

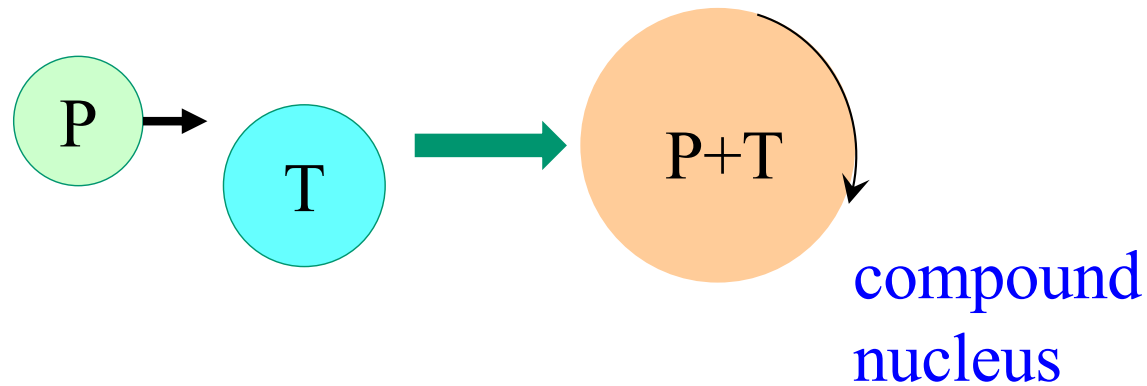
# Theoretical modeling of C+C fusion reactions : Open issues

Kouichi Hagino  
Kyoto University, Kyoto, Japan

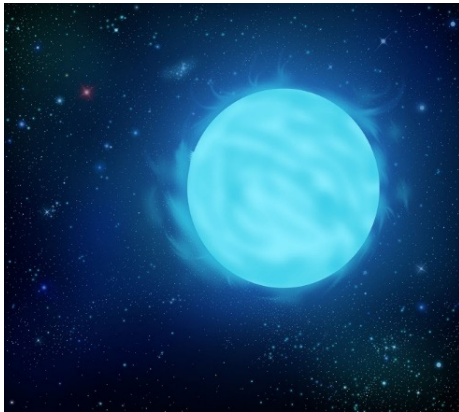


1. Introduction: fusion reactions
2. Summary of theoretical calculations for  $^{12}\text{C}+^{12}\text{C}$  fusion
3. Role of CN resonances
4. Fusion hindrance
5. Summary

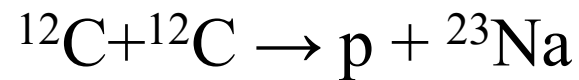
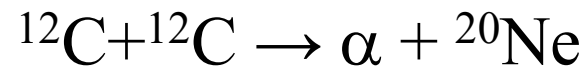
# Introduction: Fusion reactions



$^{12}\text{C}+^{12}\text{C}$  fusion : a key reaction in nuclear astrophysics



✓ Carbon burning in massive stars

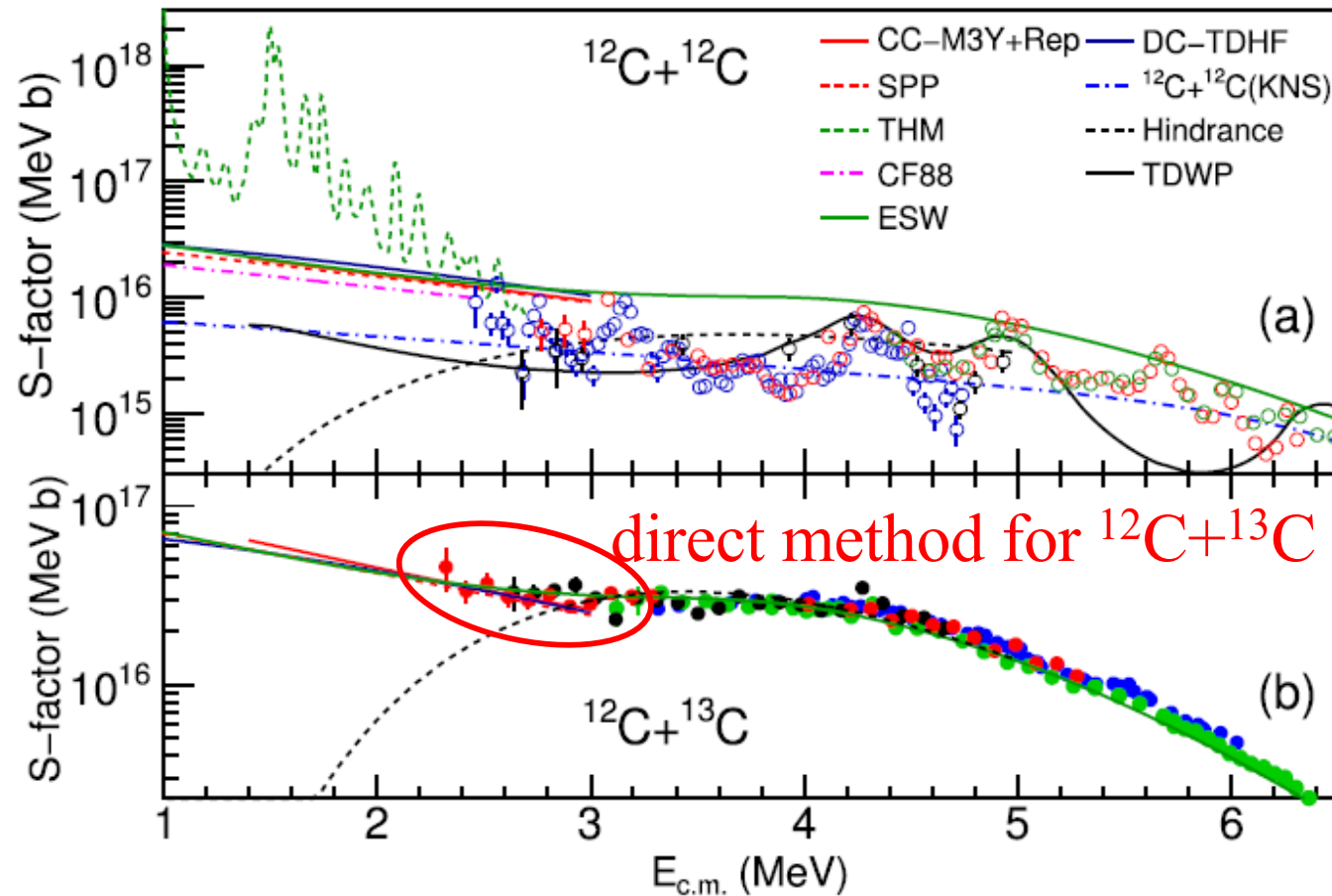


✓ Type Ia supernovae

✓ X-ray superburst

**Fusion and fission:** large amplitude motions & many-body tunneling  
← microscopic understanding: **an ultimate goal of nuclear physics**

# $^{12}\text{C}+^{12}\text{C}$ fusion reaction



N.T. Zhang, et al., Phys. Lett. B801 (2020) 135170

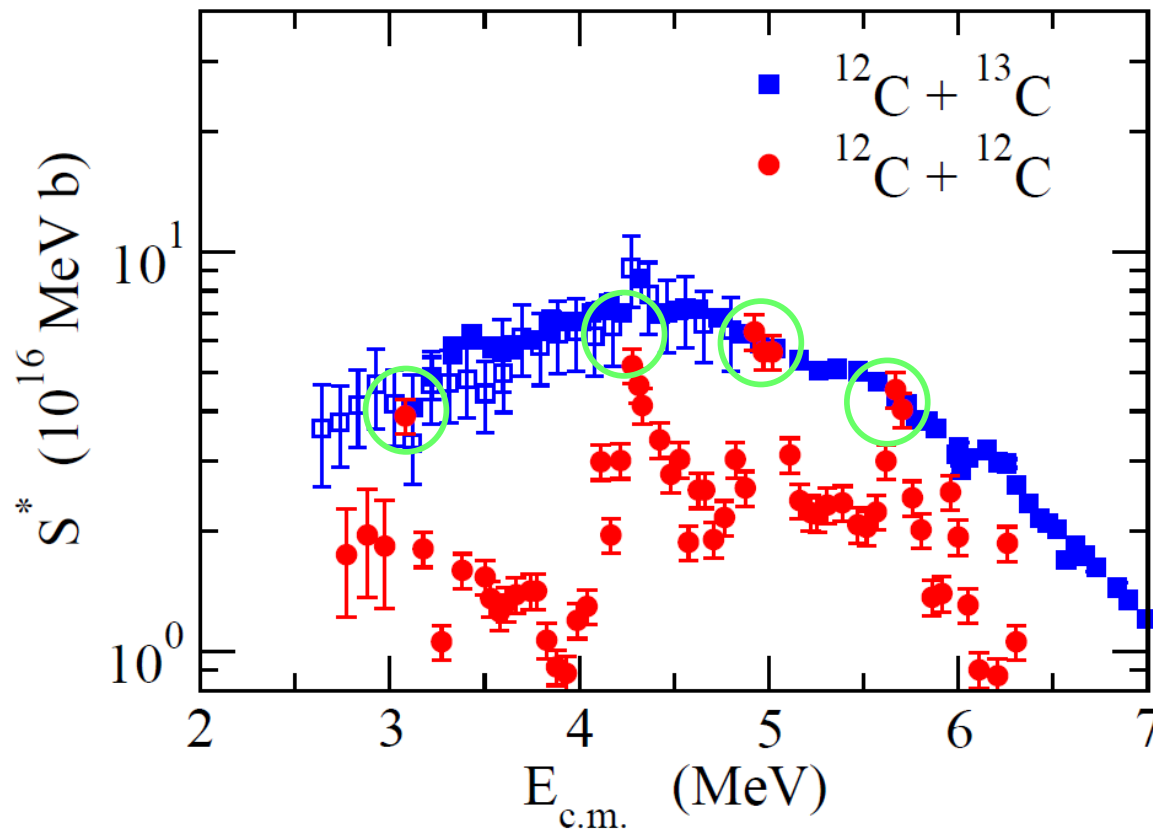
**an important issue:** can one understand  $^{12}\text{C}+^{12}\text{C}$  and  $^{12}\text{C}+^{13}\text{C}$  in a unified manner?

