Micro-physics of stellar interiors

Werner Däppen

Department of Physics and Astronomy
University of Southern California
Los Angeles

How does micro-physics enter?

Forward problem

■ Nobody would do it in this way (!), but ...

$$\begin{bmatrix} \frac{\partial \mathbf{v}}{\partial t} + \mathbf{v} \cdot \nabla \mathbf{v} \end{bmatrix} = -\frac{1}{\rho} \nabla p - \nabla \phi$$
$$\begin{bmatrix} \frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \mathbf{v}) \end{bmatrix} = 0$$
$$\begin{bmatrix} \frac{\partial s}{\partial t} + \mathbf{v} \cdot \nabla s \end{bmatrix} = -\frac{1}{T} \epsilon_{\mathsf{nucl}} - \frac{1}{\rho T} \; \mathsf{div} \mathbf{F}$$
$$\Delta \phi = 4\pi G \rho$$

... let's illustrate stellar evolution and pulsation by using the same equations, both for the equilibrium and the oscillation problem

First: the equilibrium solution

- Put all red stuff to 0!
- Obtain the usual equilibrium equations (for simplicity assume (i) spherical symmetry; (ii) no convection)

$$\frac{dp}{dr} = -\frac{GM_r\rho}{r^2}$$

$$L_r = 4\pi r^2 F$$

$$\frac{dL}{dr} = 4\pi r^2 \rho \epsilon_{\text{nuc}}$$

$$\frac{dT}{dr} = -\frac{3\kappa\rho L}{64\pi\sigma r^2 T^3}$$

Material properties are mandatory!

- So far all equations the same for all stars
- □ Richness of variety of stars only enters with the constituent equations

$$\begin{split} & \rho = \rho(T, p, \mathbf{X}) \\ & \kappa = \kappa(T, \rho, \mathbf{X}) \\ & \epsilon = \epsilon_{\mathsf{nuc}}(T, \rho, \mathbf{X}) \\ & s = s(T, \rho, \mathbf{X}) \\ & \mathbf{F} = \mathbf{K}(\nabla T; T, \rho, \mathbf{X}) \\ & (\mathsf{Diffusive radiation}: \ F \ = -\frac{16\sigma T^4}{3\kappa\rho} \nabla T \) \end{split}$$

Microphysics is not only for the variety of stars:

Material properties are also essential for stellar (parametric) evolution!

This is a seismology conference, therefore go to next step:

Second: oscillation equations

Once equilibrium model found, solve red part by putting black part = 0

$$\begin{bmatrix} \frac{\partial \mathbf{v}}{\partial t} + \mathbf{v} \cdot \nabla \mathbf{v} \end{bmatrix} = -\frac{1}{\rho} \nabla p - \nabla \phi = 0$$
$$\begin{bmatrix} \frac{\partial \rho}{\partial t} + \operatorname{div}(\rho \mathbf{v}) \end{bmatrix} = 0$$
$$\begin{bmatrix} \frac{\partial s}{\partial t} + \mathbf{v} \cdot \nabla s \end{bmatrix} = -\frac{1}{T} \epsilon_{nuc} - \frac{1}{\rho T} \operatorname{div} \mathbf{F} = 0$$

 \square If appropriate, use linearization, spherical harmonics, adiabatic $\Rightarrow \nu_{n\ell}$

