

Study of weak-coupled states in ^{23}Al via inelastic scattering

-- Using low-energy RI beam at CNS --

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Outline

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- **Experiment**

Method, RIB prod. & Setup

- **Results**

Spectra, inelastic analysis, new level scheme

- **Model Calculations**

R-matrix analysis & Shell model calc.

- **Summary**

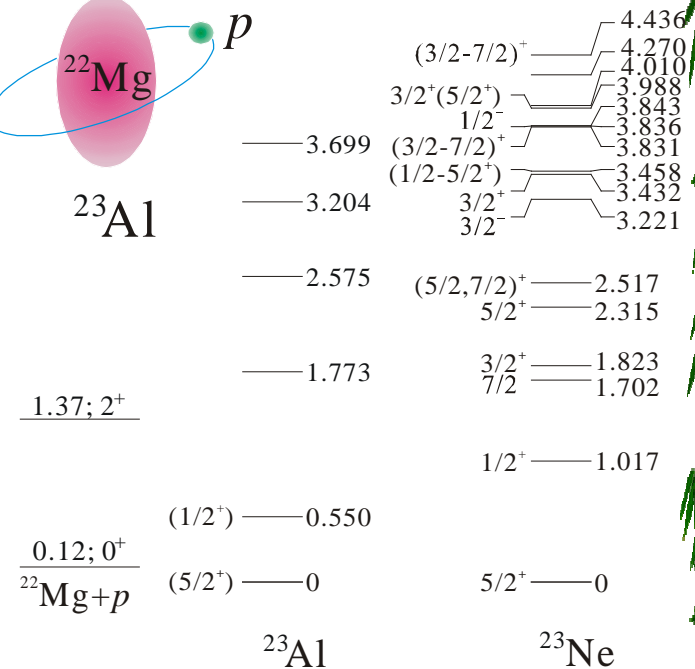
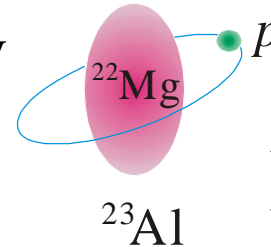


Introduction

(1) Aim: Weak-coupled structure in ^{23}Al

$$S_p(^{23}\text{Al}) = 0.125 \text{ MeV}, S_p(^{22}\text{Mg}) = 5.502 \text{ MeV}$$

$$\beta_2(^{23}\text{Al}) \approx 0.3 \sim 0.6, \beta_2(^{22}\text{Mg}) \approx 0.56$$



(2) Status: Structure in ^{23}Al

(a) $^{24}\text{Mg}(^7\text{Li}, ^8\text{He})$ @MSU;

(b) β -delayed proton decay of ^{23}Si @GANIL;

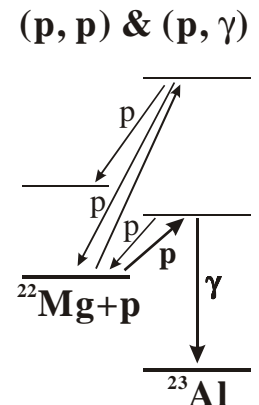
(c) Coul. Dissoc. of ^{23}Al @ RIKEN

Conclusion: Several levels were observed without spin-parity assignments.

(3) Present work

E_{cm} (^{23}Al) up to 4 MeV (by resonant elastic & inelastic scattering)

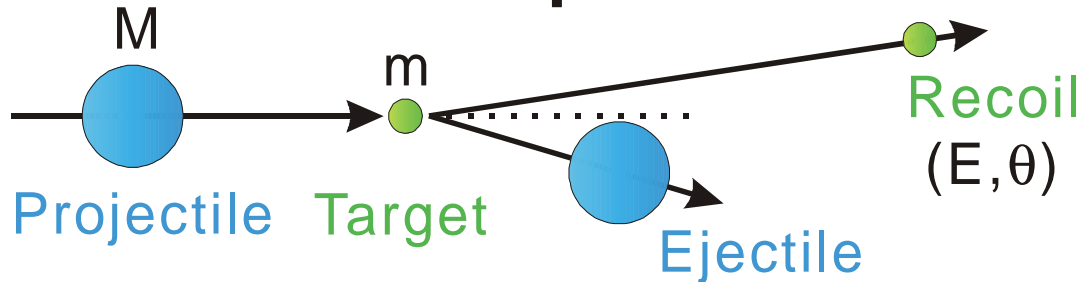
$E_R, J^\pi, \Gamma_p (\Gamma_p')$ \rightarrow Weak-coupled structure of ^{23}Al



