

· So how	does to	dilute Fern	i sustem	with	8-Punction
interaction	on behav	<u>e? (In?</u>	o ed also	1 dim	enslon.)

For a fixed number of Fermions N, it is relevant to consider the energy per particle [E(g) = EN]
as a function of density [p= NV].
. We've suntched to V for volume instead of Ω to

evoid confusion, below! (Sorry!)

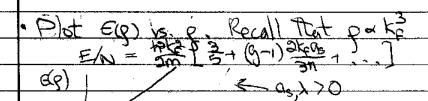
• E(P) is an intensive quantity: if he double

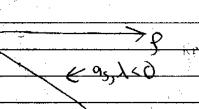
N and V, E(P) stays the same.

• The other possibility would have been the

energy density: [Etp) = E/V, which is also intensive but not as useful here.

· Note that [E(p) = E(p)/p





$$\alpha b = b \frac{gb}{gb}$$

so positive or negative slope corresponds to the sign of the pressure.

V ...

Review of Hermodynamics Statistical Mechanics

· To go to even the next order of perturbation theory is awknown and, in general, we will next to develop non-perturbative approximations.

> reed a new Framework.

We will use path integral methods.
Later will show the connection to operator bused approaches.

· The most convenient function to calculate with many-body path integrals is the grand canonical partition function.

> revue some thermo and stat. mech.

NOTE: Volume is V in this section, not I.

Recall: three ensembles

i) microcanonical: Fixed E and N

ii) canonical: system exchanges energy with heat both = Fixed N and fixed exercise = aft

Probability of system alone with E is e-E/ST = e-BE B= For

ili) grand cononical: system exchanges energy and particles with heat porticle path. Neither N mr E fixed.

Probability of system afone have E and N is

Probability of system afone have E and N is

Probability of system afone have E and N is

· Suppose we have the complete set of energy and particle number eventations: AINJUZ = EJMINJUZ; NINJUZ = NINJU

Hen Re grand portition Runction present potential $\Omega(T,V,\mu)$ by

 $\int \Omega(T,V,\mu) = -\frac{1}{2} \ln Z_G$

and the second	1/13/03
	Let's recall the laundry list of thermodynamic functions intensive variables: TPP µ
	intensive variables: TP p
	extensión variables: S V N
	energy t= t(S,V,N)
	energy $F = E(S, V, N)$ Helmholtz free energy $F = F(T, V, N) \cdot minimized for a mechanically isolated system at constant T Gibbs free energy G = G(T, P, N) \cdot equilibrium at constant T and P when minimized$
	Coibbs free green G = G(T, P, N), earlibrary at anstart T
	and P when minimized
×	Themodynamic potential DLT, V, M)
.	First land of thomas inco executive has the internal
	First law of thermodynamics specifies how the internal every changes with changes in independent variables:
	$dE = TdS - PdV + \mu dN$
<u> </u>	which implies T= (3E) VIN P= (3E) SIN U= (
	· At T=0, in the ground state, S=0 => / u=(3/2)v
	To obtain the other thermodynamic functions from E, we perform Legendre transformations
	3 change the independent variables.
	schange the independent variables. "It's important to understand these, because we'll be doing the functional generalization later
	be doing the tractional generalization later
	[F=E-(\$), S=E-TS => SF=-SST-PNV+pdN]
·	G=F+PY=E-TS+PV => dG=-SOT+VAP+udN



_	V13/63 V13/63
	So we see that we can generate all the equilibrium thermodynamics in we know $\Omega(T,V,\mu)$.
	Return to the partition function!
	$Z_{G} = \sum_{N \in N} e^{-\beta(E_{N} - \mu N)}$ since diagonal in this
	$= \sum_{N \in \mathbb{N}} \langle N $
	$= Tr(e^{-\beta(\widehat{H}-\mu\widehat{n})})$
	Here To is the trace over any complete basis => much more general Than first expression.
	$\mathcal{L}(T,V,p) = -\frac{1}{5}\ln z_G \Rightarrow \left[z_G = e^{\beta \mathcal{L}(T_0V_1p)} - T_r(e^{-\beta(\widehat{H}-p\widehat{N})})\right]$
j	The ensemble occurre or expectation value of an operator is $ \begin{array}{c} \hline $
-	
1	· Consider on exchape: 0= 1) (R) = Tr(R) = P(H-µN) = P(E-µN) = P(
	Jeck, puse = by by = -30 = N
	This is a prototype of things we will do with path integrals to find expectation valves.