# $^{12}\text{C}+^{12}\text{C}$ fusion reaction

**An open issue:**

Does  $^{12}\text{C}+^{13}\text{C}$  provide the upperlimit of  $^{12}\text{C}+^{12}\text{C}$  cross sections?

→ this has to be answered theoretically



off-resonance:

fusion inhibition

on-resonance:

match with  $^{12}\text{C}+^{13}\text{C}$

M. Notani, X.D. Tang  
et al.,

PRC85('12)014607

## Previous theoretical calculations for $^{12}\text{C}+^{12}\text{C}$

### □ Coupled-channels approaches

- ✓ B. Imanishi, PL 27B ('68) 267, NPA125 ('69) 33.
- ✓ Y. Kondo, T. Matsuse, Y. Abe, PTP59 ('78) 465.
- ✓ W. Scheid, W. Greiner, R. Lemmer, PRL25 ('70) 176.
- ✓ H. Esbensen, X. Tang, and C.L. Jiang, PRC84, 064613 (2011).
- ✓ M. Assuncao and P. Descouvemont, PLB723, 355 (2013).

### □ Time-dependent wave packet approach

- ✓ A. Diaz-Toress and M. Wiesher, PRC97, 055802 (2018).

### □ TDHF-based approaches

- ✓ Imaginary-time evolution: A. Bonasera and J.B. Natowitz, PRC102
- ✓ ATDHF: P.-H. Heenen, PL99B, 298 (1981).
- ✓ Density-Constrained Hartree-Fock: K. Godbey et al., PRC100

### □ Anti-symmetrized Molecular Dynamics (AMD) approach

Y. Taniguchi and M. Kimura, PLB823, 136790 (2021).

AMD + R-matrix

## Previous theoretical calculations for $^{12}\text{C}+^{12}\text{C}$

### □ Coupled-channels approaches

- ✓ B. Imanishi, PL 27B ('68) 267, NPA125 ('69) 33.

#### Nogami-Imanishi model

- ✓ Y. Kondo, T. Matsuse, Y. Abe, PTP59 ('78) 465.

#### Band-crossing model

- ✓ W. Scheid, W. Greiner, R. Lemmer, PRL25 ('70) 176.

#### Double resonance model

- ✓ H. Esbensen, X. Tang, and C.L. Jiang, PRC84, 064613 (2011).

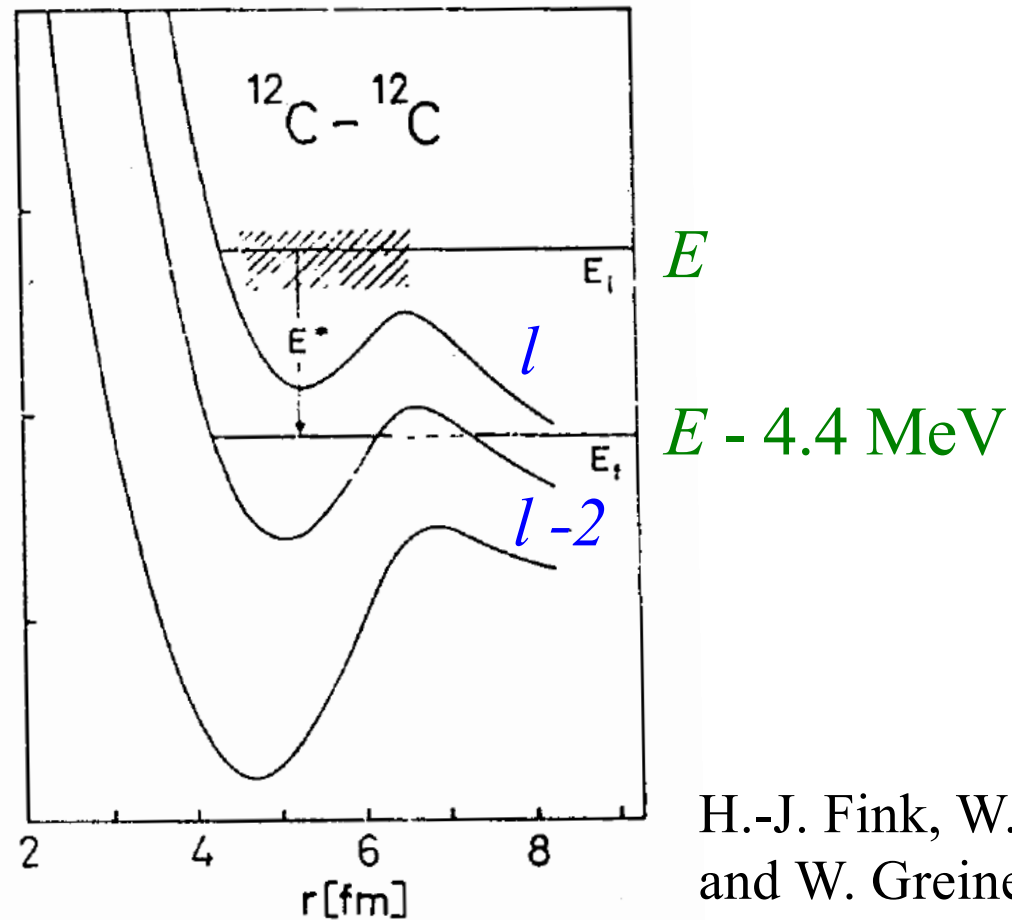
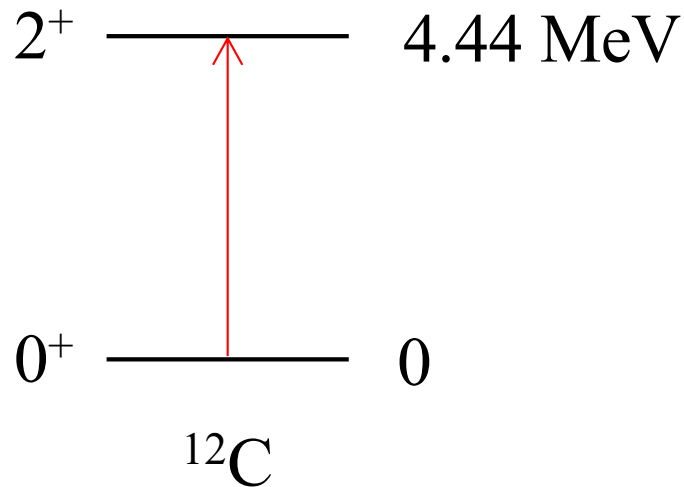
#### Systematic studies for three C+C systems

- ✓ M. Assuncao and P. Descouvemont, PLB723, 355 (2013).

#### Role of the Hoyle state

## Previous calculations for $^{12}\text{C}+^{12}\text{C}$

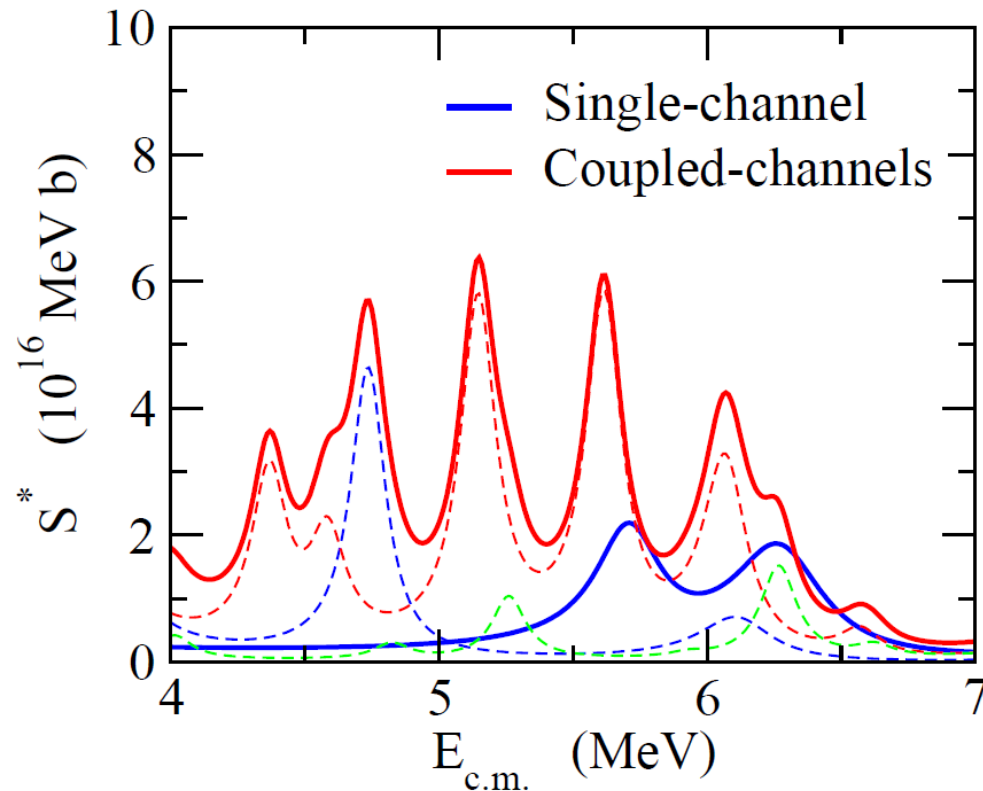
- Nogami-Imanishi model (B. Imanishi, PL 27B ('68) 267, NPA125 ('69) 33)
  - Band-crossing model (Y. Kondo, T. Matsuse, Y. Abe, PTP59 ('78)465)
  - Double resonance model (W. Scheid, W. Greiner, R. Lemmer, PRL25 ('70) 176)
- \* the basic concept is all same



