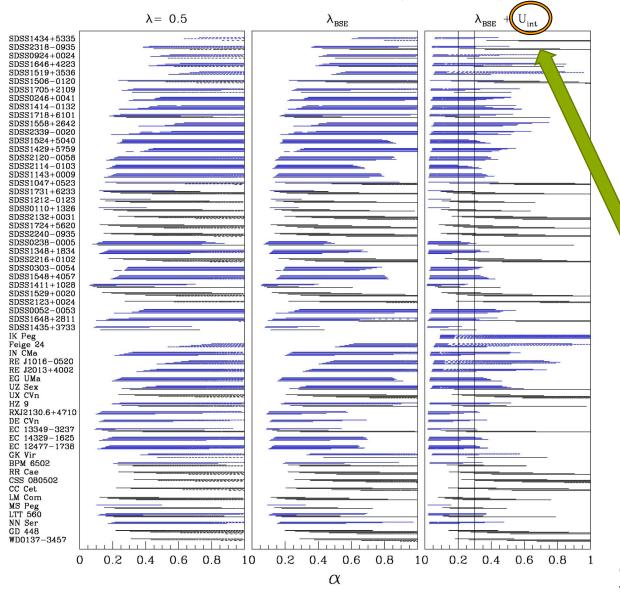
$$E_{\text{bind}} = \frac{GM_1M_{1,e}}{0.5R_1}$$

• Middle:  $\lambda$  from BSE code (Hurley et al 2002) without  $U_{rec}$ 

• Right:  $\lambda$  from BSE code with a fraction  $\alpha_{rec} = \alpha_{CE}$  of  $U_{rec}$ 

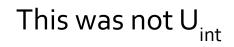
$$\alpha_{CE} \sim 0.25$$

Zorotovic et al. (2010)





"I suppose I'll be the one to mention the elephant in the room."

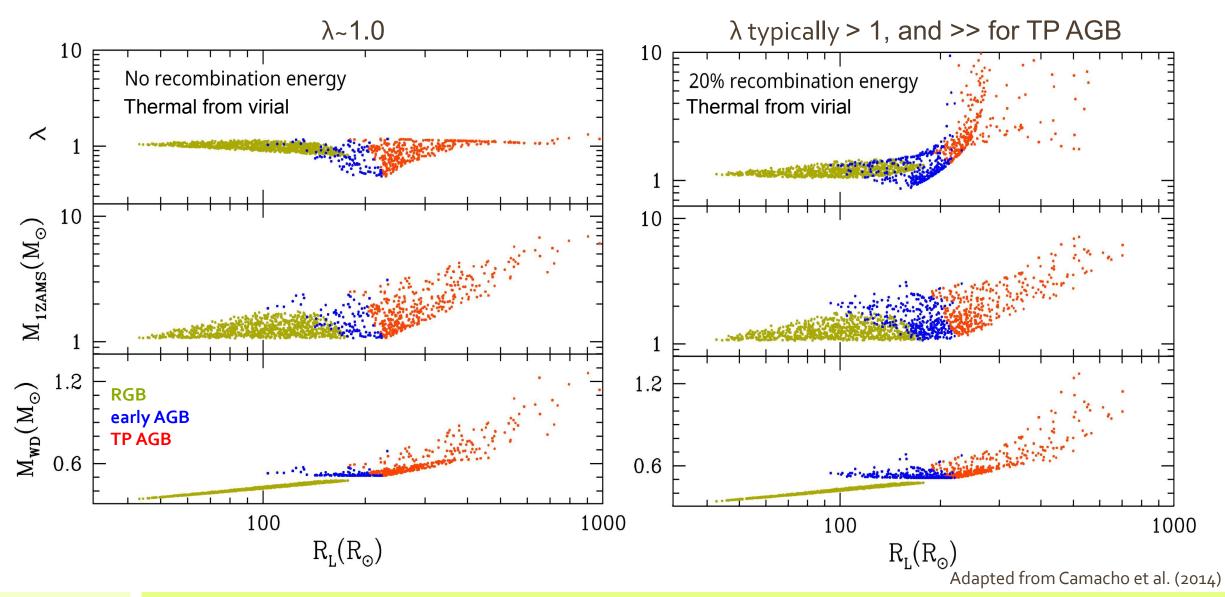


It was  $\mathbf{U}_{\text{rec}}$ 

U<sub>th</sub> included (virial)\*

\*not clarified in the BSE code, nor in Robert Izzards's PhD thesis (2004) or Claeys et. al (2014) where the fitting formulae was published. But  $\lambda$  is multiplied by 2 at the end, consistent with virial theorem (e.g. De Marco et al. 2011)

### λ from BSE



#### Close WD+dM binaries

• From reconstruction of observed systems:  $\alpha_{CF}$  ~0.25 from Zorotovic et. al (2010) was valid for  $\alpha_{CF} = \alpha_{rec}$ 

25% of recombination energy seems unrealistic!

