

Getting the Most Out of *SnB*

Russ Miller

Hauptman-Woodward Med. Res. Inst.

Principal Contributors:

C.-S. Chang

G.T. DeTitta

S.M. Gallo

H.A. Hauptman

D.A. Langs

R. Miller

S. Potter

C.M. Weeks

Partial funding from NIH and NSF.

Overview

- *Shake-and-Bake*
 - Direct Methods Optimization Technique
 - Multiple Trial Structures
 - Real Space \Leftrightarrow Reciprocal Space
- *SnB* is the program based on *Shake-and-Bake*
- Phase Refinement Techniques
 - Parameter Shift
 - Tangent Formula
 - Simulated Annealing
 - Genetic Algorithms

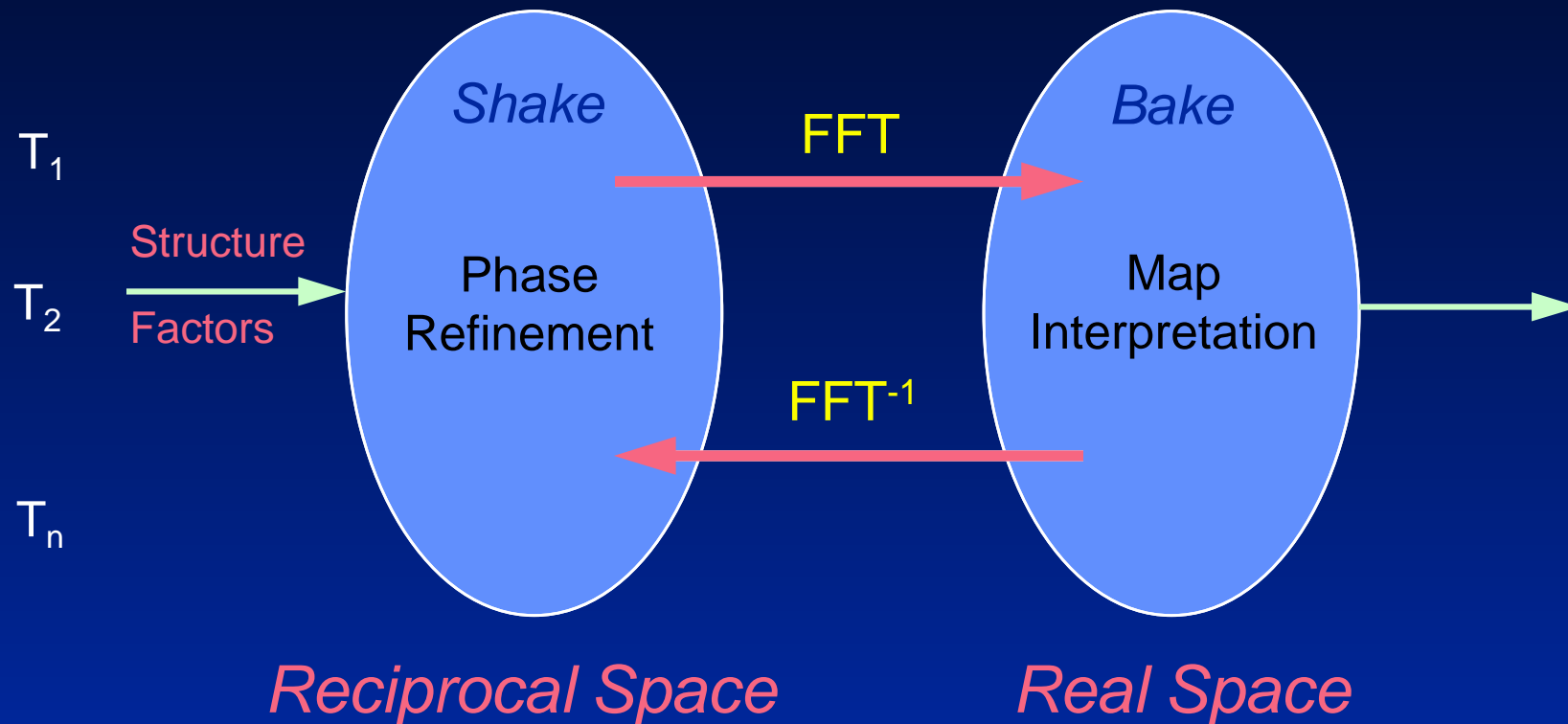
Computing Platforms

- Unix Workstations
 - SGI
 - Sun
 - DEC/Alpha
 - IBM RS6000
- Thinking Machines Corp. CM-5
- Cray T3D
- Cray C90
- IBM SP2

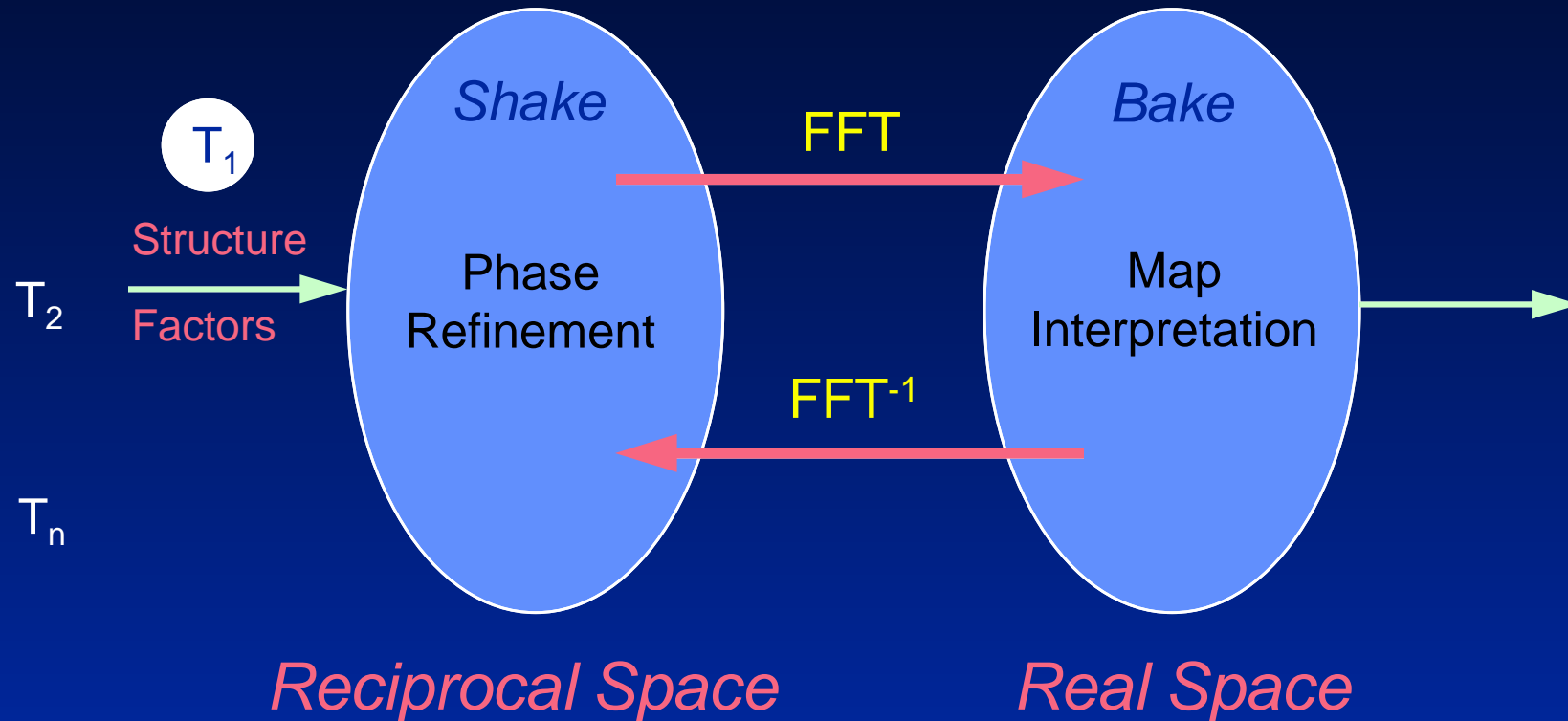
Availability of *SnB*

- MSC's teXsan 1-7.2
- MacScience
- Pittsburgh Supercomputing Center
- Hauptman-Woodward Inst.
- Cornell Theory Center