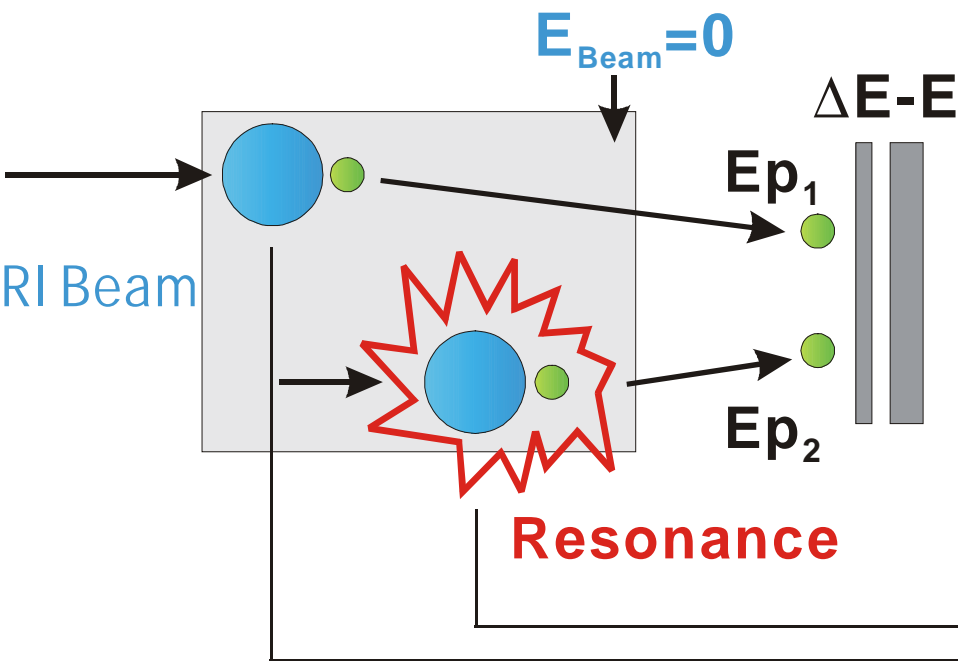
Experiment

(1) Experimental Method

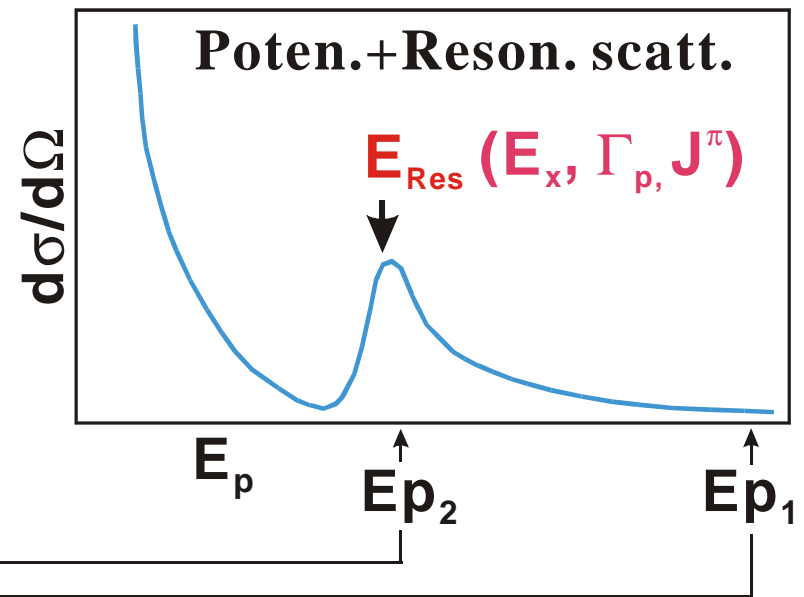
As for proton elastic scattering



$$E_{cm} = \frac{A+1}{4A \cos^2(\theta_{lab})} * E_p$$
$$E_x = E_{cm} + Q$$



Excitation Function



(2) RI Beam Production

CNS RI Beam Separator (CRIB)