H.-J. Fink, W. Scheid,  
and W. Greiner,  
NPA188 ('72) 259

## Previous calculations for $^{12}\text{C}+^{12}\text{C}$

- Nogami-Imanishi model (B. Imanishi, PL 27B ('68) 267, NPA125 ('69) 33)
  - Band-crossing model (Y. Kondo, T. Matsuse, Y. Abe, PTP59 ('78)465)
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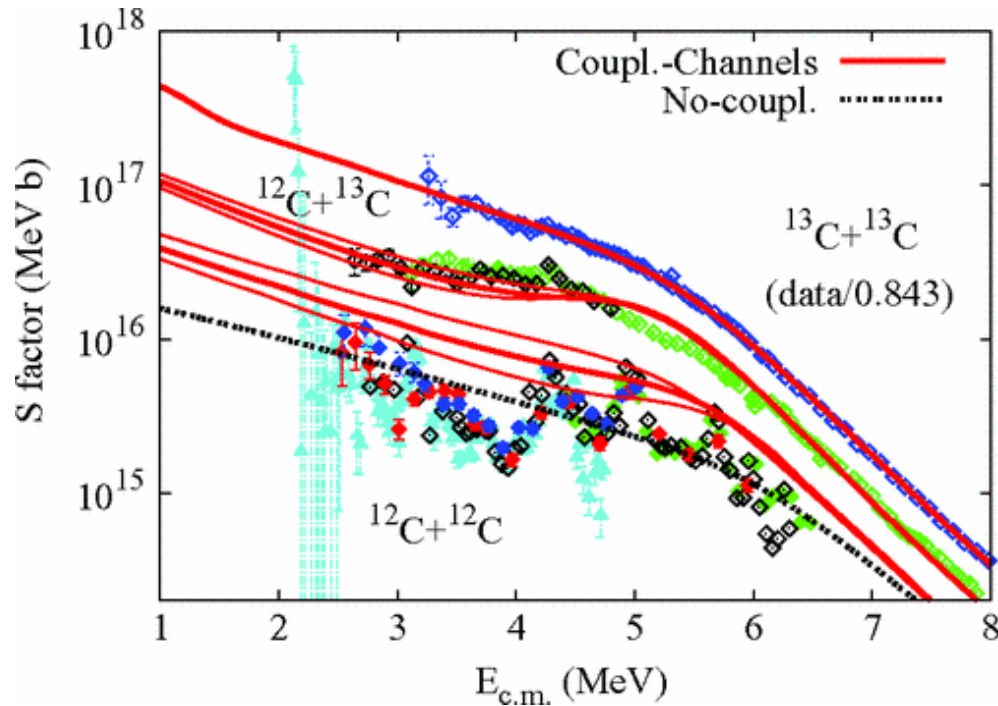
Analogy to Feshbach resonances  
in cold atoms!

Y. Kondo, T. Matsuse, and Y. Abe,  
PTP 59 ('78) 465



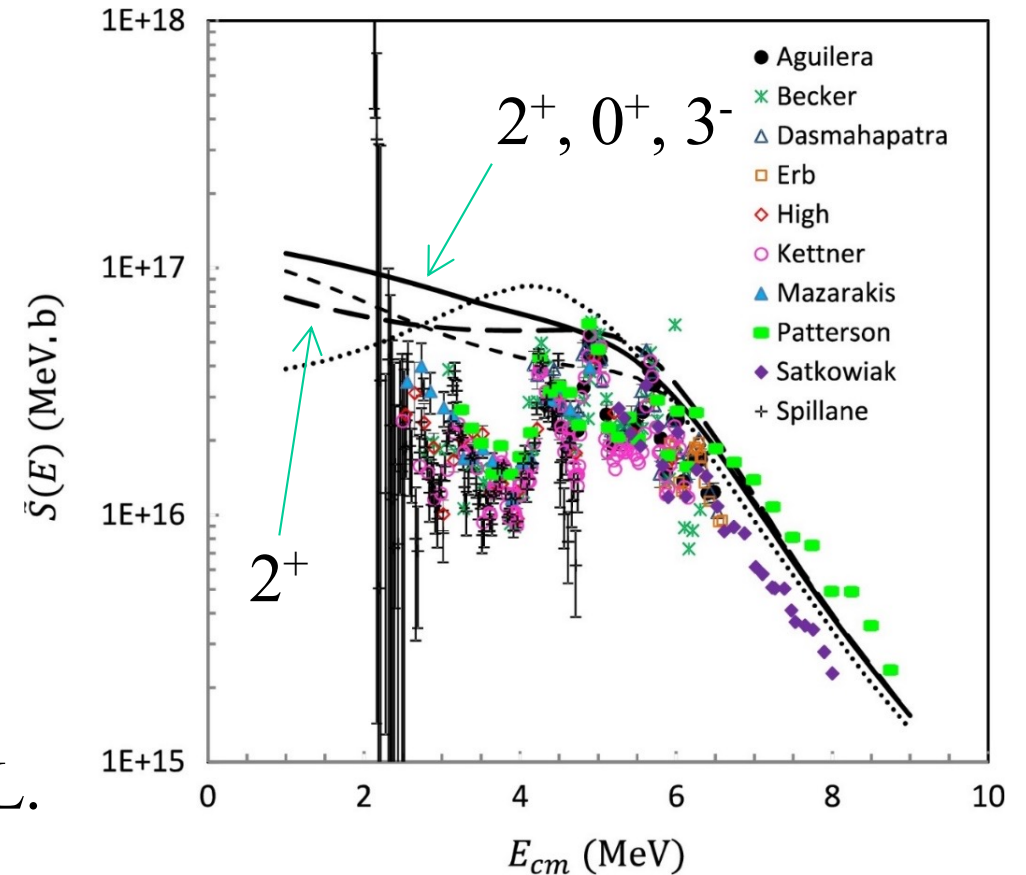
# More recent coupled-channels calculations

## Systematic study



H. Esbensen, X. Tang, and C.L. Jiang, PRC84, 064613 (2011).

## role of the Hoyle state

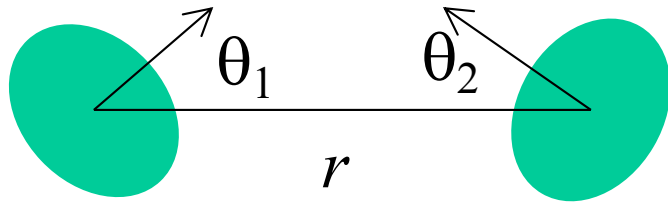


M. Assuncao and P. Descouvemont, PLB723, 355 (2013).

# Previous theoretical calculations for $^{12}\text{C}+^{12}\text{C}$

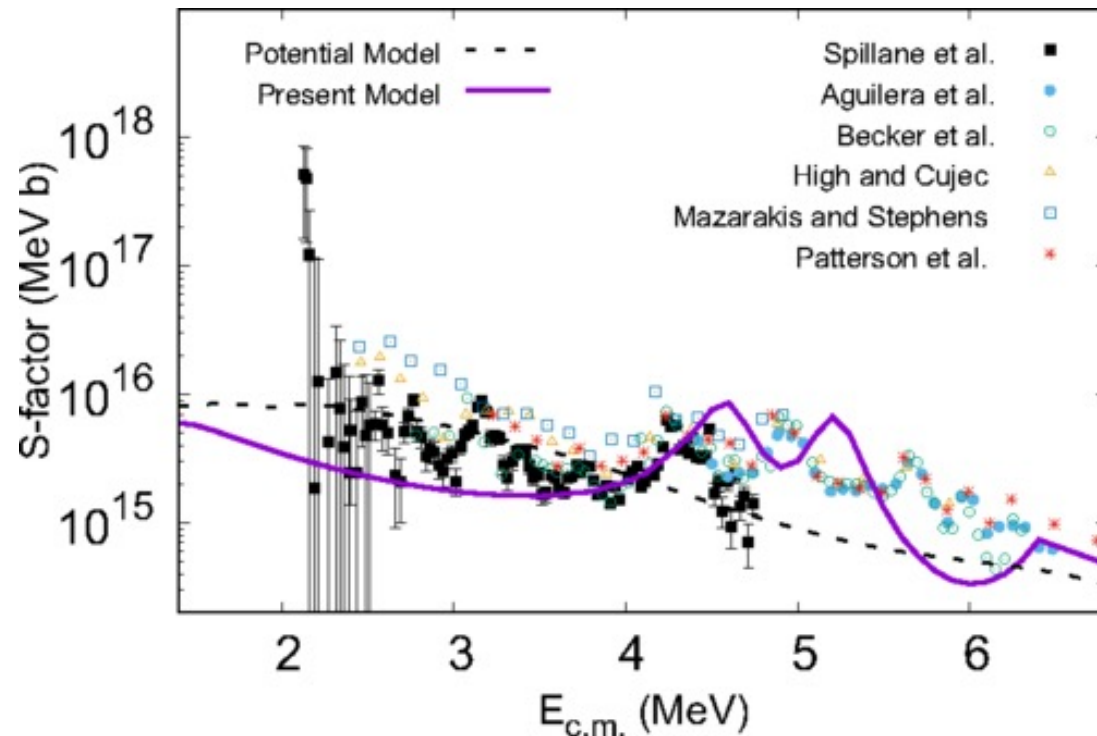
## □ Time-dependent approach

- ✓ A. Diaz-Toress and M. Wiesher, PRC97, 055802 (2018).  
Time-dependent wave packet approach



$$i\hbar \frac{\partial}{\partial t} \psi(r, \theta_1, k_1, \theta_2, k_2; t) = H \psi(r, \theta_1, k_1, \theta_2, k_2; t)$$

$$\psi(r, \theta_1, k_1, \theta_2, k_2; t = 0) = \chi_0(r) \psi_0(\theta_1, k_1, \theta_2, k_2)$$



Gaussian rotor  
wave packet

## Previous theoretical calculations for $^{12}\text{C}+^{12}\text{C}$

### □ TDHF-based approaches

✓ P.-H. Heenen, PL99B, 298 (1981).