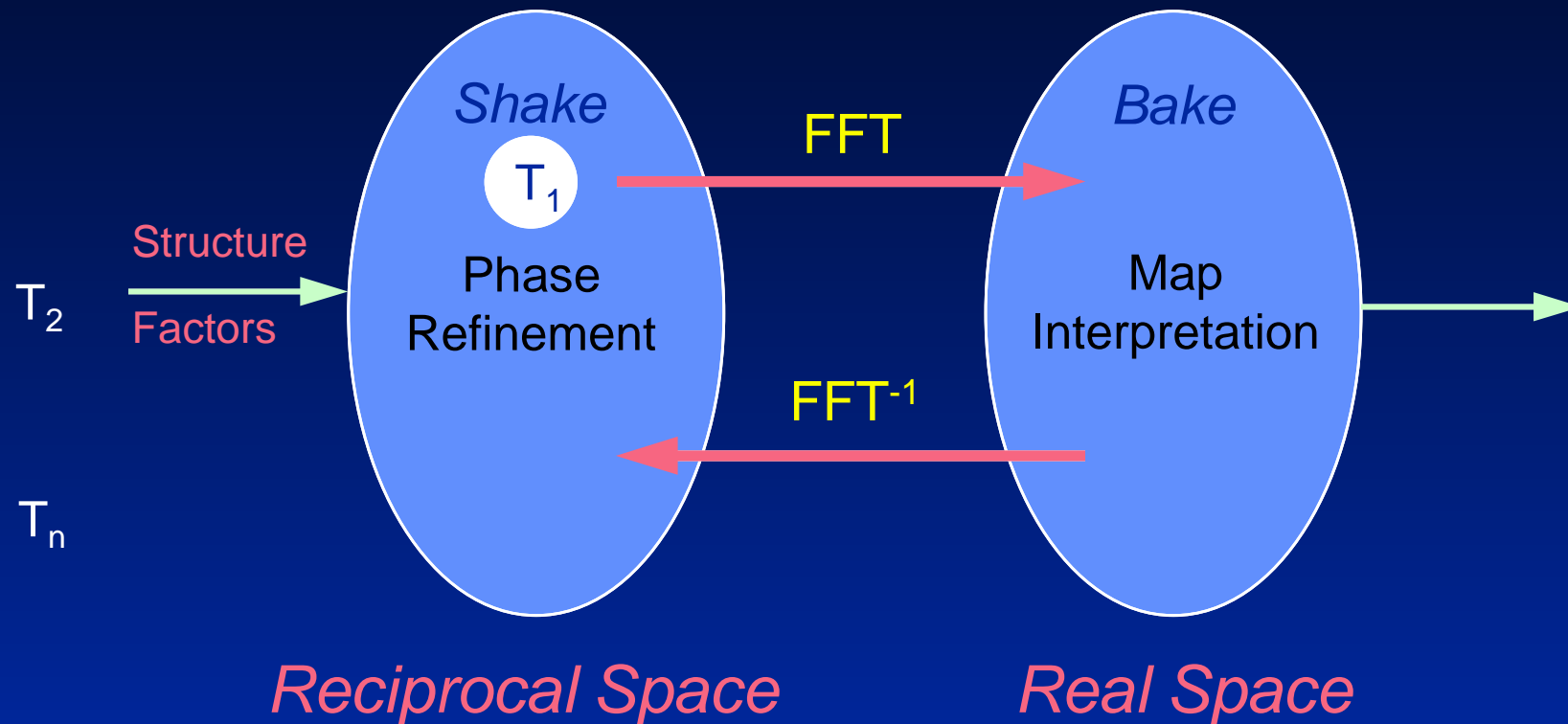
Shake-and-Bake



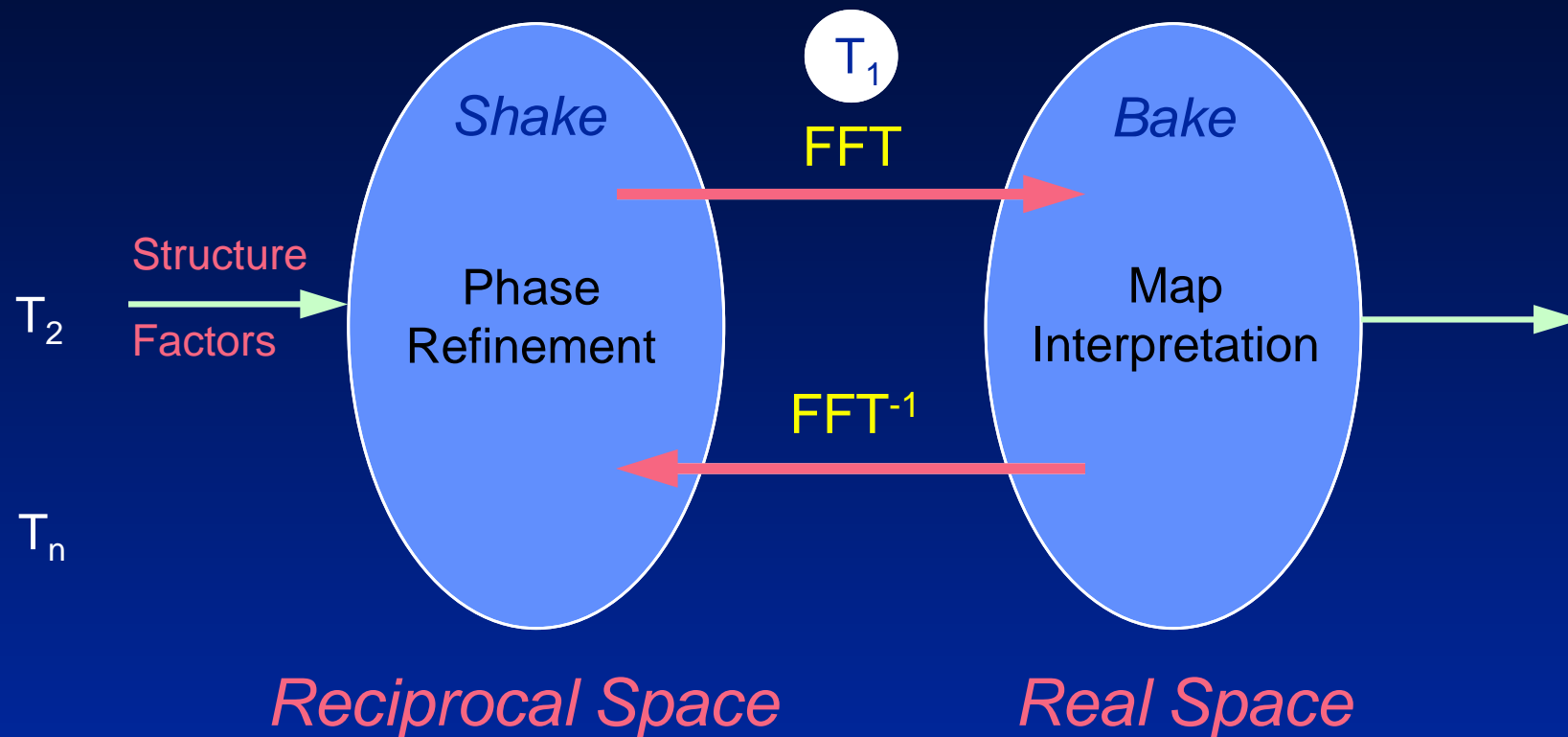
Shake-and-Bake



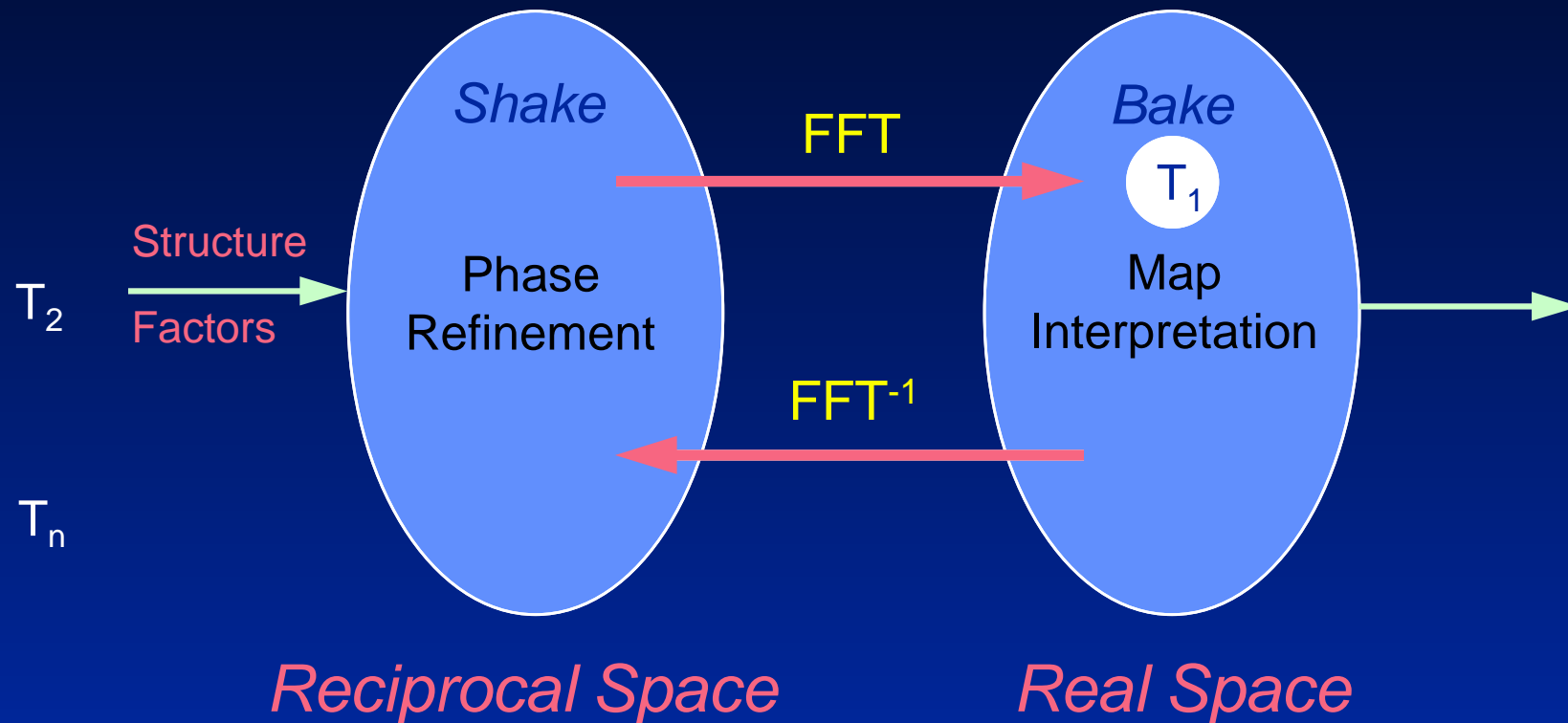
Shake-and-Bake



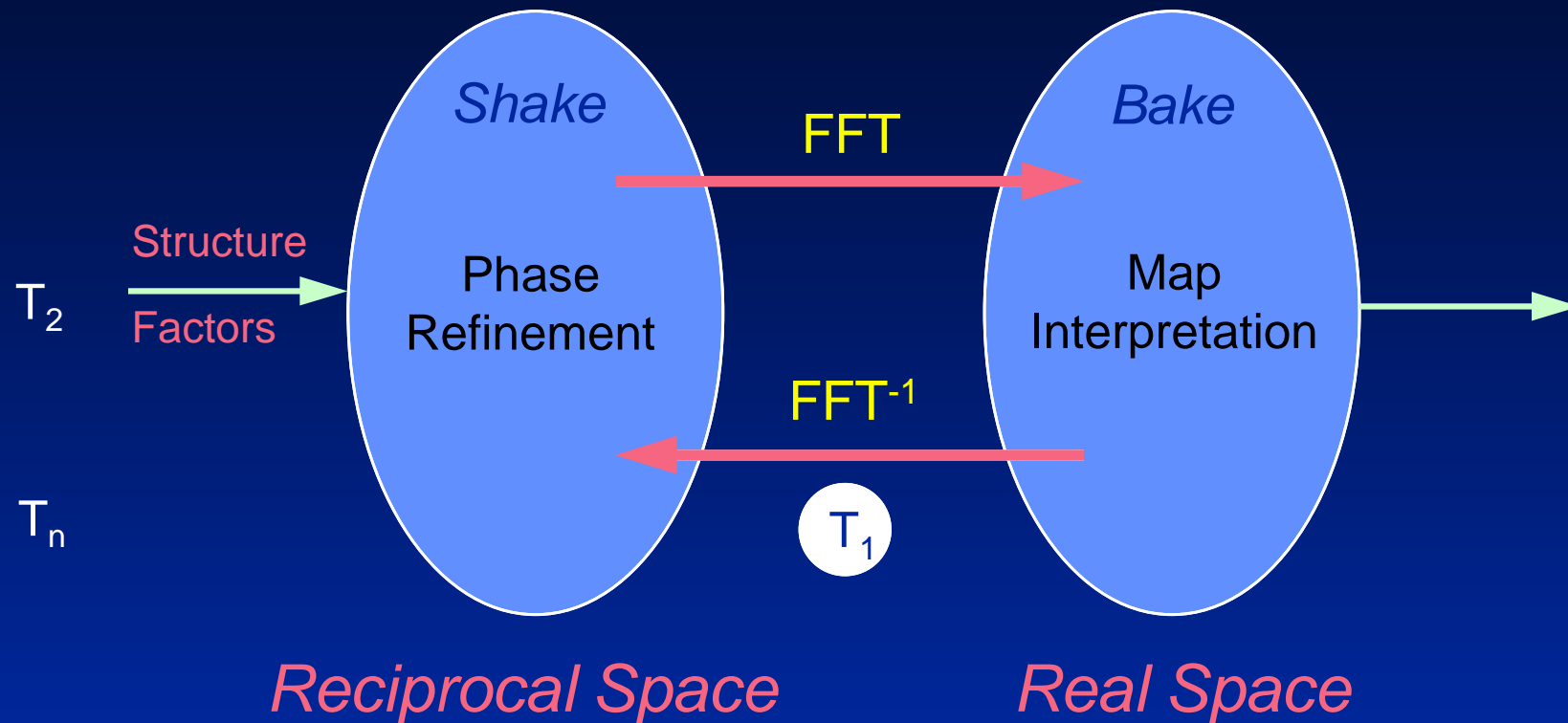
Shake-and-Bake



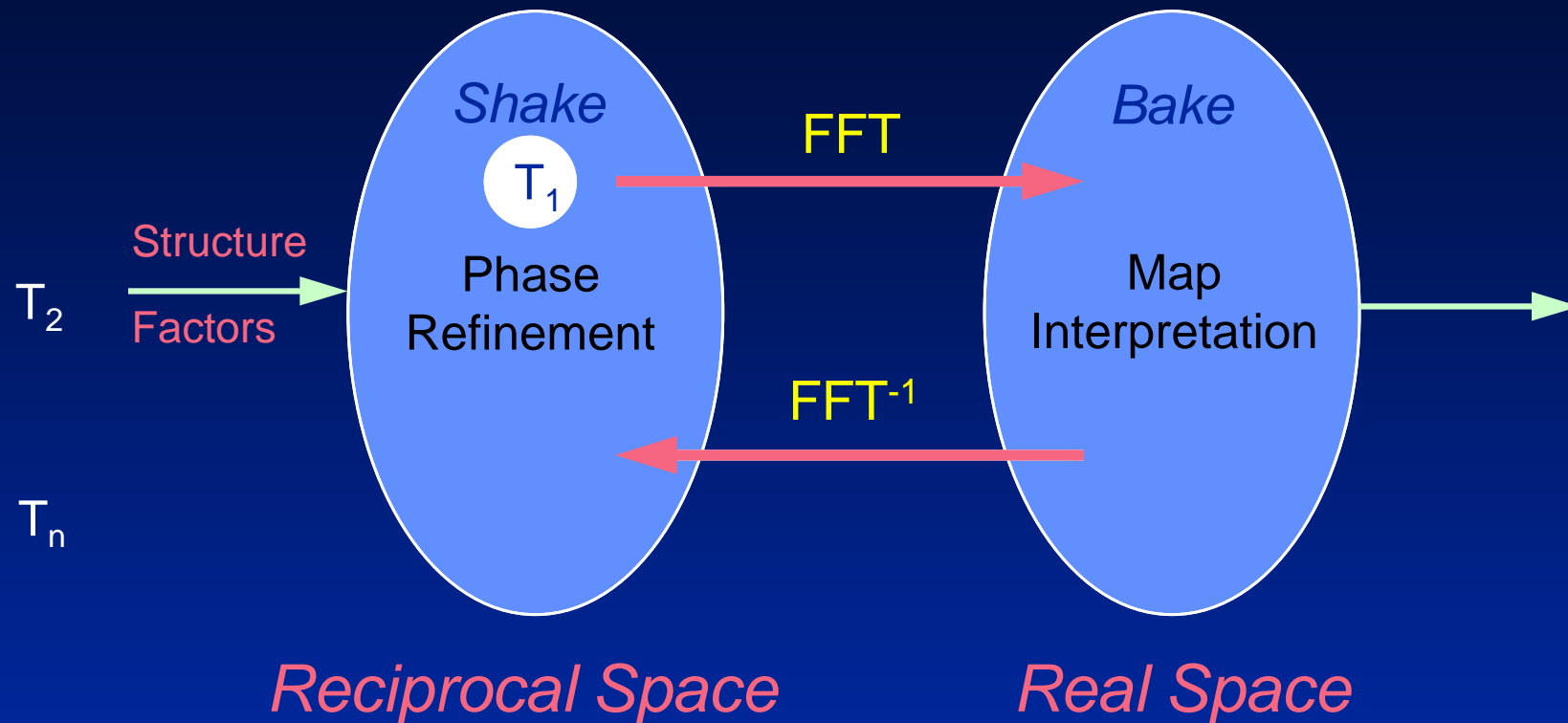
Shake-and-Bake



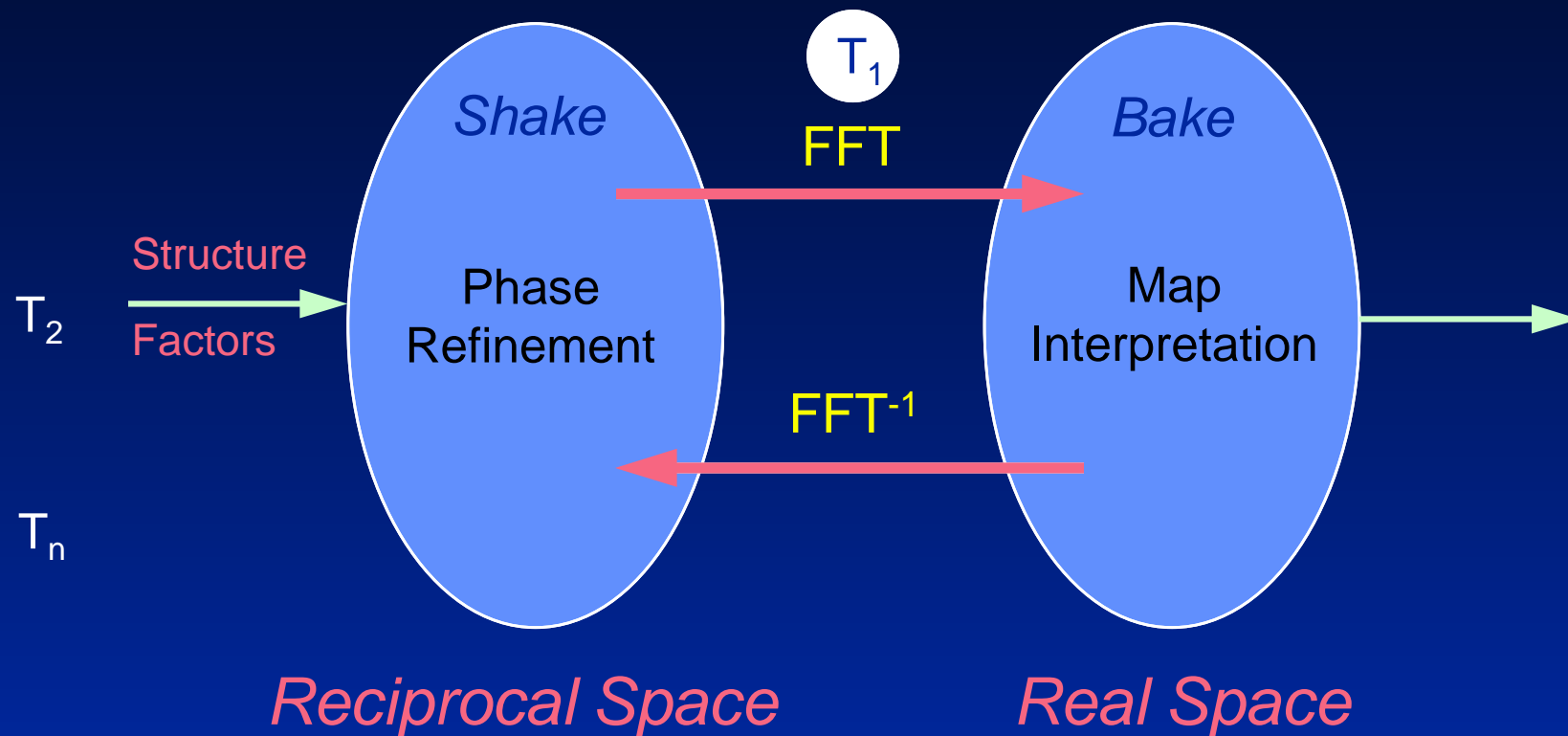
Shake-and-Bake



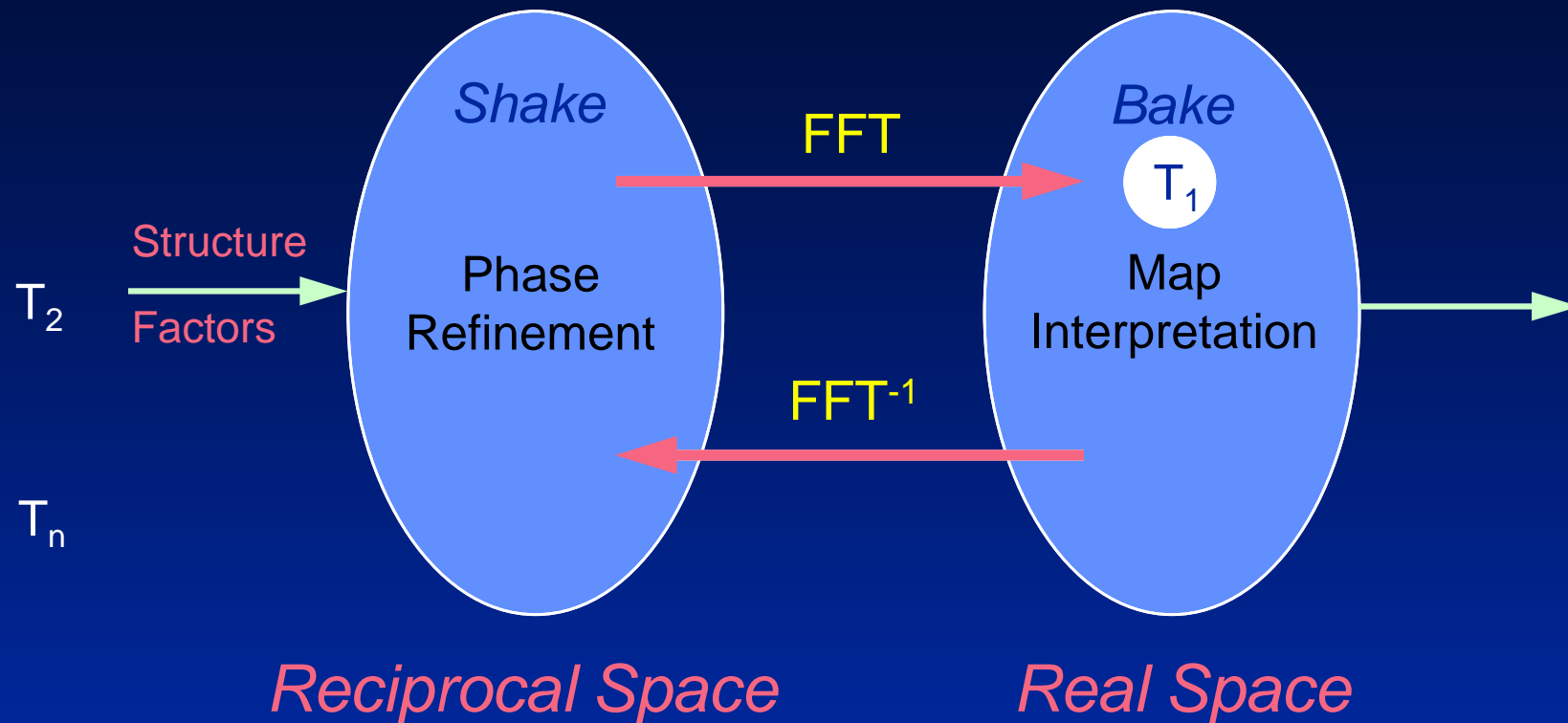
Shake-and-Bake