Production Reactions

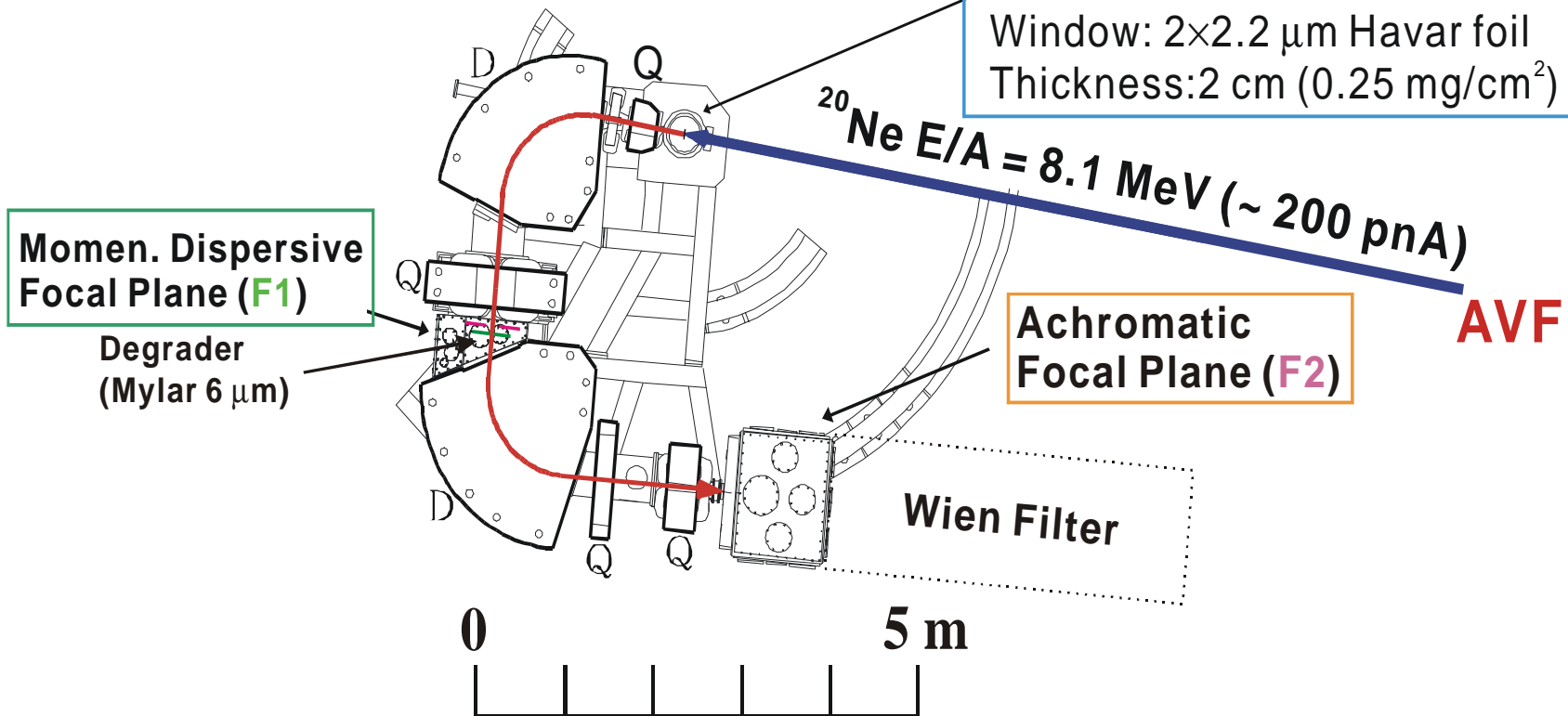


${}^3\text{He}$ gas target

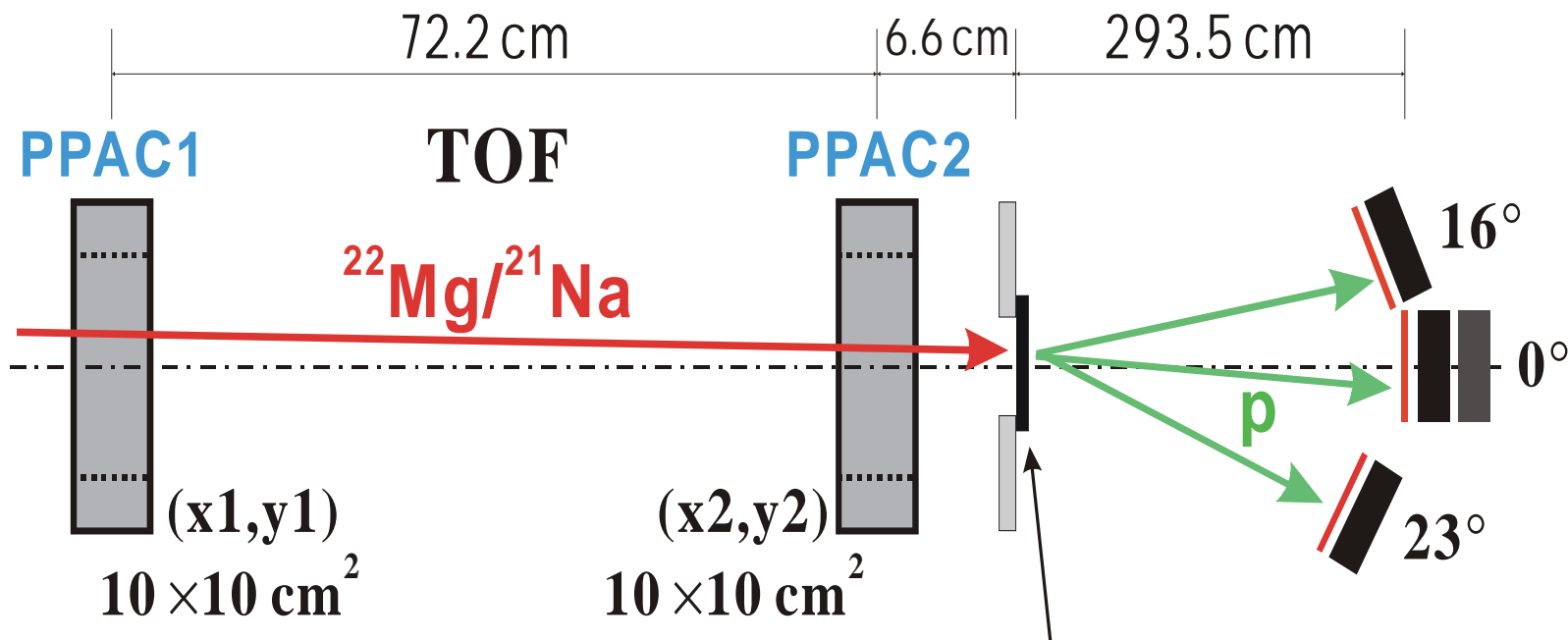
Pressure: 1 atm.

Window: $2 \times 2.2 \mu\text{m}$ Havar foil