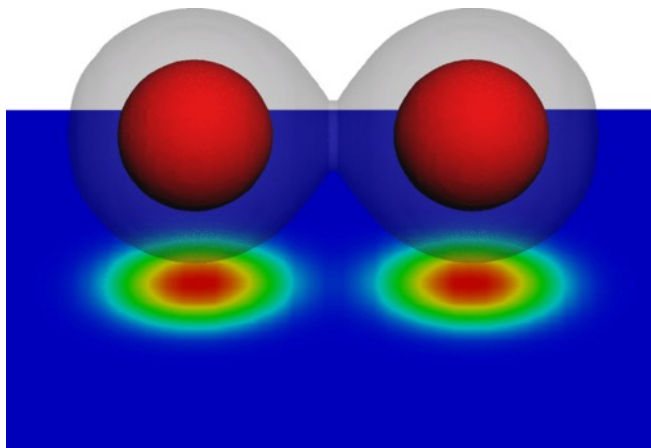
#### Adiabatic TDHF

✓ K. Godbey, C. Simenel, and A.S. Umar, PRC100, 024619 (2019).

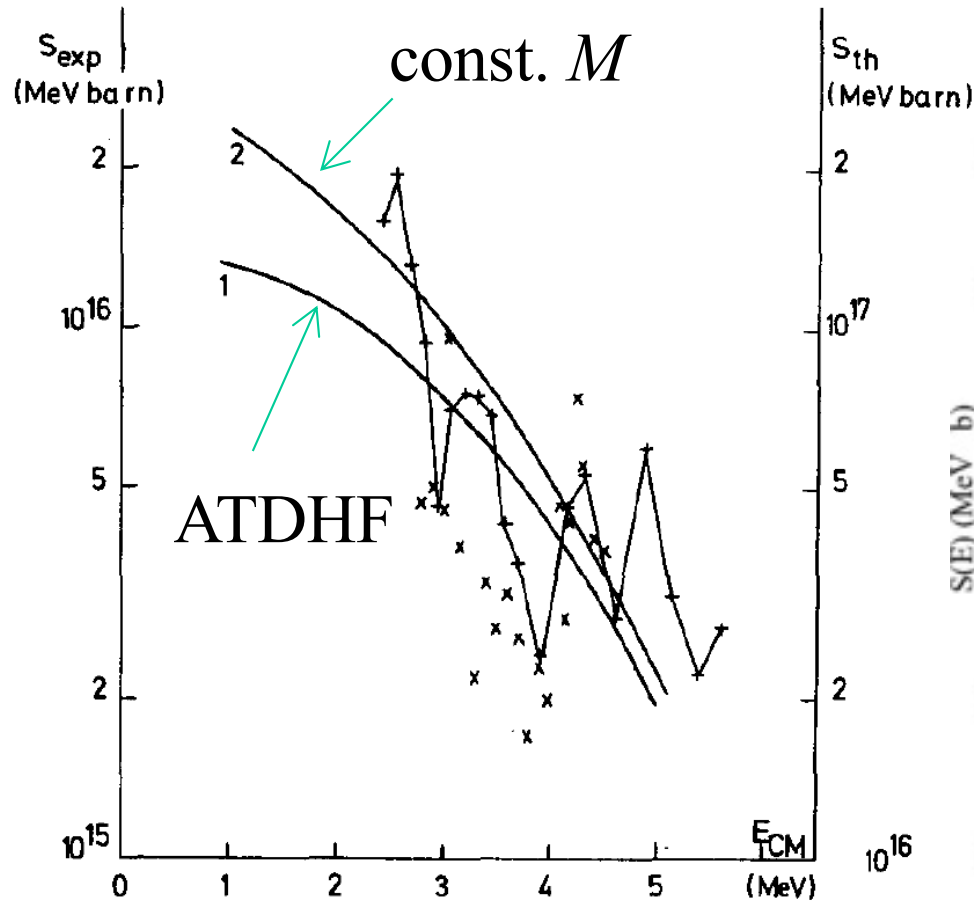
#### Density-Constrained Hartree-Fock

collective Hamiltonian from TDHF

$$H_{\text{coll}} = \frac{1}{2}M(R)\dot{R}^2 + V(R) \rightarrow P_{\text{fus}}(E)$$

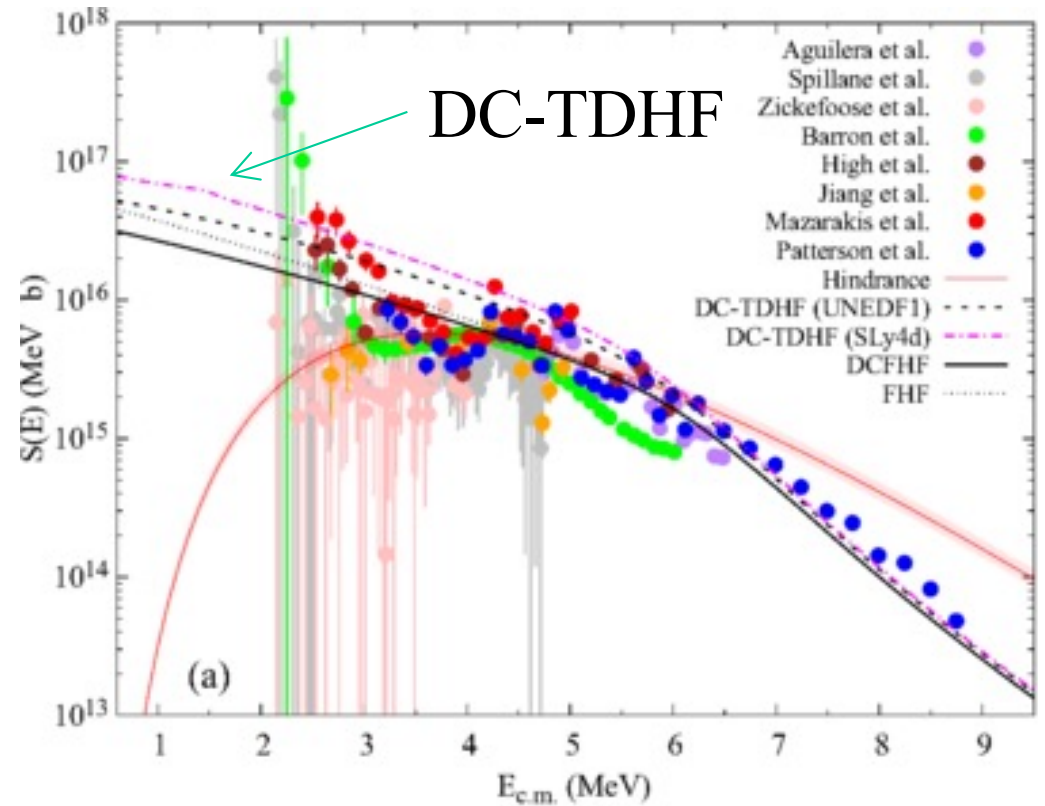


# ATDHF



P.-H. Heenen,  
 PL99B, 298 (1981)

# DC-TDHF



K. Godbey, C. Simenel, and  
 A.S. Umar,  
 PRC100, 024619 (2019).

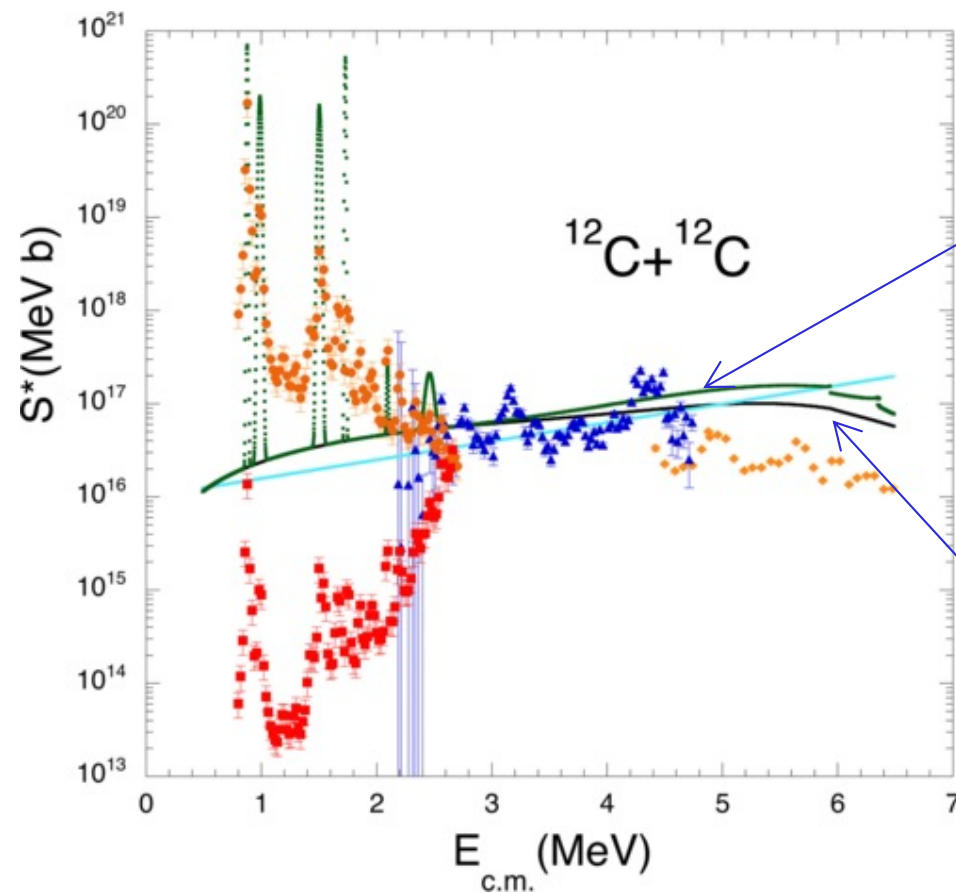
## Previous theoretical calculations for $^{12}\text{C}+^{12}\text{C}$

### □ TDHF-based approaches

- ✓ A. Bonasera and J.B. Natowitz, PRC102, 061602(R) (2020).