2010A&A...520A..86Z

2010/09

Post-common-envelope binaries from SDSS. IX: Constraining the common-envelope

efficiency

Zorotovic, M.; Schreiber, M. R.; Gänsicke, B. T. and 1 more

#### Close WD+dM binaries

• From reconstruction of observed systems:  $\alpha_{CE}$  ~0.25 from Zorotovic et. al (2010) was valid for  $\alpha_{CE}$  =  $\alpha_{rec}$ 

25% of recombination energy seems unrealistic!

#### However....

• From simulations vs observations (WD+dM SDSS sample):

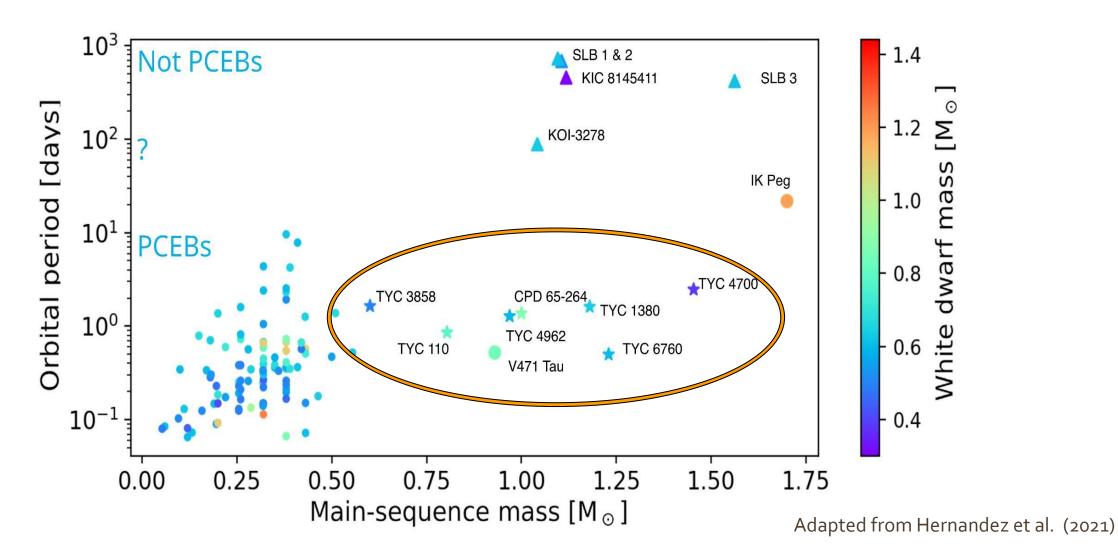
Camacho et al. (2014) favored models with  $\alpha_{CE} \sim 0.3$  and no  $U_{rec}$  (or at least a very small fraction)

Toonen & Nelemans (2013) also favored low  $\alpha_{CE}$  ( $\alpha_{CE}\lambda$  ~ 0.25, consistent if  $\lambda$  ~ 1, i.e. no recombination)



#### **Close WD+AFGK binaries**

The white dwarf binary pathways survey



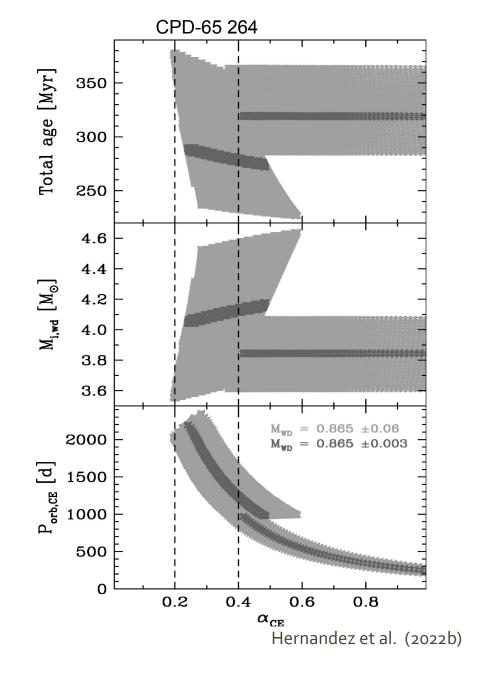
TYC 6760, "The first pre-supersoft X-ray binary" (Parsons et al 2015)  $\alpha_{CF} = 0.08-0.89$  for AGB progenitor (most likely) and > 0.9 for RGB