Thickness: 2 cm (0.25 mg/cm^2)



(3) Setup for Elastic Scattering



After PPAC2

^{22}Mg : 4.4 AMeV
6.6 kaps (3%)

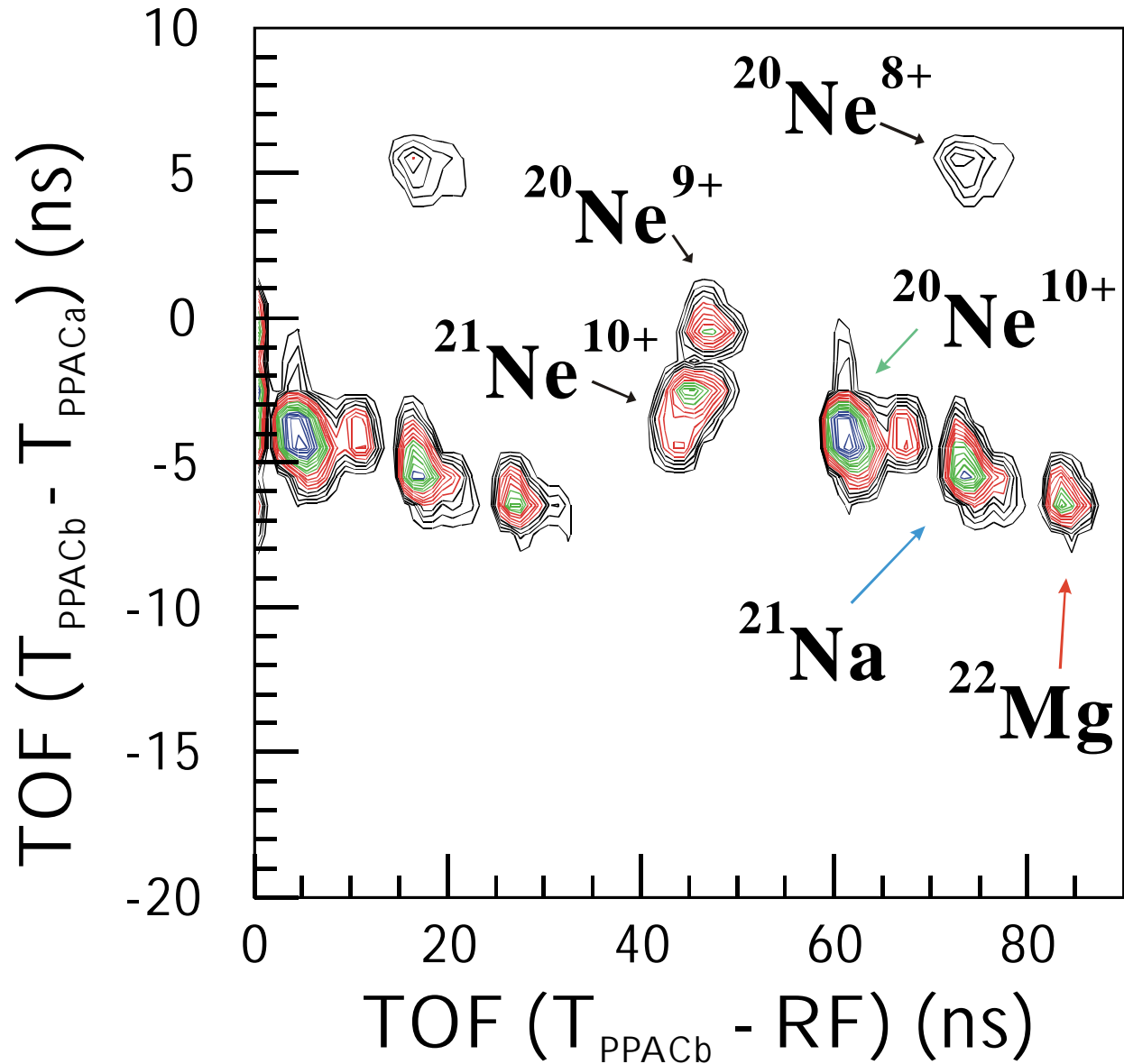
^{21}Na : 4.0 AMeV
24.3 kaps (12%)