Result: model frequencies

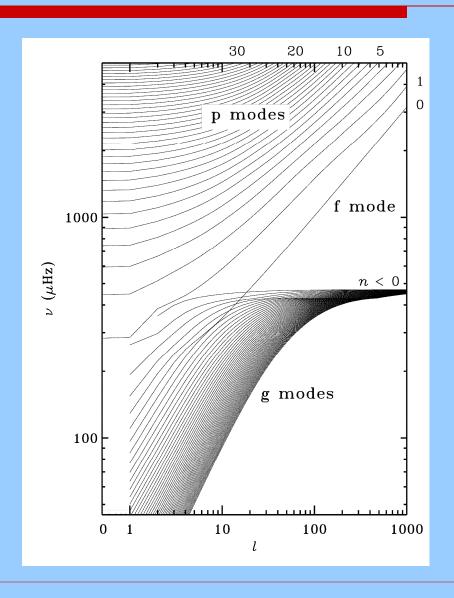


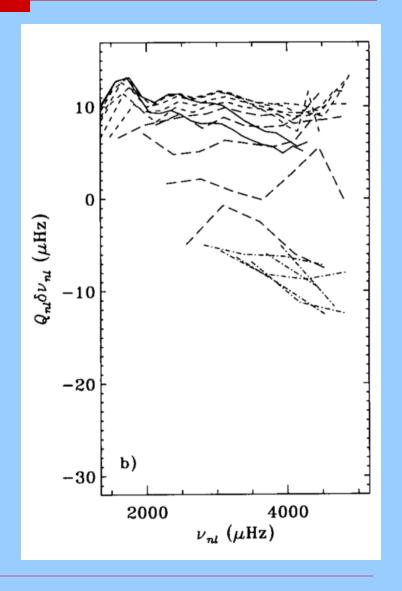
Figure:

J. Christensen-Dalsgaard

Allow "O-C Diagrams"

- \square O = "Observation"
- \square C = "Computation"

(Q_{nl} is a scale factor. Details see: Christensen-Dalsgaard & Däppen 1992, A&A Rev. **4**, 267)



Tools to test physics

Since the equation of state is the most basic ingredient, let's go into more detail!

In the Sun and solar-like stars: fortunate situation

- In the convection zone, the stratification is (largely) adiabatic, its structure is mainly determined by thermodynamics.
- Little "contamination" from opacity
- Helioseismology can probe locally (to some degree also asteroseismology, e.g., in abundance determination)

Elementary stellar thermodynamics

In stellar physics, before 1975, normal (non-degenerate) stars were successfully modeled by

$$pV = (\Sigma_i N_i)kT$$

lacktriangle Provided: N_i from Saha equation:

Good to 90% accuracy!

In early helioseismology

- From 1975-1985, more refined equations of state, mainly
- Detailed chemical composition
- □ Fermi-Dirac electrons
- Debye-Hückel screening

good to 95-99% accuracy!

The **dominant** non-ideal effect turned out to be...**DH!**

Screened Coulomb potential

$$V_D = \frac{Q_1 Q_2}{r} \exp^{-r/r_D}$$

(=the [static] Debye-Hückel approximation)

Impact: two similar solar models...

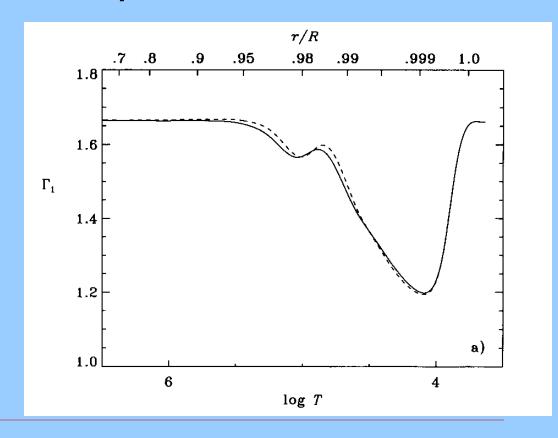
□ Identical models, except for their equations of state. One is with Debye-Hückel screening, one without. Their adiabatic exponents are:

Dashed: with screening

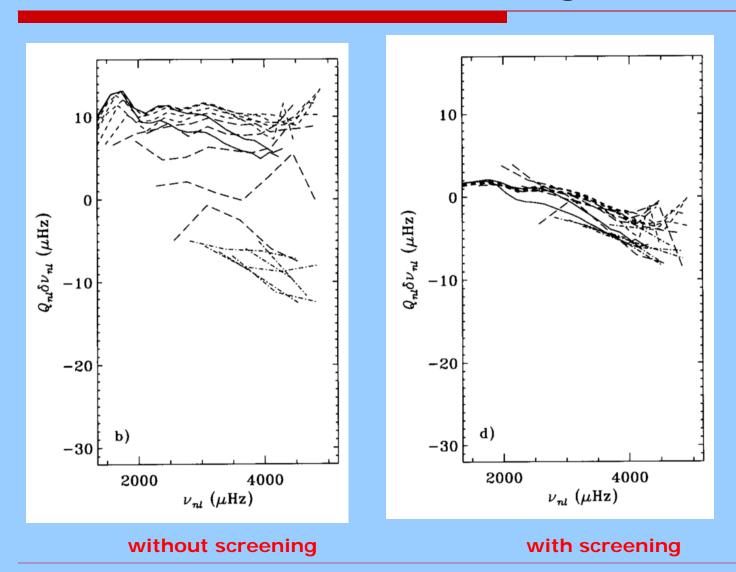
Solid: without screening

From: Christensen-Dalsgaard & Däppen

1992, A&A Rev. 4, 267



...and their O-C diagrams



from: Christensen-Dalsgaard & Däppen 1992, A&A Rev. 4, 267

To my chagrin, the physicists were not overwhelmed!

A tangent Seismic abundance determination

Acoustic modes (largely) governed by the adiabatic sound speed

$$c_{\rm ad}^2 = \gamma_1 \frac{p}{\rho} \propto \frac{T}{\mu}$$

$$\gamma_1 = (\frac{\partial p}{\partial \rho})_s \text{ (often denoted } \frac{\Gamma_1 \text{ to Douglas Gough's chagrin!)}}$$

Intuitive reasoning

$$\gamma_1pprox\gamma=rac{c_p}{c_V} \ c_p-c_Vpprox1$$
 (1st Law; molar units)

 c_p , c_V each grow in ionization zones (analogous to latent heat!)

$$\gamma = \frac{c_p}{c_V}$$
 drops in ionization zones: down from 5/3 to about 1.2 for a hydrogen plasma

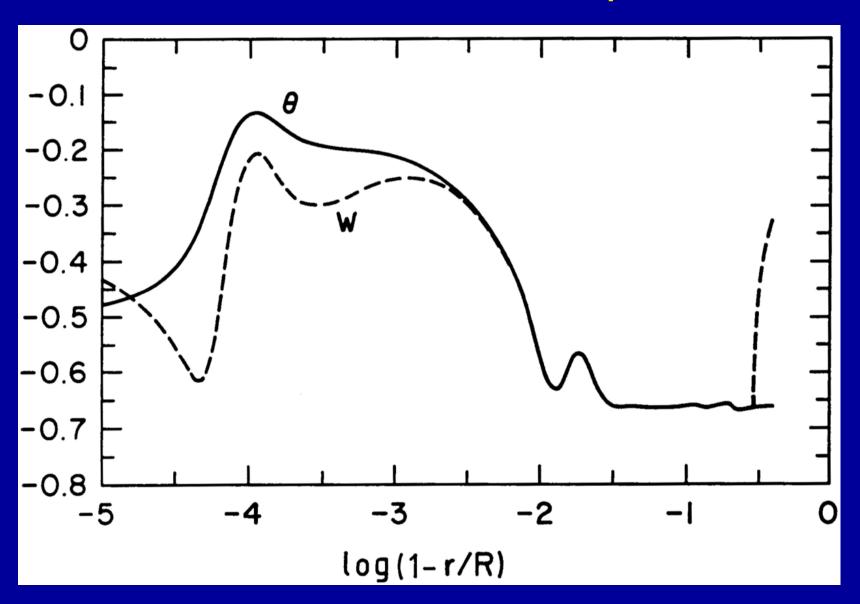
Supertool for abundances

Douglas Gough (Catania 1983) applies adiabatic constraint to the hydrostatic equation:

$$W = \frac{dc^2}{dr} \frac{1}{g(r)} = \frac{1 - \gamma_1 - \gamma_{1,\rho}}{1 - \gamma_{1,c^2}} = \Theta$$

What would that look like inside the Sun?

The helium bump!



Result of the helium hump method (and similar techniques, e.g., by Dziembowski, Thompson, Vorontsov, Antia, Basu...)

$$Y = 0.24...0.25$$

inside the solar convection zone

Whatever the value of Y thus obtained, the result is entangled with the uncertainty in the equation of state!

Practical tools for modelers I

Phenomenological options

Capparage of the control of the c

On the plus side

Useful computational tool, remarkably accurate Broad range of applicability Even includes relativistic electrons

On the minus side

Misses dominant non-ideal term (later fixed, see next) Ad-hoc pressure ionization term Can lead to artificial phase transitions

There are improved descendants of EFF, great tools for modelers!