Object	$\alpha_{\mathrm{CE}}$	WD mass $(M_{\odot})$	$M_{1,\mathrm{i}}~(\mathrm{M}_\odot)$	$P_{\text{orb,i}}$ (d)	Age (Gyr)
TYC 4700	0.2-0.3	0.40-0.44	1.30-1.53	203–386	2.78-4.81
	0.18 - 1.0	0.38-0.44	1.29-1.90	109-395	
TYC 1380	0.2 - 0.3	0.64-0.85	2.54-3.94	706-2222	0.22 - 0.77
	0.10 - 1.0	0.64-0.85	2.36-4.06	126-2222	
TYC 4962	0.2 - 0.3	0.59-0.77	2.28-3.36	617-1917	0.35-1.04
	0.07-1.0	0.59-0.77	1.75-3.46	103-1917	
				Hernandez e	et al. (2021

Object	$\alpha_{\mathrm{CE}}$	$WD \; mass \; (M_{\odot})$	$M_{1,\mathrm{i}}\;(\mathrm{M}_{\odot})$	$P_{\text{orb, i}}$ (d)	CE age (Gyr)
TYC 110 (HST)	0.2-0.3	0.75-0.82	3.24-3.71	929-1947	0.26-0.38
	0.19-1.0	0.75 - 0.82	3.15-3.75	147-2015	0.26-0.42
TYC 3858	0.2-0.3	0.34-0.62	0.98-2.11	103-1477	1.29-13.29
	0.03-1.0	0.27-0.68	0.98 - 2.98	22-1591	0.49-13.29

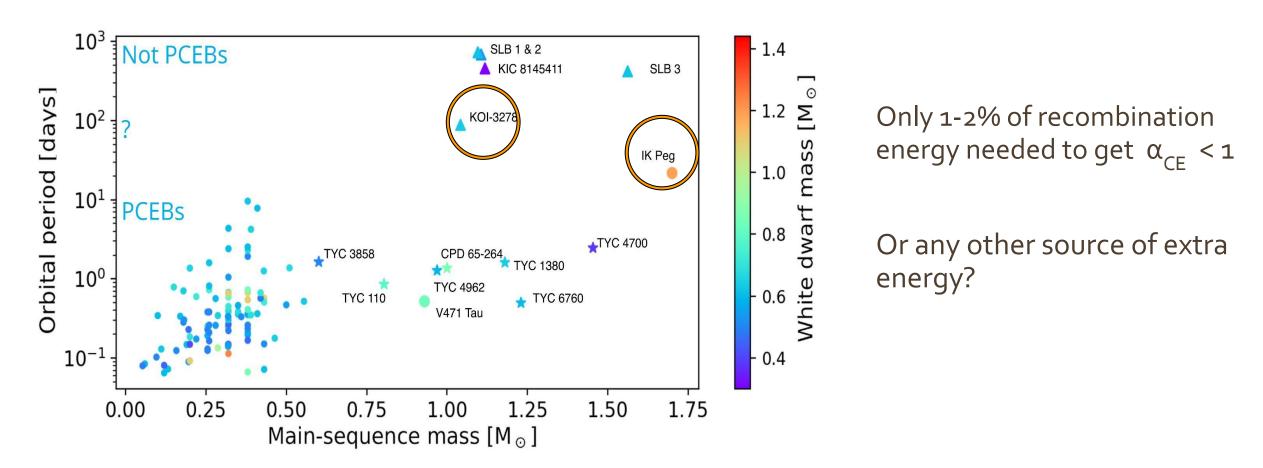
Hernandez et al. (2022a)

All systems wit P orb < 3 days No recombination energy needed (thermal included)  $\alpha_{CF} \gtrsim \text{0.2 works for all}$ 