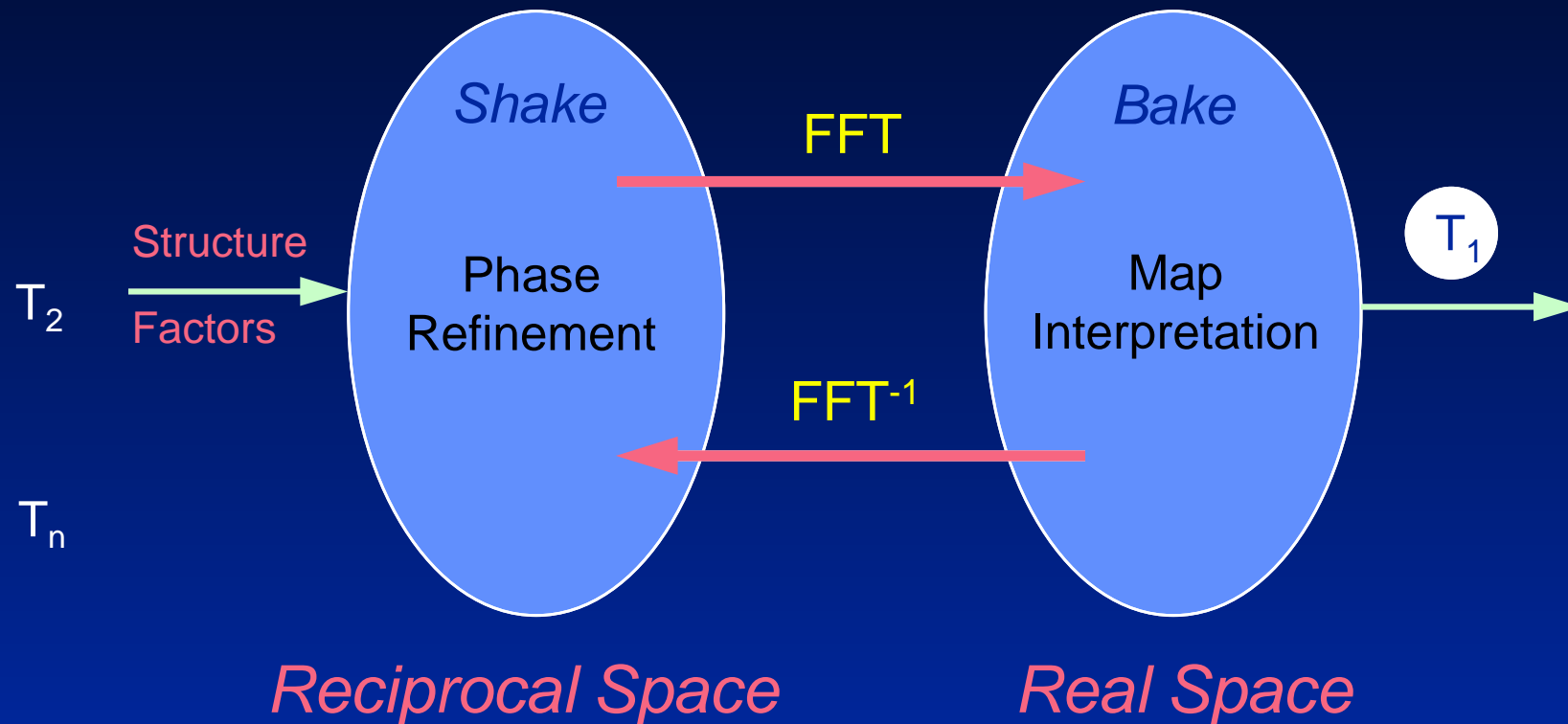
Shake-and-Bake



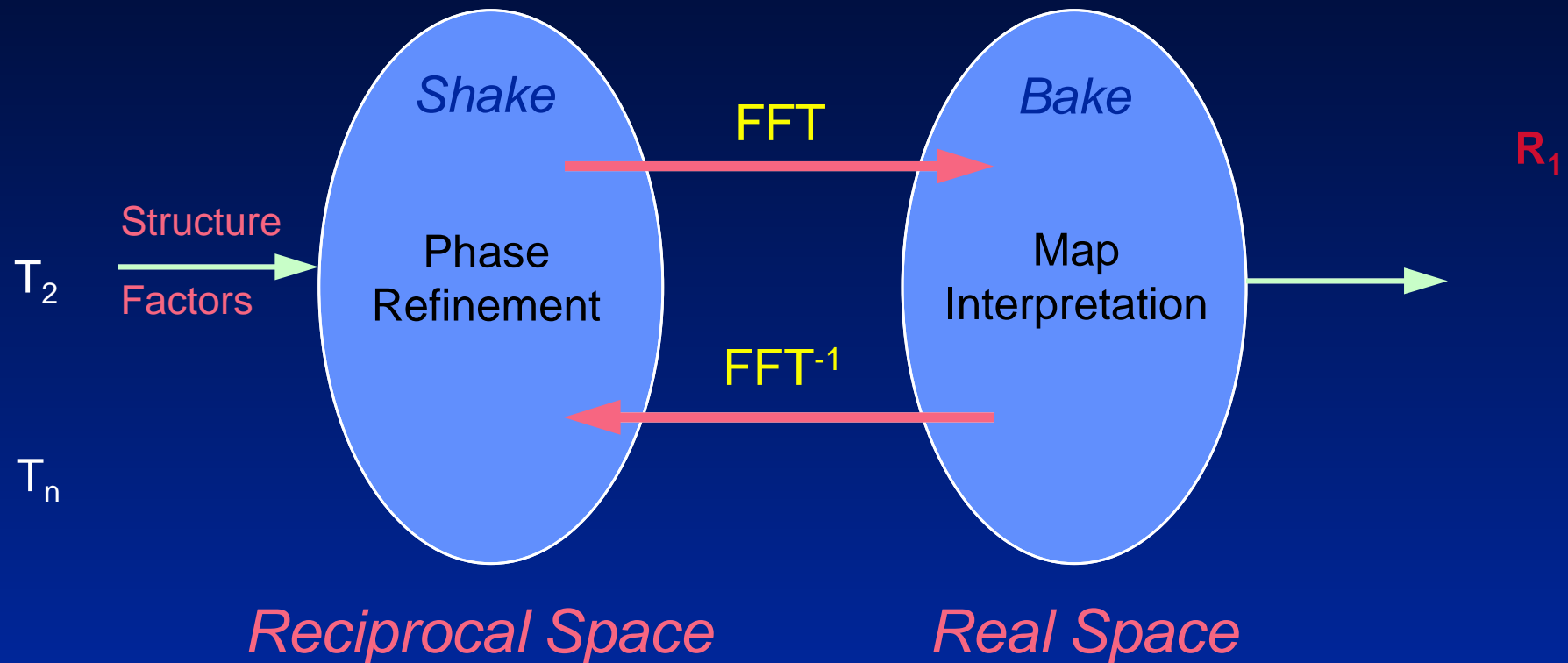
Shake-and-Bake



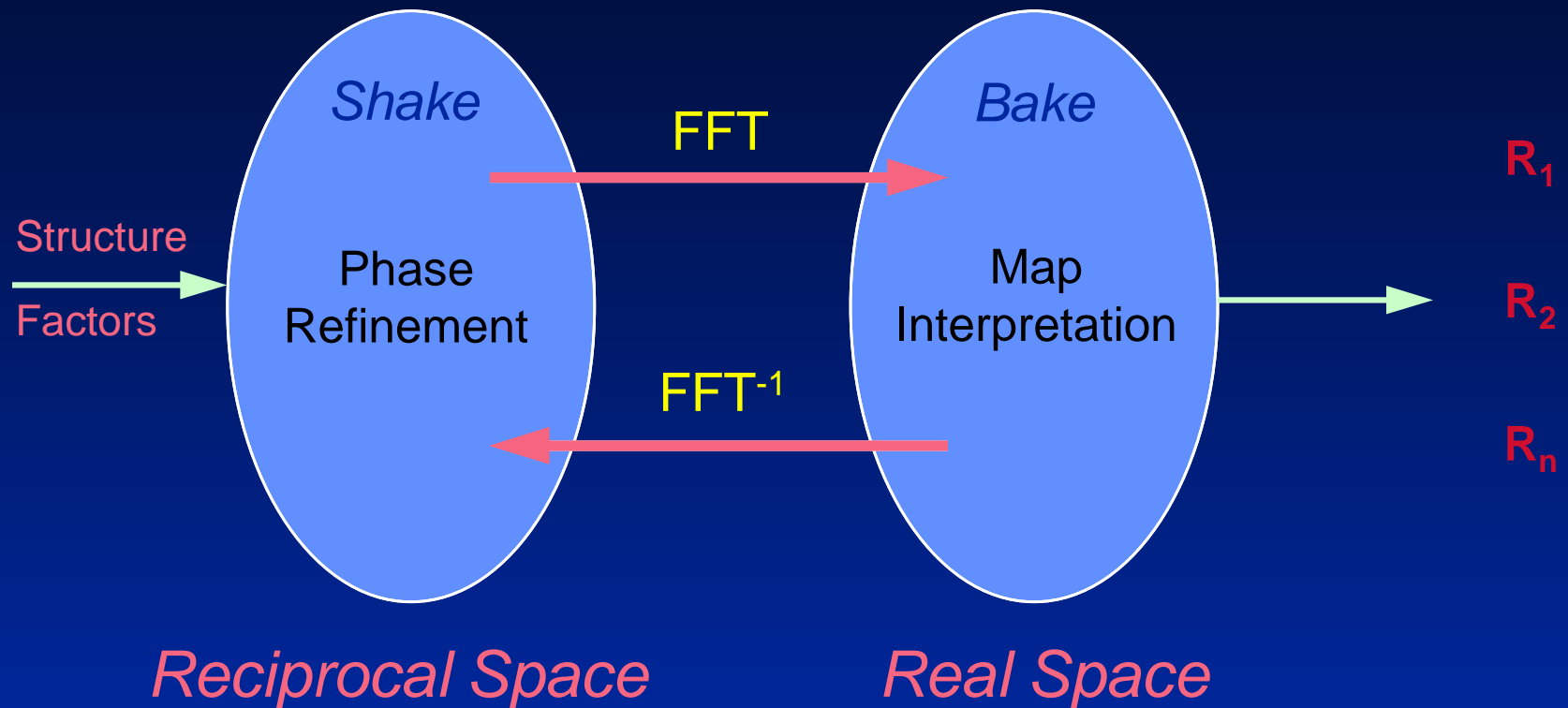
Shake-and-Bake

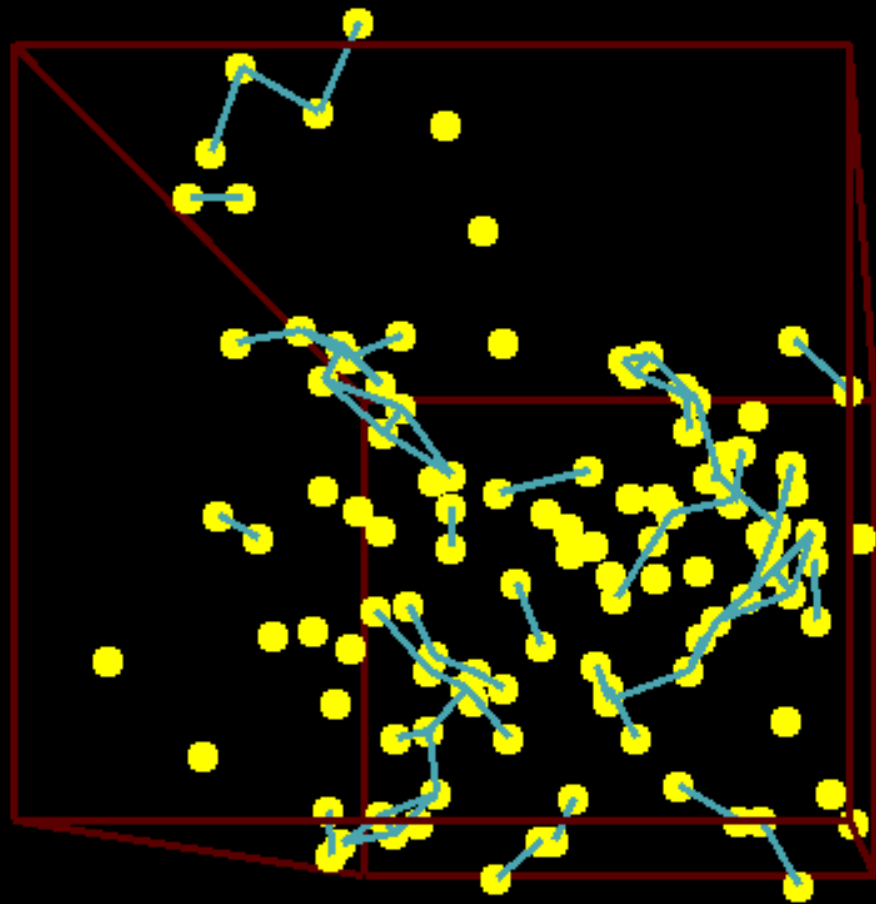


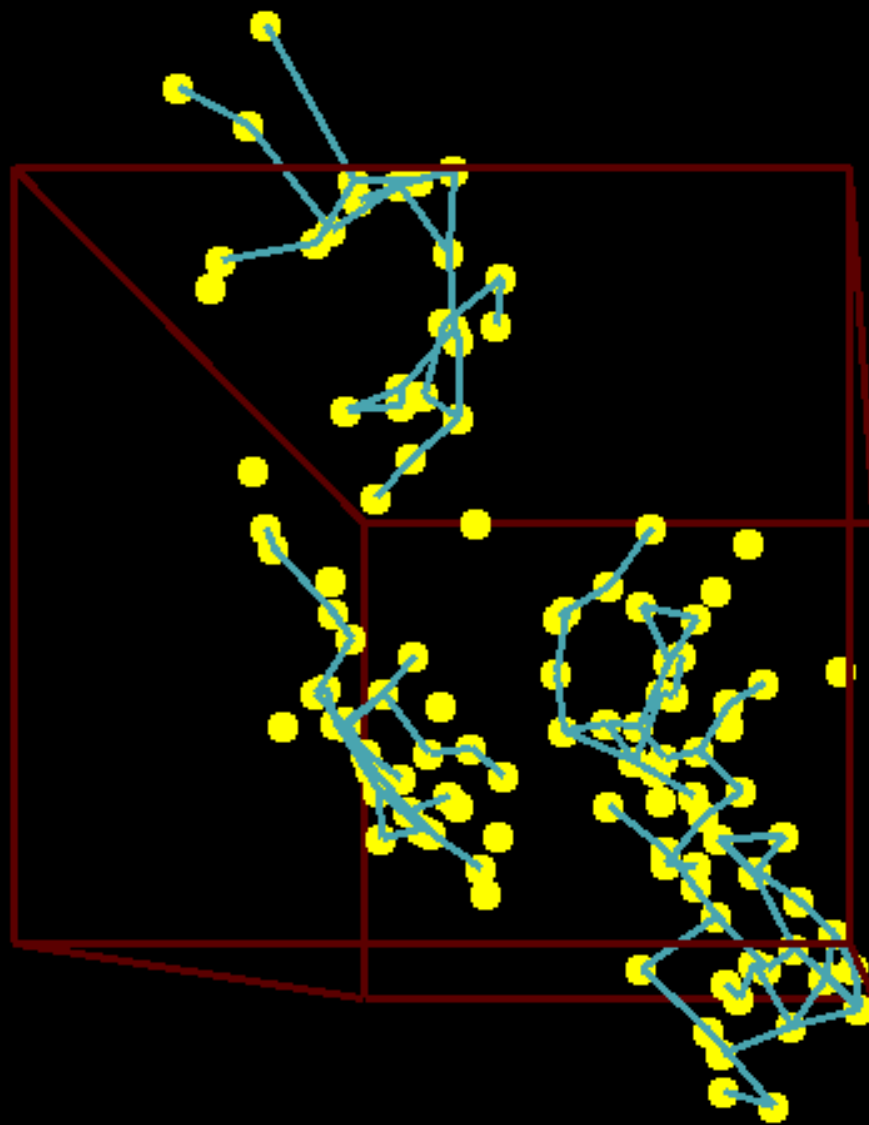
Shake-and-Bake

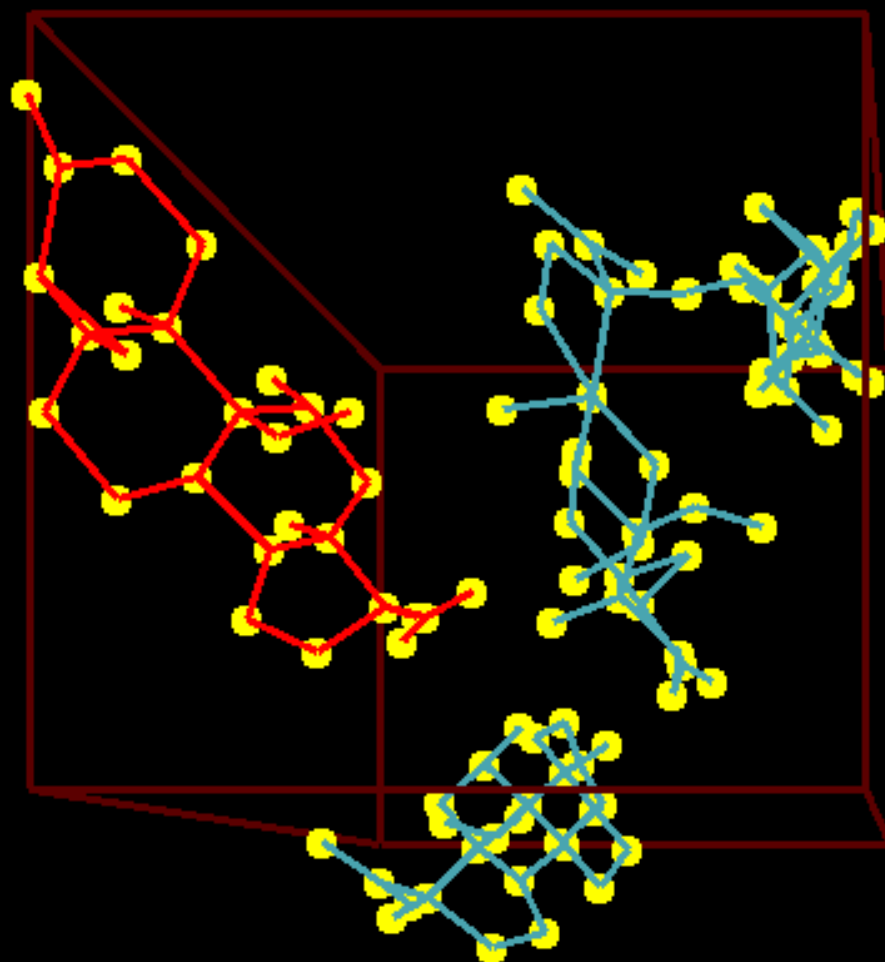


Shake-and-Bake

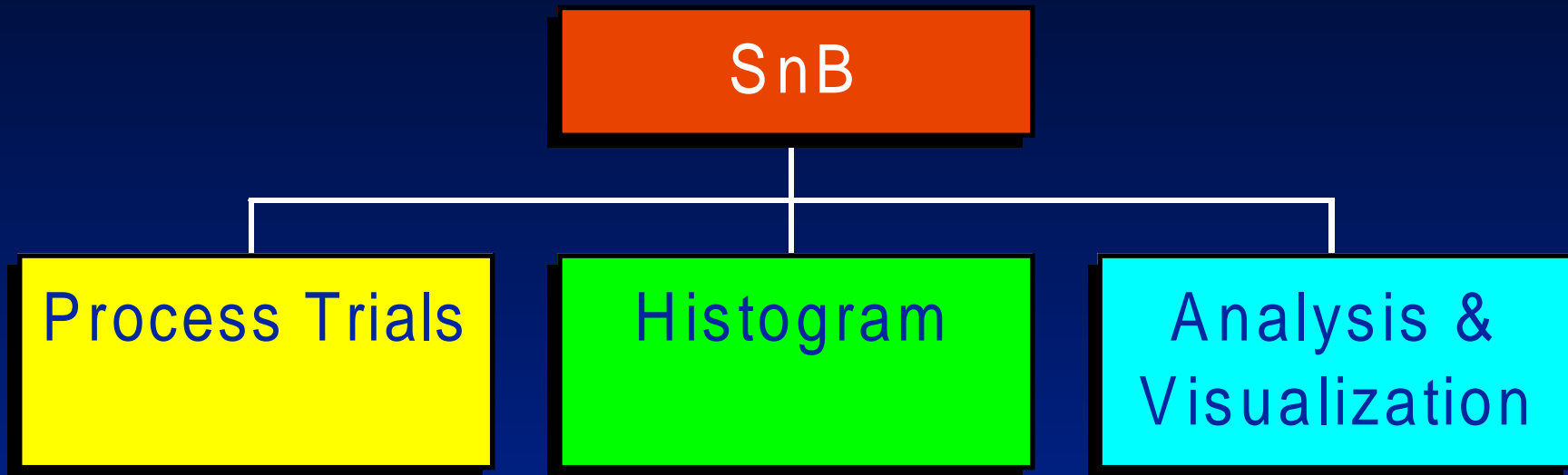


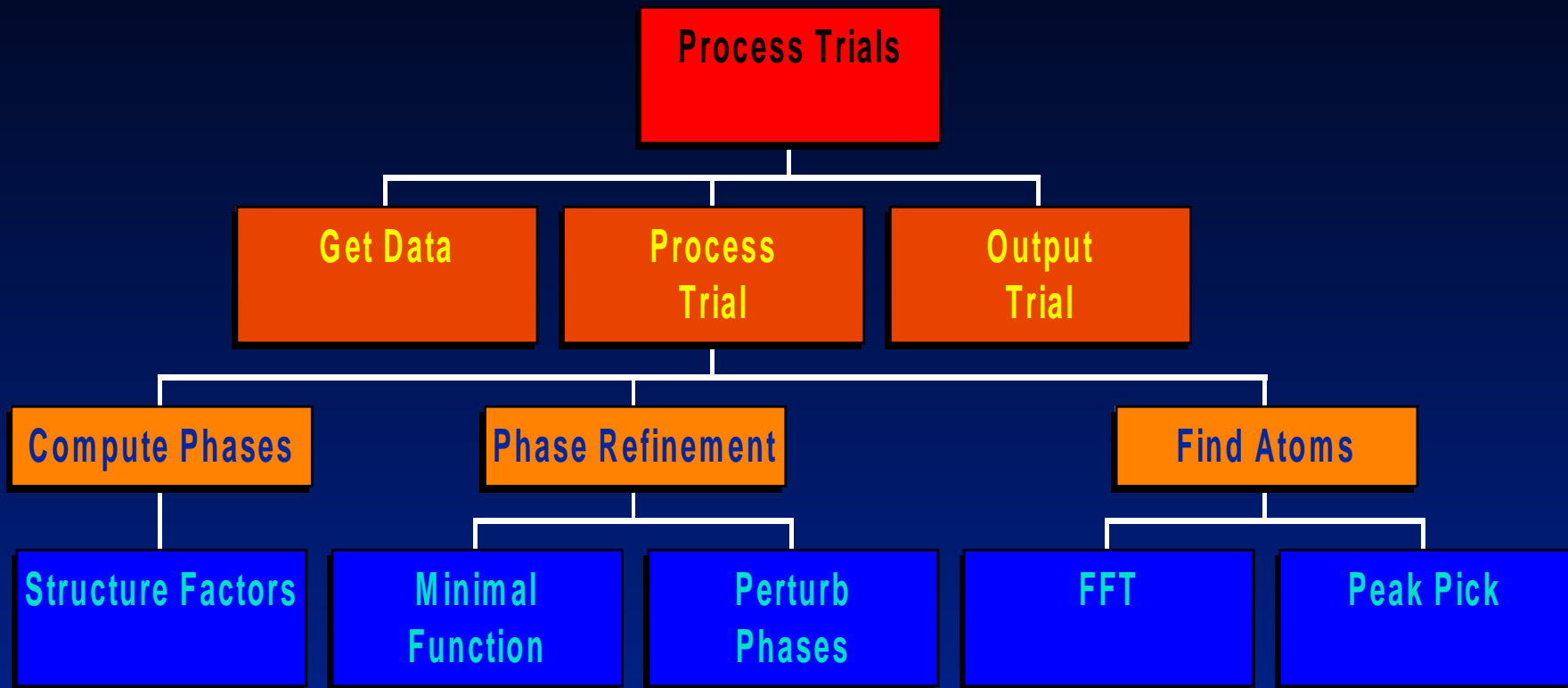






Structure of *SnB*