• ... CEFF

... by J. Christensen-Dalsgaard and WD (1991), who added a Coulomb (C) term, following progress in helioseismology

... SIREFF

... by Swenson, Irwin and Rogers (1996) who added excited-states terms, later extended into:

... FreeEOS

... by Irwin (excellent tool, freely downloadable from [http://freeeos.sourceforge.net/], newest version: 2008)

WARNING:

BORING STUFF!

(even more of it in the Appendix, if interested...)

Two main approaches: introduction

- Free-energy minimization chemical picture intuitive, but highly practical
- Grand-canonical expansions Physical picture systematic method for non-ideal corrections

Chemical picture

- Treat compounds as fundamental entities
- \square Reactions (H \leftrightarrow H⁺ + e⁻, etc.)
- \square Constraints $(N_{\mathsf{H}} + N_{\mathsf{p}} = \mathsf{const.}\ , \ \mathsf{etc.})$
- \square Minimize $F(T, V, N_{\mathsf{H}}, N_{\mathsf{p}}, N_{\mathsf{e}^{-}}, ...)$
 - !!!subject to constraints!!!
- □ In practice, cook a free energy (intuition!)

$$F_{\text{tot}} = F_{\text{nuc}} + F_{\text{e}} + F_{\text{interactions}} + \dots$$

 \square Consistent (formally!) $p = -(\frac{\partial F}{\partial V})_T$, etc.

Example: MHD

- Fairly conventional realization (chemical picture)
- Key ingredient: occupation probabilities

Hummer, D.G. & Mihalas, D.M. 1988, *ApJ* **331**, 794; Mihalas, D.M., Däppen, W. & Hummer, D.G. 1988, *ApJ* **331**, 815 Däppen, W., Mihalas, D.M., Hummer, D.G. & Mihalas, B.W. 1988, *ApJ* **332**, 261

$$Z_{jk}^{\mathrm{int}} = \sum_{i} w_{ijk} \ g_{ijk} \mathrm{exp} \left(-\beta E_{ijk} \right)$$

$$\begin{split} &(w_{ijk})_{\text{neutral}} = \exp\left[-(4\pi/3V)\sum_{j',k'}N_{j'k'}(r_{ijk} + r_{1j'k'})^3\right] \\ &(w_{ijk})_{\text{charged}} = \exp\left\{-\left(\frac{4\pi}{3V}\right)16\left[\frac{(Z_{jk}+1)^{1/2}e^2}{K_{ijk}^{1/2}\chi_{ijk}}\right]^3\sum_{j',k'}N_{j'k'}Z_{j'k'}^{3/2}\right\} \end{split}$$

Physical picture

- Only electrons and nuclei are fundamental
- No reactions
- Quantum mechanics and statistical mechanics dealt with simultaneously
- Nothing to minimize
- Consistent (not just formally!)

Example: OPAL/ACTEX

 First successful stellar modeling with an equation of state in the physical picture (LLNL)

> Rogers, F.J. 1986, *ApJ* **310**, 723; Rogers, F.J., Swenson, F.J. & Iglesias, C.A. 1996, *ApJ* **456**, 902 Rogers, F.J. & Nayfonov, A. 2002, *ApJ* **576**, 1064

Key points: systematic expansions
 (z = activity)

$$\frac{p}{k_B T} = z + z^2 b_2 + z^3 b_3 + \dots \; ; \quad \rho = \frac{z}{k_B T} (\frac{\partial p}{\partial z})$$

Planck-Larkin Partition Function

$$PLPF = \sum_{nl} (2l+1) \left[exp(-\frac{E_{nl}}{kT}) - 1 + \frac{E_{nl}}{kT} \right]$$

Domain of validity of OPAL/ACTEX equation of state for stellar models

> 0.1 solar masses

MHD vs. OPAL in the Sun

c^2 Inversions (numerical; Sun-Model)