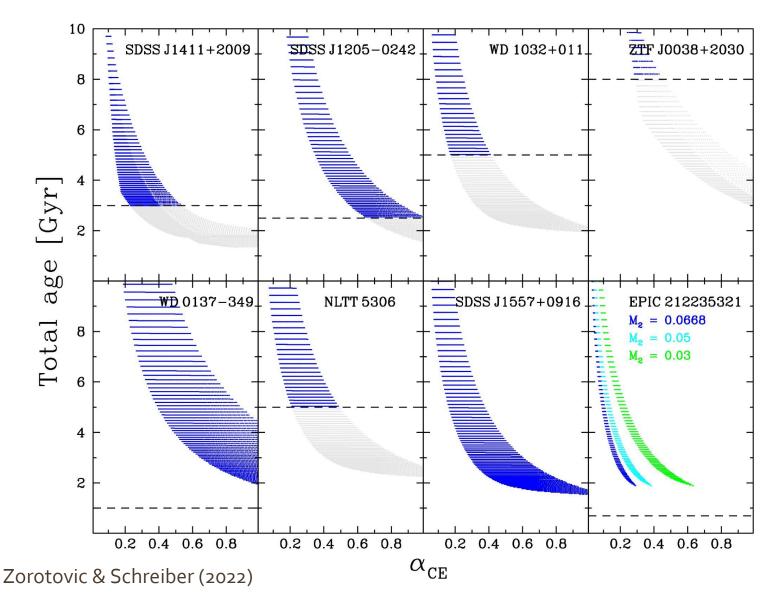
## IK Peg and KOI 3287

**Outliers?** 



Adapted from Hernandez et al. (2021)

#### Close WD+BD binaries

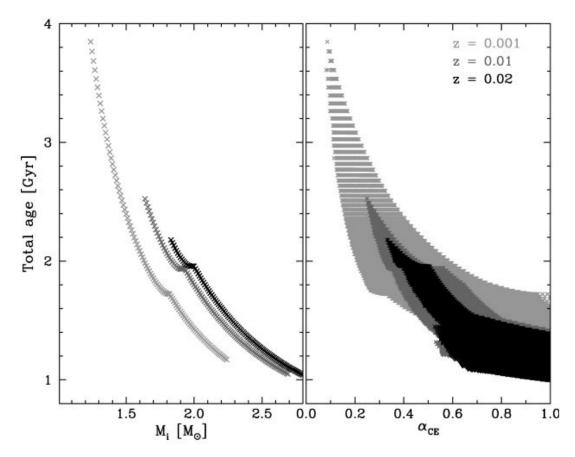


7 systems with  $P_{orb}$  < 0.1 days 1 system with  $P_{orb}$  ~ 0.43 days

No recombination energy needed (thermal included)

 $\alpha_{CE} = 0.24 - 0.41$  works for all when considering age restrictions (dashed lines)

## **Metallicity effects**



**Figure 6.** Total age of the system as a function of initial mass of the progenitor of the white dwarf (*left*) and common-envelope efficiency  $\alpha_{CE}$  (*right*) derived from our reconstruction for three different assumptions of the initial metallicity (z). A white dwarf mass of  $0.6 - 0.7 M_{\odot}$  and a brown dwarf mass of  $0.085 M_{\odot}$  were assumed in this figure.

Casewell et al. (in prep)

Work in progress for GD 1400 Another WD+BD with P<sub>orb</sub> ~ 0.4 days

Lower z allows for smaller initial masses

- $\rightarrow$  smaller  $\alpha_{CF}$  (less mass in the envelope)
- $\rightarrow$  longer evolutionary timescale (older systems).

## **Summary**

- We need to be consistent in the treatment of  $\,\lambda$  in order to constrain  $\alpha_{_{CE}}$
- Recombination energy is not needed for the vast majority of systems (only 1-2% needed to explain outliers, could be something else)
- WD+dM PCEBs:  $\alpha_{CF} \sim 0.3$
- WD+AFGK PCEBs:  $\alpha_{CE} \gtrsim$  0.2
- WD+BD PCEBs:  $\alpha_{CE} = 0.24 0.41$
- WD+WDs:  $\alpha_{CF} \sim 1/3$  ? (see Peter Scherbak's talk this afternoon)
- Metallicity matters!