### Imaginary-time evolution

$$i\hbar \frac{\partial}{\partial t} \psi_i(\mathbf{r}, t) = h[\rho] \psi_i(\mathbf{r}, t) \rightarrow -\hbar \frac{\partial}{\partial \tau} \psi_i(\mathbf{r}, \tau) = h[\rho] \psi_i(\mathbf{r}, \tau)$$



im.-time TDHF  
+resonances  
(justification?)

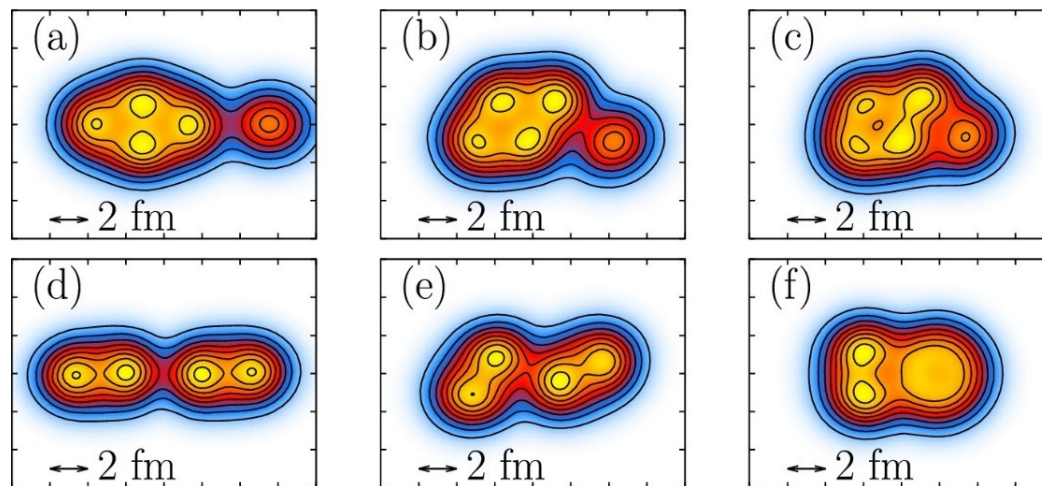
im.-time TDHF

# Previous theoretical calculations for $^{12}\text{C}+^{12}\text{C}$

## □ Anti-symmetrized Molecular Dynamics (AMD) approach

Y. Taniguchi and M. Kimura, PLB823, 136790 (2021).

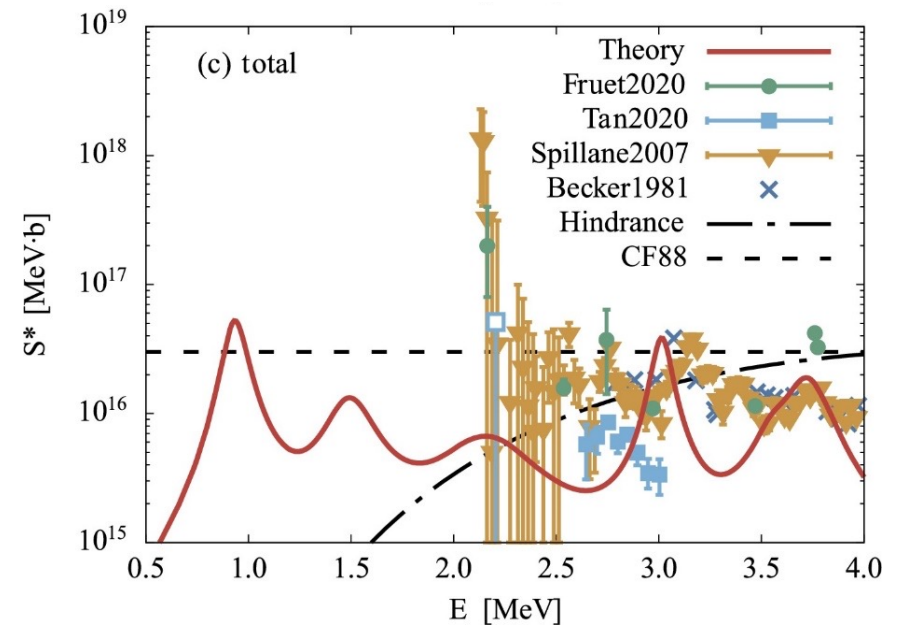
### AMD + R-matrix



$$\Psi_M^{J\pi} = \sum_{iK} c_{iK} P_{MK}^J \Phi_i^\pi$$

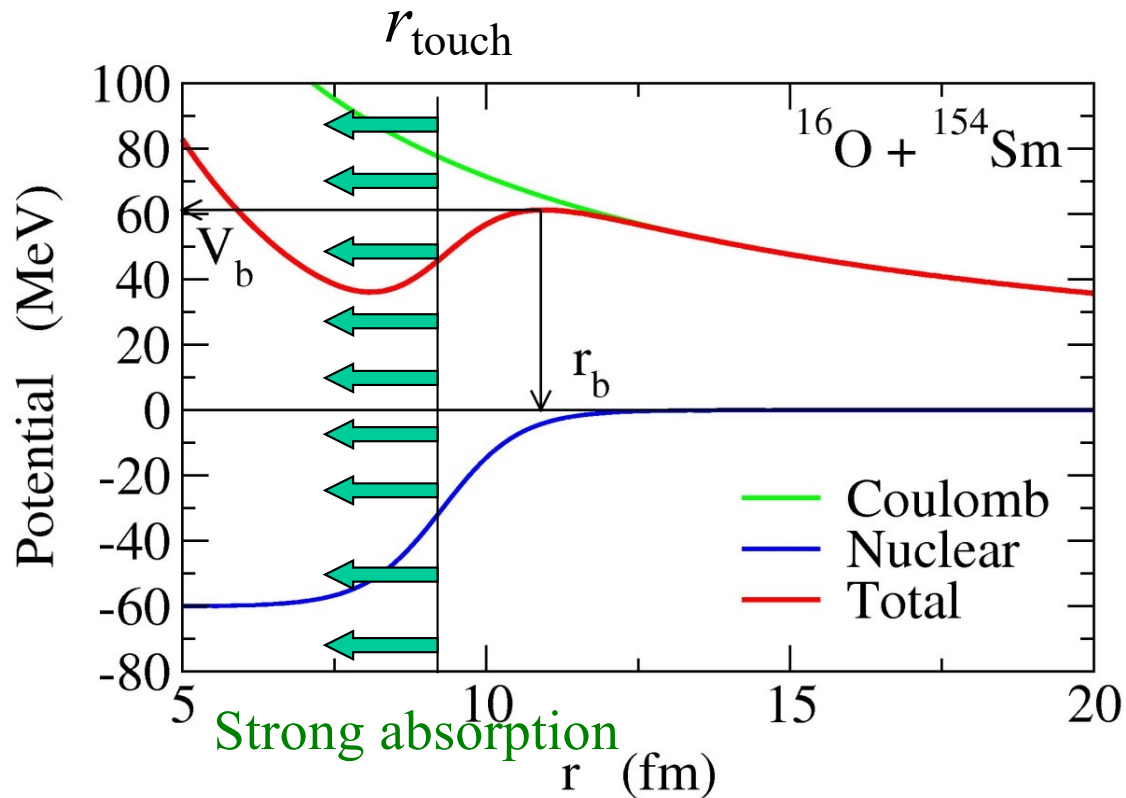
→ R-matrix to match the outer  
Coulomb wave functions

→  $\sigma_{\text{fus}}(E)$

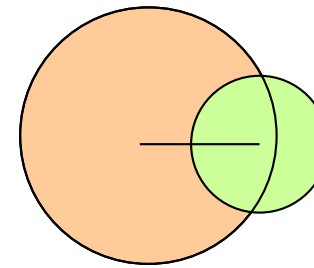




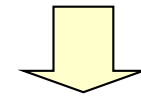
# Strong absorption model



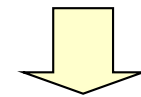
the region of large overlap



• High level density (CN)



The relative energy is quickly lost and converted to internal energies



Formation of hot CN (fusion reaction)