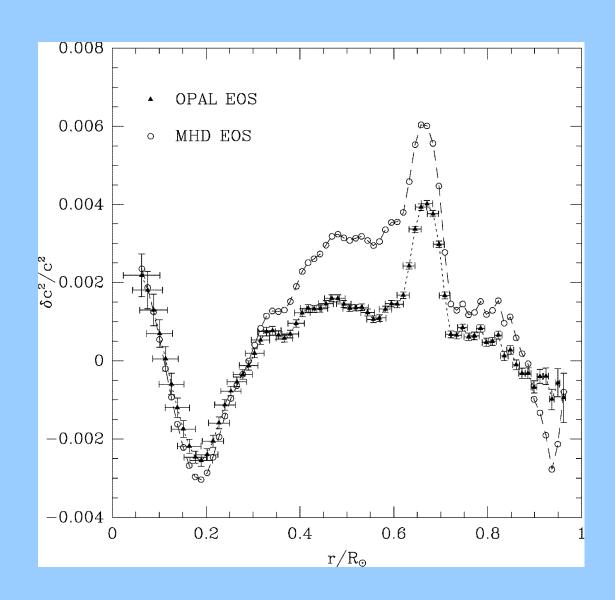


Figure from: S. **Basu**

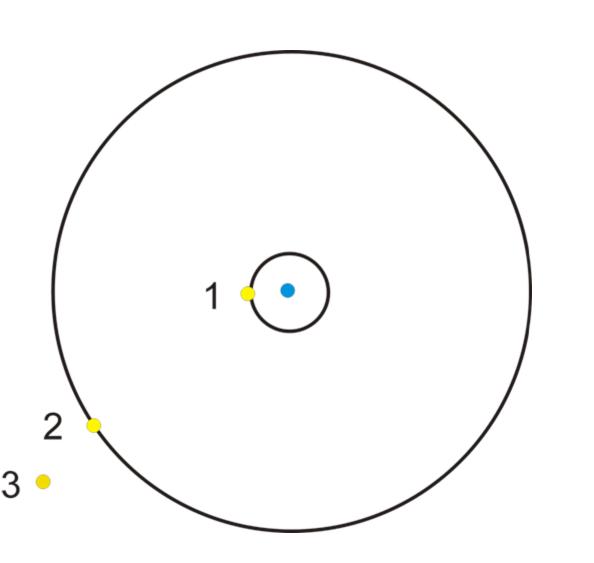
OPAL fares better than MHD...

□ Why? Likely answer:

- ☐ There is no PLPF in MHD
- □ There are no scattering states in MHD

□ Open question: is it fundamentally impossible to find PLPF entirely from within the chemical picture?

More precisely: consider 3 states of a H atom...



1: ground state

2: weakly bound

3: continuum

force law for electrons (C=Coulomb, F=free)

	CEFF	MHD	OPAL
1: ground state 2: weakly bound 3: continuum	C F F	C C F	C C

Treatment of excited states:

OPAL good, CEFF at least consistent, MHD inconsistent (as are all chemical picture formalisms with excited states but without PLPF)

FYI: more of it in an appendix!

Further possibilities towards very high precision...

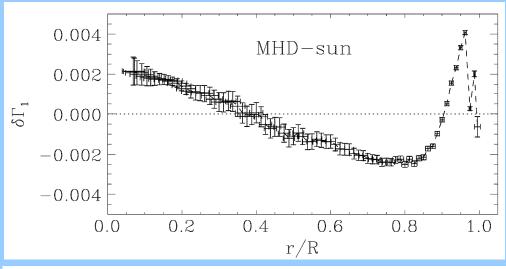
To illustrate: a small effect - relativistic electrons in the Sun

Relativistic corrections are expected to be small, central temperature

$$kT \approx 1 \text{ keV} \ll 511 \text{ keV}$$

And yet: the effect can be observed!! (Elliot & Kosovichev 1998, ApJ, 500 L199)

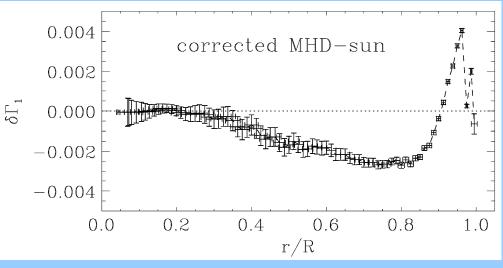
Models with and without relativistic electrons



Figures from:

Elliot, J.R. & Kosovichev, A.G.