SnB Default Parameters

Atoms in asymmetric unit	n
Number of Phases	$8n - 10n$
Number of Triples	$70n - 100n$
Number of Quartets	0
Number of Cycles (Parameter Shift, Cent. Search)	$n/2$
Number of Cycles (Tangent Refinement)	$n/4$
E-Fourier Recycling Steps	C

Crambin

Atoms in asymmetric unit	400
Number of Phases	4,000
Number of Triples	40,000
Number of Quartets	0
Number of Cycles (Parameter Shift)	200
E-Fourier Recycling Steps	5

Structures

<i>STRUCTURE</i>	<i>ATOMS</i>	<i>SPACE GROUP</i>	<i>SUCCESS RATE</i>
Ternatin	110	P2₁2₁2₁	2%
Scripps	144	P1	
MSC	180	P1	(8 years with traditional methods)
UCLA	302	C2	
Gramicidin A	317	P2₁2₁2₁	0.3%
Crambin	400	P2₁	4%
OCI / U. of T.	408	P1	5%
Rubredoxin	500	P2₁	2.7%

Cost-Effectiveness

- Tangent Formula in SnB
 - Smaller structures
 - Carefully chosen parameters
- Space Group
 - P1
 - P-1
- Simulated Annealing
- Genetic Algorithms

Running Times

(Average Seconds/Cycle on Iled)

Machine	Total	StrFac	Refine	FFT	Pk Pk
SGI Pwr Ind 2	1.247	0.210	0.426	0.485	0.068
DEC/Alpha	1.822	0.437	0.682	0.520	0.075
SGI Indigo	2.975	0.881	0.760	1.037	0.163
Sparc 10	3.140	0.821	0.919	1.111	0.162
Pentium 75	3.535	1.181	0.995	1.205	0.154
Sparc 5	3.759	0.765	1.087	1.661	0.210
Pentium 90	3.960	1.192	1.172	1.362	0.192
i486dx2-66	11.396	2.175	2.423	5.467	0.711

Additional Applications of *SnB*

- **Fragment Recycling (D.A. Langs)**
 - Large radius of convergence
 - Superior to traditional tangent formula (RANTAN)
- **Electron Diffraction Data (Dorset & McCourt)**
 - High success rates
 - Small % of data out of principle plane

Related Presentations

H.A. Hauptman	Last Night	1e.2.A
C.M. Weeks	Monday/Tuesday	M137
G. Prive	Wednesday/Thursday	W008
D. Anderson	Thursday Afternoon	4a.7.A

SnB Addresses

- <http://www.hwi.buffalo.edu/SnB>
- snb-requests@hwi.buffalo.edu
- snb-comments@hwi.buffalo.edu

Summary

- Alternative structure determination method
- Targeted at 100-800 atom structures
- Publicly available
- Recommend: LEVY / EVAL (Bob Blessing)
- Current Research
 - Parameter tuning / Code Optimization (2X)
 - Additional FOMs
 - Lower resolution data
 - SIR/SAS data