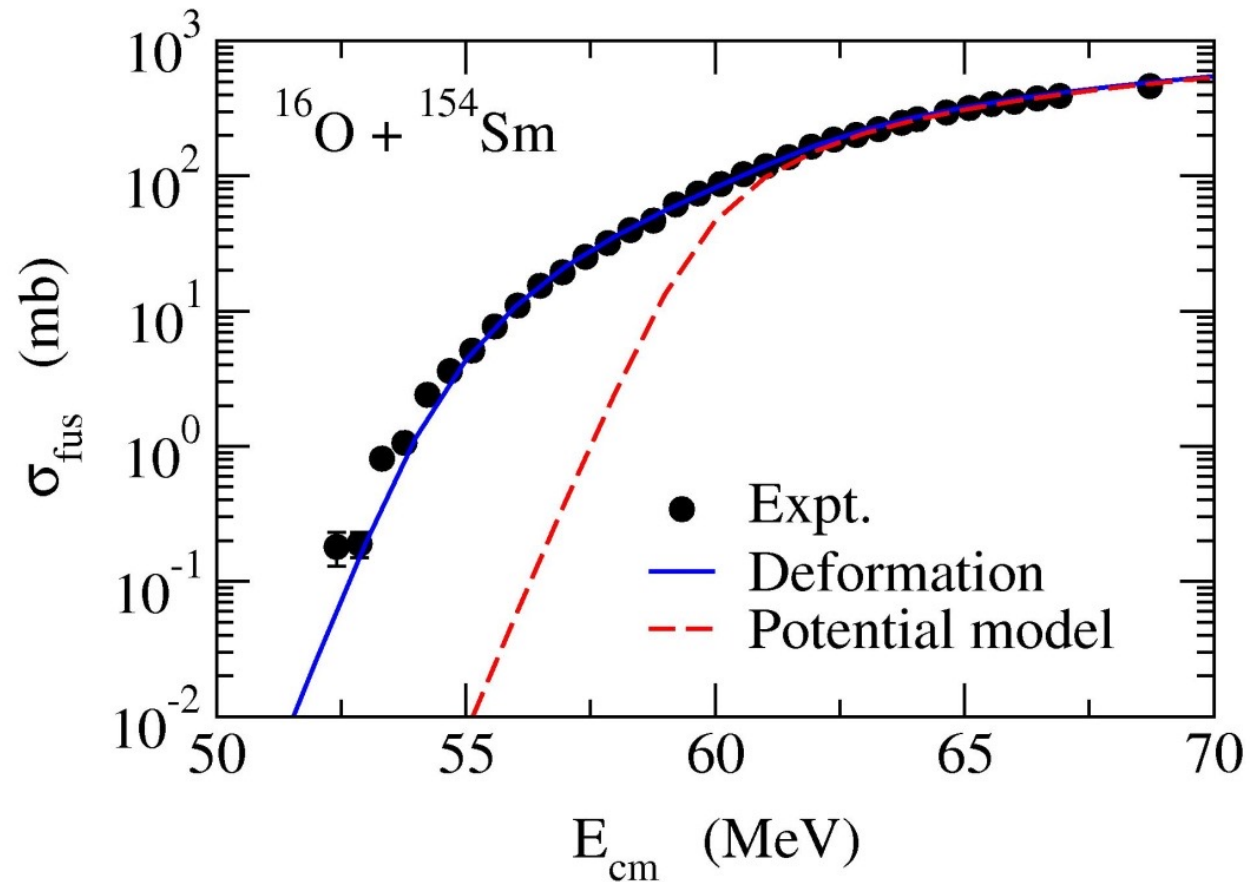
This can be simulated either with

✓ a large imaginary potential at the potential pocket

or with

✓ the incoming wave boundary condition

The assumption of strong absorption works good for heavy systems:

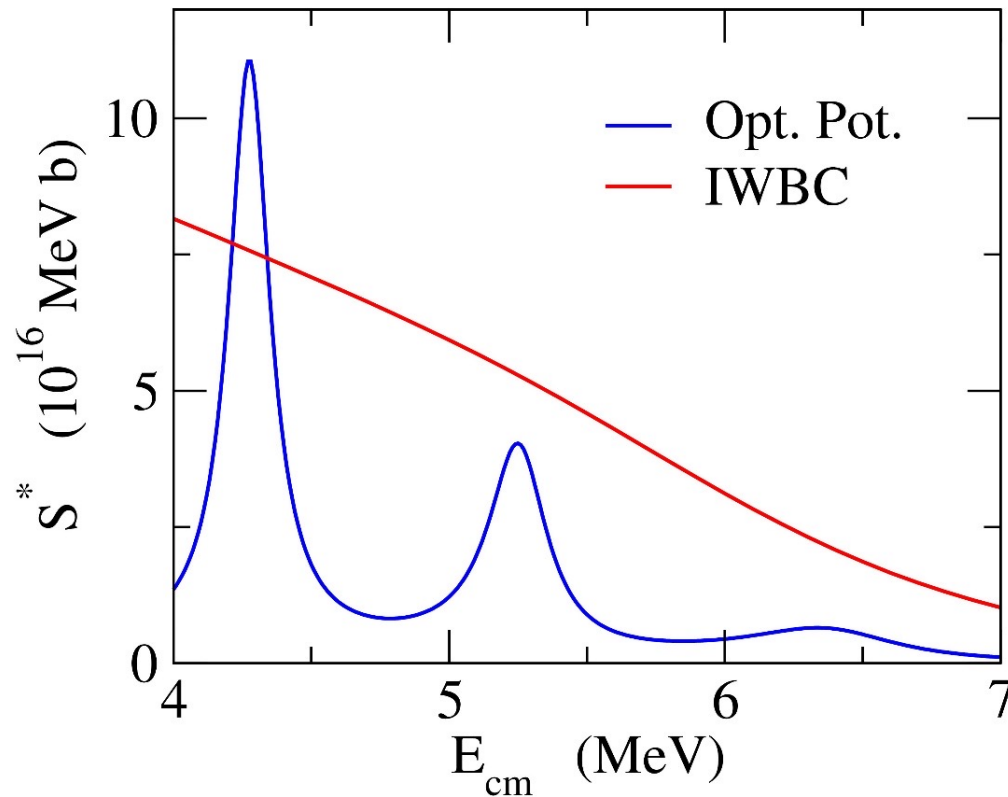


K. Hagino and N. Takigawa,  
Prog. Theo. Phys.128 ('12)1061.

....but this may not be the case for light systems such as  $^{12}\text{C}+^{12}\text{C}$



## single-channel calculation



$V_0=20$  MeV,  $r_0=1.35$  fm,  $a=0.5$  fm  
 $W=0.2$  MeV,  $r_w=1.05$  fm,  $a_w=0.2$  fm

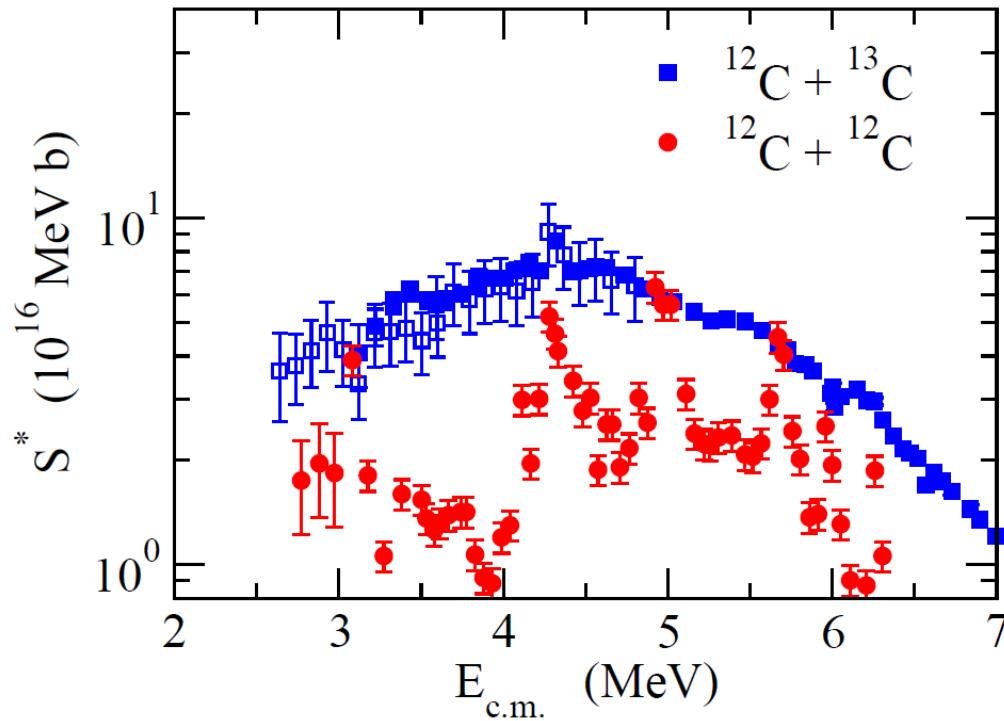
IWBC:

- ✓ no resonance structure
- ✓ “the upper limit” (but not really) of S-factors

## Effects of isolated CN resonances

C.L. Jiang et al.,  
PRL110, 072701 (2013).