1998, *ApJ*, **500** L199

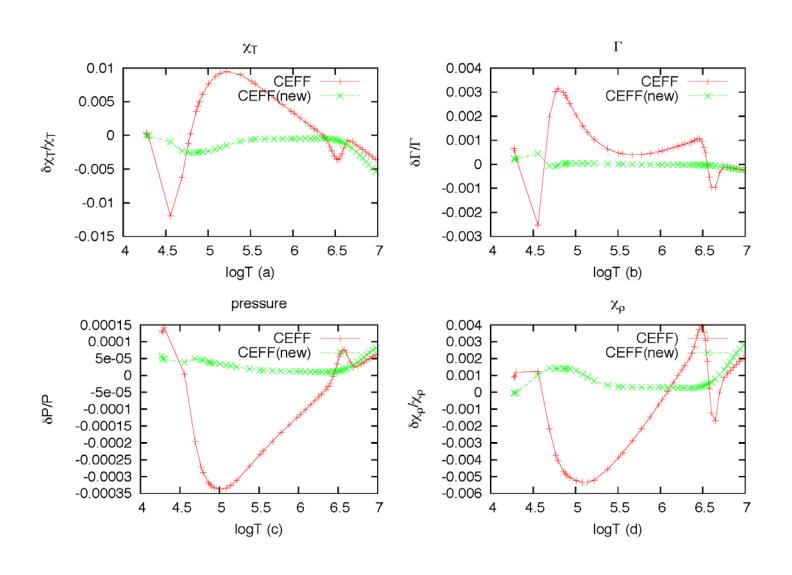


Practical tools for modelers II

Tool for the community

- OPAL emulator as an in-line formalism, starting out from CEFF (Hsiao-Hsuan Lin)
- Higher precision than FreeEOS but only for Sun, while FreeEOS has larger stellar domain
- Work in progress, in several steps, current state on following slide

Results (Hsiao-Hsuan Lin, 2011)



A further option is SAHA-S

SAHA-S is a serious alternative to OPAL. It is based on a systematic development, and it contains a modified Planck-Larkin partition function.

See:

Starostin, A.N. & Roerich, V.C., Phys. A: Math. Gen. 39(2006)4431

Truly Exact Options

- □ There are exact density expansions, for instance the Feynman-Kac [FK] path-integral computations [1] or Green-function calculations [2].
- ☐ The coefficients in these expansions are indeed exact!
- Problem is the domain of applicability!
- Why? See Appendix, if you insist!
- [1] Alastuey, A. & Perez, A. 1992, Europhys.Lett., 20, 19
- [2] Kraeft, W.-D., Kremp, D., Ebeling, W., Röpke, G., 1986, Quantum Statistics of Charged Particle Systems (Plenum)

$$\beta P = \sum_{\alpha} \rho_{\alpha} - \frac{\kappa_{D}^{3}}{24\pi}$$

$$+ \frac{\pi}{6} (\ln 2 - 1) \sum_{\alpha,\beta} \beta^{3} e_{\alpha}^{3} e_{\beta}^{3} \rho_{\alpha} \rho_{\beta}$$

$$- \frac{\pi}{\sqrt{2}} \sum_{\alpha,\beta} \rho_{\alpha} \rho_{\beta} \lambda_{\alpha\beta}^{3} Q(x_{\alpha\beta}) - \frac{\pi}{3} \beta^{3} \sum_{\alpha,\beta} \rho_{\alpha} \rho_{\beta} e_{\alpha}^{3} e_{\beta}^{3} \ln (\kappa_{D} \lambda_{\alpha\beta})$$

$$+ \frac{\pi}{\sqrt{2}} \sum_{\alpha} \frac{(-1)^{2\sigma_{\alpha}+1}}{(2\sigma_{\alpha}+1)} \lambda^{3}_{\alpha\alpha} \rho_{\alpha}^{2} E(x_{\alpha\alpha})$$

$$- \frac{3\pi}{2\sqrt{2}} \beta \sum_{\alpha,\beta} e_{\alpha} e_{\beta} \kappa_{D} \rho_{\alpha} \rho_{\beta} \lambda_{\alpha\beta}^{3} Q(x_{\alpha\beta})$$

$$- \frac{\pi}{2} \beta^{4} \sum_{\alpha,\beta} \rho_{\alpha} \rho_{\beta} e_{\alpha}^{4} e_{\beta}^{4} \kappa_{D} \ln (\kappa_{D} \lambda_{\alpha\beta})$$