	V13/03
	· Let's evaluate the trace in Zo lising The occupation
	· Let's evaluate the trace in Zo Using the occupation number bosis. In general, these are not eigenstates of A, but let's start with
	$H \rightarrow H_0 = \Xi \in [a, a]$ $N = \Xi a, a$
1 44.00	where i runs over the single-particle quantum
	where i runs over the single-particle quantum numbers (e.g. R, x for fermions in a box),
	We'll see that the trace is easy to evaluate in this case. This in turn, will lead to a strategy for evaluating the more general case.
- m- L/M d	evaluating the more general case.
<u> </u>	"In the occupation number basis, we don't sum over Esm and N, but over 12, 12, the mades" (single-particle states) 1, 2, 3,
	$ \frac{1}{26 = 10} = \frac{1}{6} \frac{\left[\frac{1}{100} - \frac{1}{100} \right] \left[\frac{1}{100} - \frac{1}{100} \right] \left[\frac{1}{100} - \frac{1}{100} \right] \left[\frac{1}{100} - \frac{1}{100} \right]}{\left[\frac{1}{100} - \frac{1}{100} \right] \left[\frac{1}{100} - \frac{1}{100} \right]} = \frac{1}{100} = \frac{1}{100} \left[\frac{1}{100} - \frac{1}{100} \right] \left[\frac{1}{100} - \frac{1}{100} - \frac{1}{100} - \frac{1}{100} \right] \left[\frac{1}{100} - \frac{1}{100} - \frac{1}{100} - \frac{1}{100} - \frac{1}{100} \right] \left[\frac{1}{100} - \frac{1}$
	Recall Plat Ing-now)=(n,7 in 2). For bosons, each n;=0,1,2,
	$\Rightarrow \left[\widehat{H}_{0} n_{1} \cdots n_{\infty} \right] = \left[\underbrace{\varepsilon_{i} \alpha_{i}^{\dagger} \alpha_{i}^{\dagger} n_{i} \cdots n_{\infty} \right]}_{\varepsilon_{i} \varepsilon_{i} \varepsilon$
	$ \hat{N} _{\Omega_1,\dots,\Omega_\infty,7} = \sum_{i=1}^\infty \hat{N}_i _{\Omega_1,\dots,\Omega_\infty,7}$
	· Recall Plat [\hata_i, \hata_j] = 0 for all i', j, which means we can exponentiate Plese results.
	<u>.</u>

V13/03 Plugging in boson or fermion only 1 μ ! $\frac{1}{26} = \sum_{n=0}^{\infty} \left(\frac{1}{2} - \frac{1}{2} \right) \cdot \left(\frac{1}{2} - \frac{1}{2} - \frac{1}{2} \right) \cdot \left(\frac{1}{2} - \frac{1$ Plugging in, bosons: n=0,1,2, = 20 = ple;-pln; = 1+ = ple;-pl = 2ple;-pl) (bosons) = 1- polery) geometric series, Fermions: n=0,) \(\frac{1}{2} \in \text{ple:-ple:-pln:} = 1 + \in -\text{p(E:-p)} \) (Fermions Drott, V, p) = - pln 2 =)+1 2 ln (1- = p(e-p)) - 1 S In (1+Ep(E, - M)) from Line => just two little sign differences! = \(\frac{1}{2}\) \(\frac{1}{2 where no is the mean occupation number in the its state Fermions: (1) = - (= = 1+ E p(E;-h) E P = + \(\frac{1}{p\left(\frac{1}{p\reft(\frac{1 no = (ep(Et-p)+1)-1 for fermions