cross section deficit

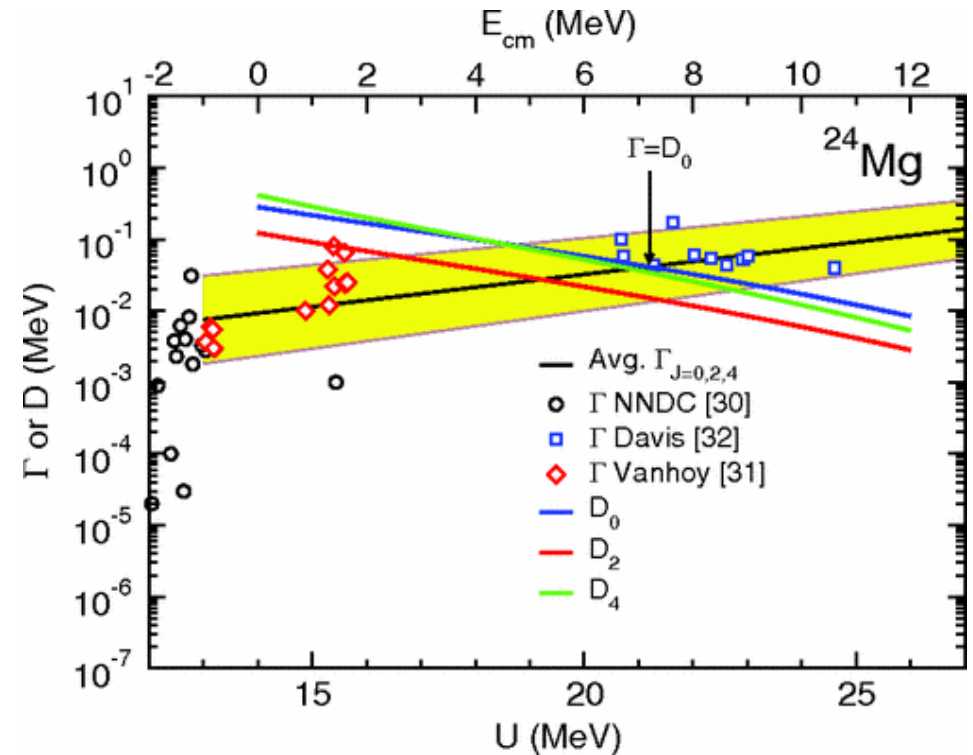


$$E_{c.m.} < 7 \text{ MeV}$$

$\Gamma < D \rightarrow$  fusion inhibition

$$E_{c.m.} > 7 \text{ MeV}$$

$$\Gamma > D \rightarrow P_J \sim 1$$



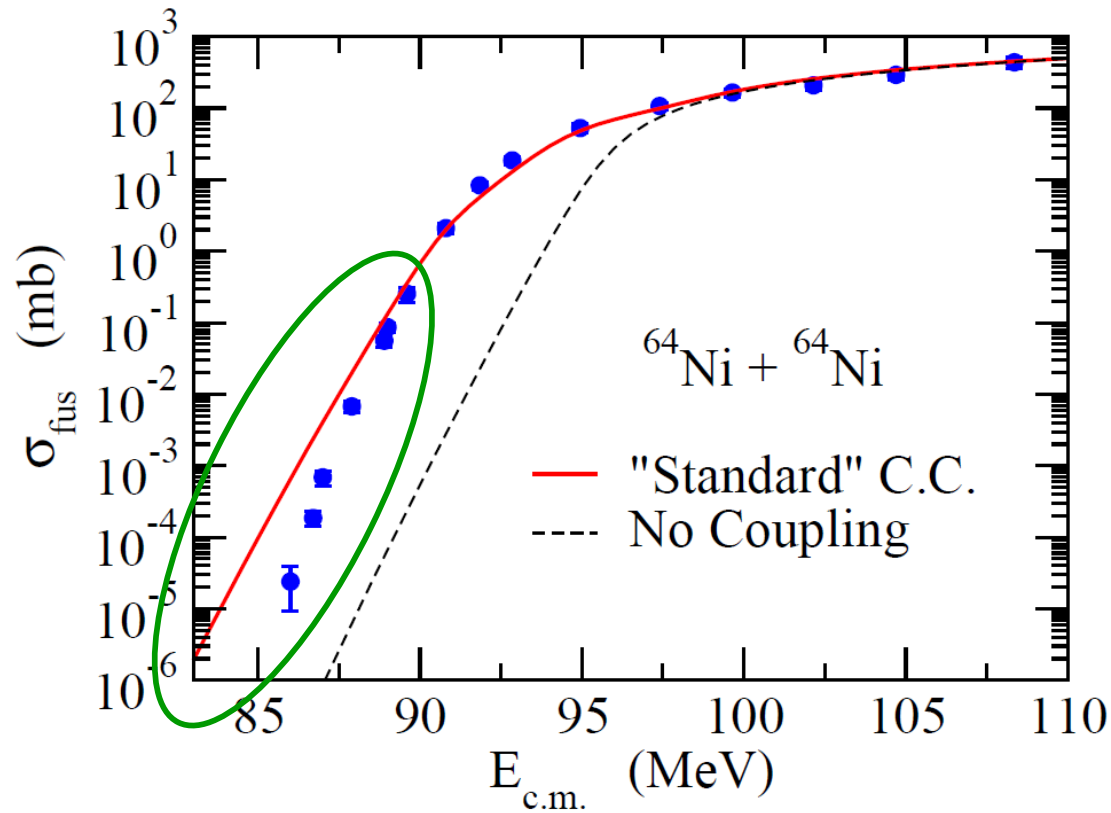
Moldauer factor

$$P_J = 1 - \exp(-2\pi\Gamma_J/D_J)$$

$$\rightarrow \sigma = \sum_J \sigma_{CC}^J P_J$$

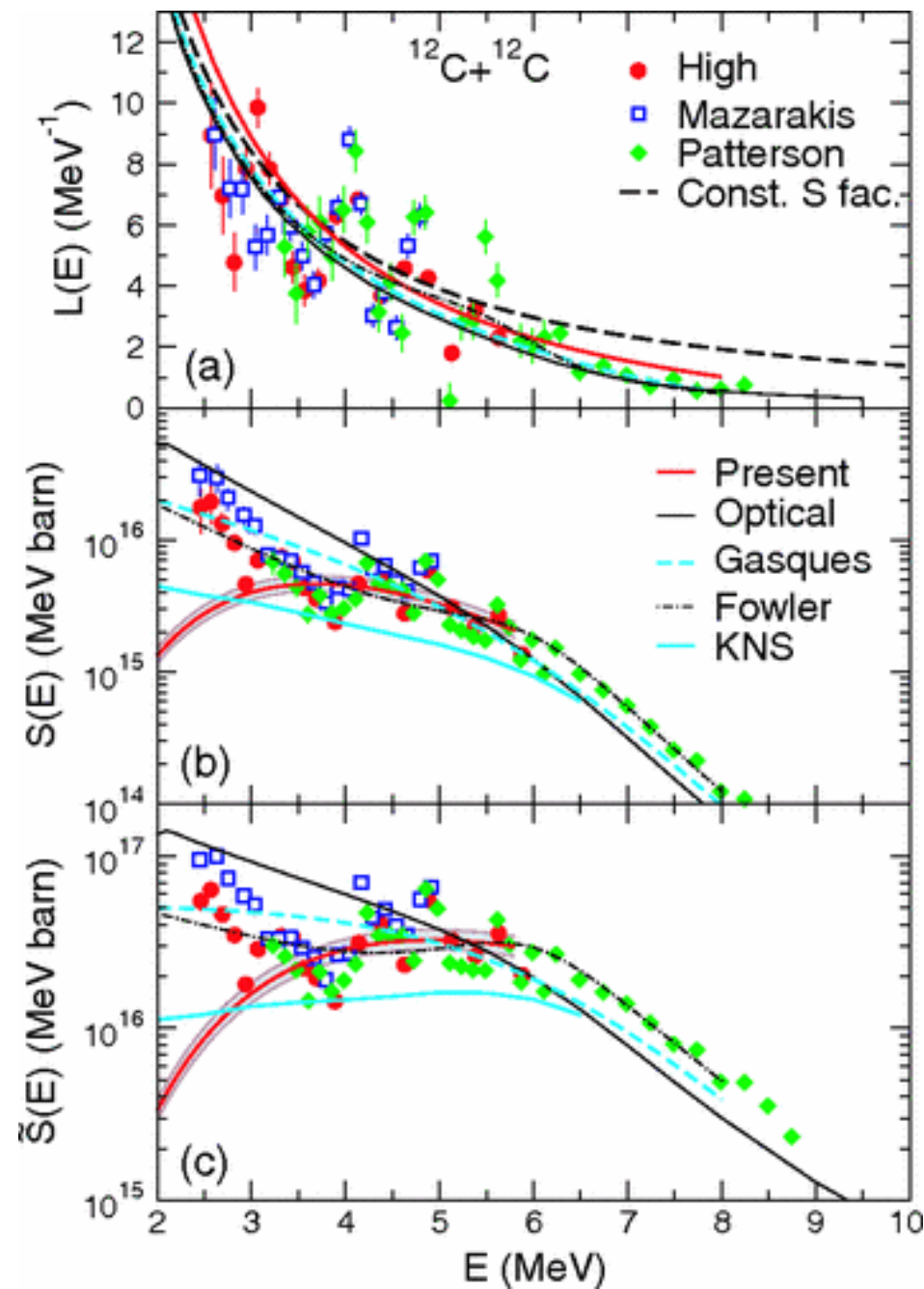
# Hindrance

Deep sub-barrier hindrance in medium heavy systems



C.L. Jiang et al., PRL89('02)052701;  
PRL93('04)012701

Is there a large hindrance in  $^{12}\text{C}+^{12}\text{C}$  fusion?