$$+ \frac{3\pi}{2\sqrt{2}} \beta \sum_{\alpha} \frac{(-1)^{2\sigma_{\alpha}+1}}{(2\sigma_{\alpha}+1)} \lambda_{\alpha\alpha}^{3} \rho_{\alpha}^{2} e_{\alpha}^{2} \kappa_{D} E(x_{\alpha\alpha})$$

$$+ \frac{1}{16} \sum_{\alpha} \frac{\beta^{2} \hbar^{2} e_{\alpha}^{2}}{m_{\alpha}} \kappa_{D}^{3} \rho_{\alpha} + \pi (\frac{1}{3} - \frac{3}{4} \ln 2 + \frac{1}{2} \ln 3) \times$$

$$\sum_{\alpha,\beta} \beta^{4} e_{\alpha}^{4} e_{\beta}^{4} \kappa_{D} \rho_{\alpha} \rho_{\beta}$$

$$+ C_{1} \sum_{\alpha,\beta,\gamma} \beta^{5} e_{\alpha}^{3} e_{\beta}^{4} e_{\beta}^{3} \kappa_{D}^{-1} \rho_{\alpha} \rho_{\beta} \rho_{\gamma}$$

$$+ C_{2} \sum_{\alpha} \beta^{6} e_{\alpha}^{3} e_{\beta}^{3} e_{\beta}^{3} e_{\beta}^{3} \kappa_{D}^{-3} \rho_{\alpha} \rho_{\beta} \rho_{\gamma} \rho_{\delta}$$

More in the Appendix!

Finally, let's not forget the physical issues in the equation of state

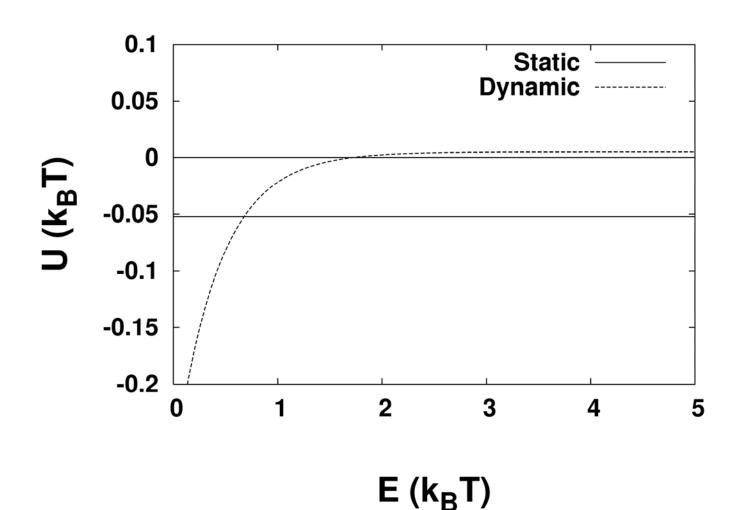
Various smaller competing - but relevant effects, among other:

- Existence and population of excited states
- Diffraction and exchange terms
- □ Parametric "size" in hard-spheres
- Relativistic correction for electrons

One current issue in nuclear reactions

Dynamic screening in solar and stellar nuclear reactions

Dynamic screening energy at the turning point for pairs of protons with a given relative kinetic energy [Mao et al. ApJ 701(2009)1204]



Screening energies and the ratio of screened to unscreened nuclear reaction rates for solar p-p reactions [Mussack & Däppen, ApJ, 729(2011)96]

Case	Screening energy U	Reaction-rate correction
Unscreened	0	1
Statically screened	$U_0 = -Z_1 Z_2 e^2 / R_D$	1.042
Dynamically screened	$U_0(E) = k_B T (0.005 - 0.281 \exp(-2.35E/k_B T))$	0.996

Our simulations suggest that dynamic effects obliterate the static-screening enhancement. In other words, they imply that at the the most probable collision energy (of 3-4 kT) the screening cloud is virtually absent and the naked Coulomb potential has to be tunneled through.

But,

Anderegg et al. [Phys. Plasmas 17(2010)055702] have used an apparent "duality" to infer the solar screening results from a system under totally different physical conditions, in their case magnetization experiments on a lasercooled ionic system. They think that their experiments confirm Salpeter's view of static screening enhancement.