C.L. Jiang et al.,  
PRC75, 015803 (2007)

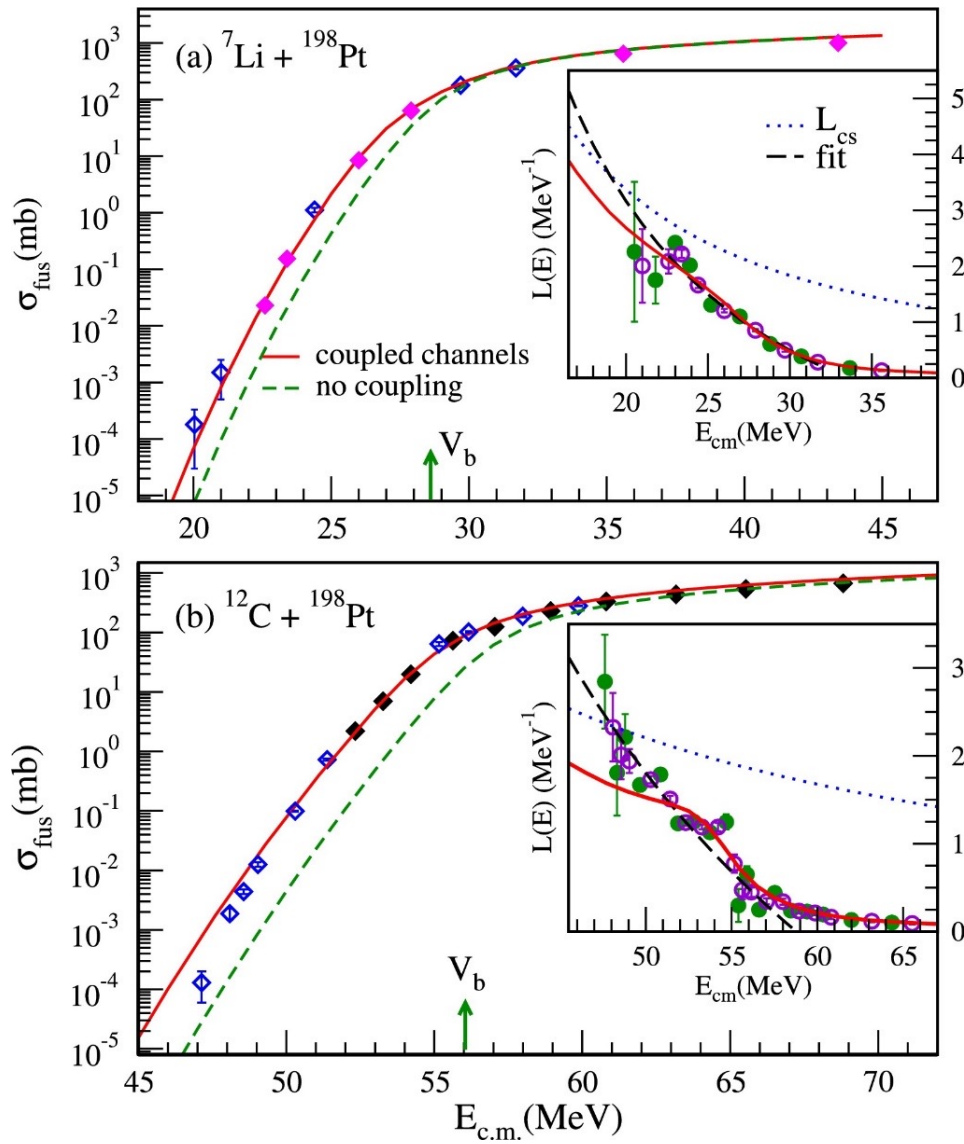
$$L(E) = \frac{d}{dE} \ln(E\sigma)$$

fit with:

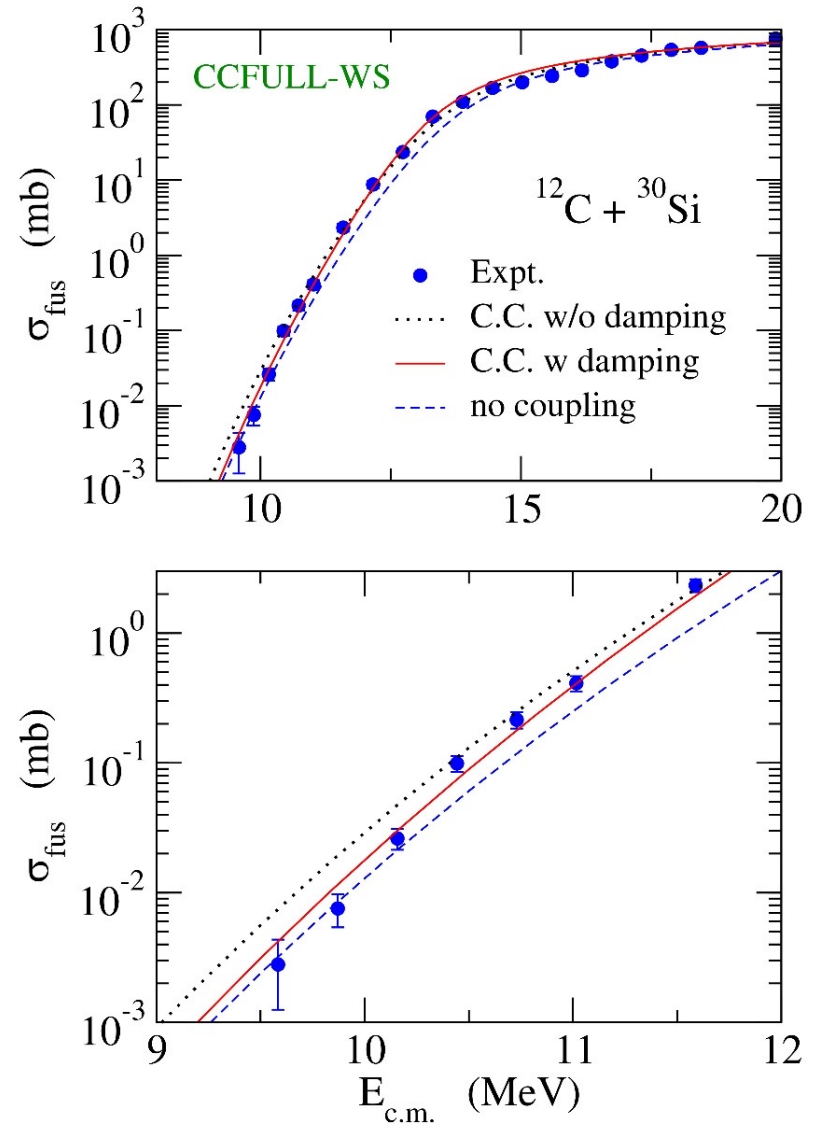
$$L(E) = A_0 + B_0/E^n$$

→ large hindrance

no (or only tiny) hindrance has been observed in light systems



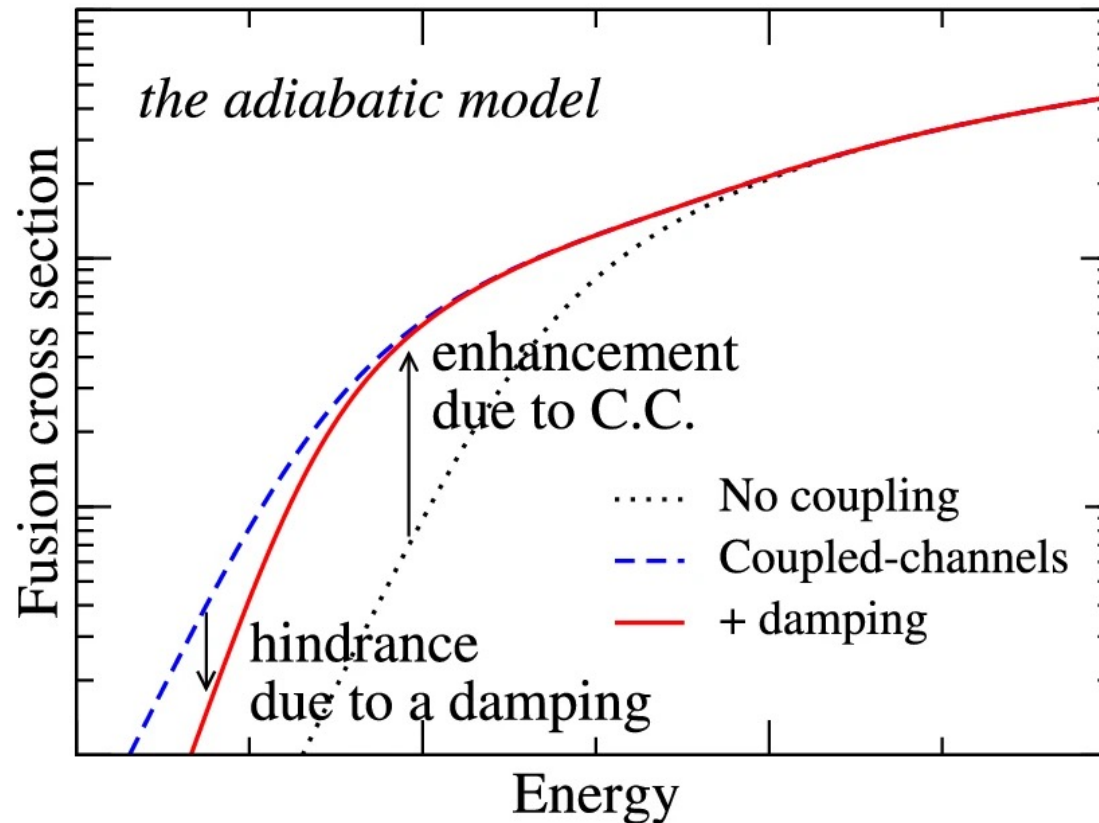
A. Shrivastava et al.,  
PLB755, 332 (2016).



G. Montagnoli et al.,  
PRC97, 024610 (2018).

## adiabatic model for fusion hindrance:

T. Ichikawa, K.H., and A. Iwamoto, PRL103('09)202701



C.L. Jiang et al., EPJA57, 235 (2021)

**light systems:** small enhancement of fusion cross sections  
→ small hindrance

# Summary

Open questions:

- \* How can one connect the resonance peaks to CN states?
  - can one have a model which reproduces all the resonance peaks?
  - how can one connect CN states to the reaction dynamics?
  - can one construct a toy model to start with?
- \* Can one describe  $^{12}\text{C}+^{12}\text{C}$  and  $^{12}\text{C}+^{13}\text{C}$  simultaneously?
  - Does  $^{12}\text{C}+^{13}\text{C}$  provide the upper limit of  $^{12}\text{C}+^{12}\text{C}$  cross sections?
- \* How much fusion hindrance can one expect for  $^{12}\text{C}+^{12}\text{C}$ ?
- \* Electron screening effects?