Not so fast!

However, since that work is based on the assumption of a faithful mapping between two entirely different physical systems, and furthermore, since the present state of those experiments corresponds to a stronger coupled plasma than in the center of the Sun, this is unlikely the last word on the issue of dynamic screening.

Conclusions

- □ There are many stars
- They have different ages and chemical compositions
- Solar and stellar constraints will improve microphysics
- □ Thus better astrophysics and better physics from the Sun and the stars

APPENDIX -

Density and activity expansions

Look at the ideal Fermi-Dirac gas: the exact solution (but implicit!) ...

$$\frac{p}{kT} = \frac{1}{\lambda^3} f_{5/2}(\tilde{z})$$

$$\frac{N}{V} = \frac{1}{\lambda^3} f_{3/2}(\tilde{z})$$

$$\lambda = \frac{h}{\sqrt{2\pi mkT}}; \quad \tilde{z} = e^{\frac{\mu}{kT}} = \text{fugacity}$$

$$f_{5/2}(\tilde{z}) = \frac{4}{\sqrt{\pi}} \int_0^\infty dx \ x^2 \ln(1 + \tilde{z}e^{-x^2})$$

$$f_{3/2}(\tilde{z}) = \tilde{z} \frac{d}{d\tilde{z}} f_{5/2}(\tilde{z})$$

... and its high-temperature virial expansion

$$\frac{pV}{NkT} = 1 + \frac{1}{2^{5/2}} \frac{N\lambda^3}{V} + \dots$$

[obtained from expansion of Fermi integrals and elimination of \tilde{z}]

This is a nice explicit classical limit (and it includes the relevant criterion)!

Expand without eliminating: fugacity/activity expansion

$$\begin{split} \frac{p}{kT} &= \frac{1}{\lambda^3} \left[\tilde{z} - \frac{1}{2^{5/2}} \tilde{z}^2 + \ldots \right] = z - \frac{\lambda^3}{2^{5/2}} z^2 + \ldots \\ \frac{N}{V} &= \frac{1}{\lambda^3} \left[\tilde{z} - \frac{1}{2^{3/2}} \tilde{z}^2 + \ldots \right] = z - \frac{\lambda^3}{2^{3/2}} z^2 + \ldots \end{split}$$

Which is based on the natural definition of the activity

$$z = \frac{1}{\lambda^3}\hat{z}$$

Any real system has an activity expansion (ACTEX) ...

$$\frac{p}{kT} = z - b_2 z^2 + \dots$$

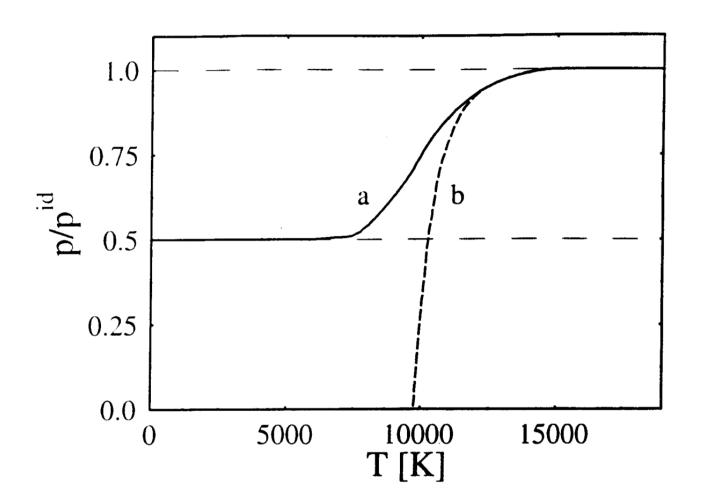
$$\frac{N}{V} = z - 2b_2z^2 + \dots$$

... because the grand-canonical partition function delivers these coefficients b_i (of course infinitely many)

Radically different behavior already to first

non-ideal order (here illustrated with H-H2 system)

(a: activity b: virial)



For reacting systems, activity expansions are much better suited than virial expansions!

See: Kremp D., Schlanges, M. & Kraeft, W.-D., Quantum Statistics of Nonideal Plasmas (Springer, Berlin, 2005)