Main Background: ^{20}Ne

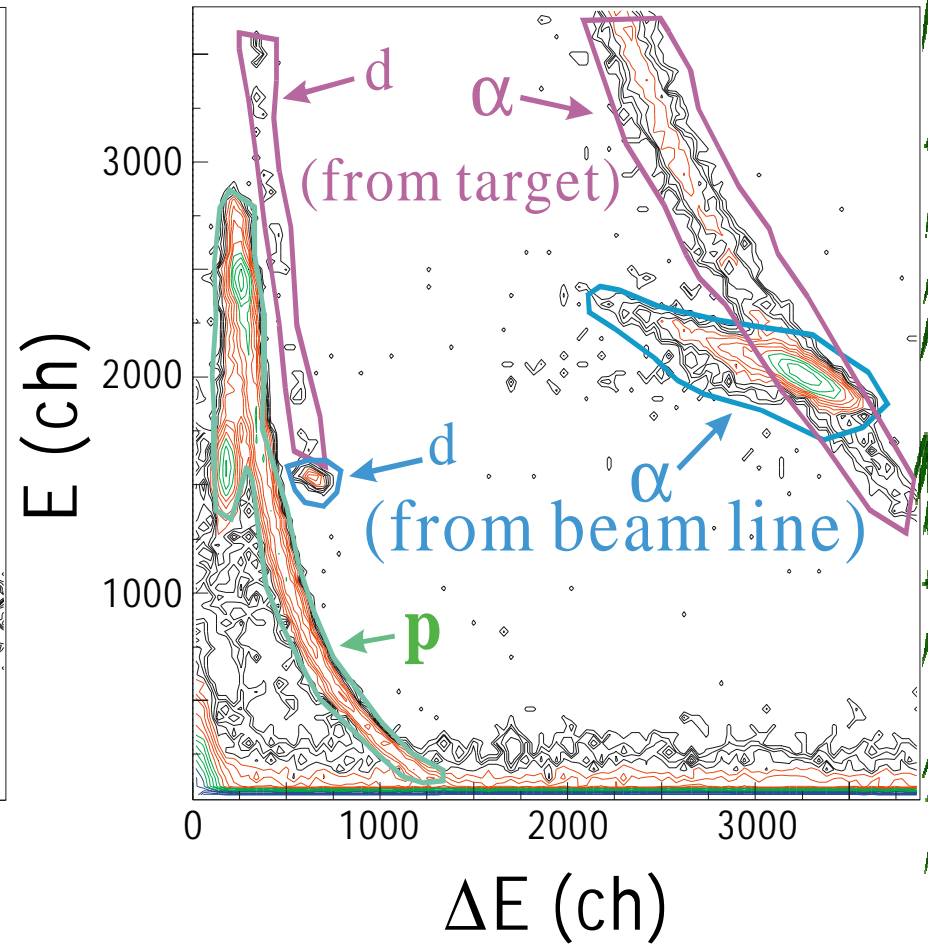
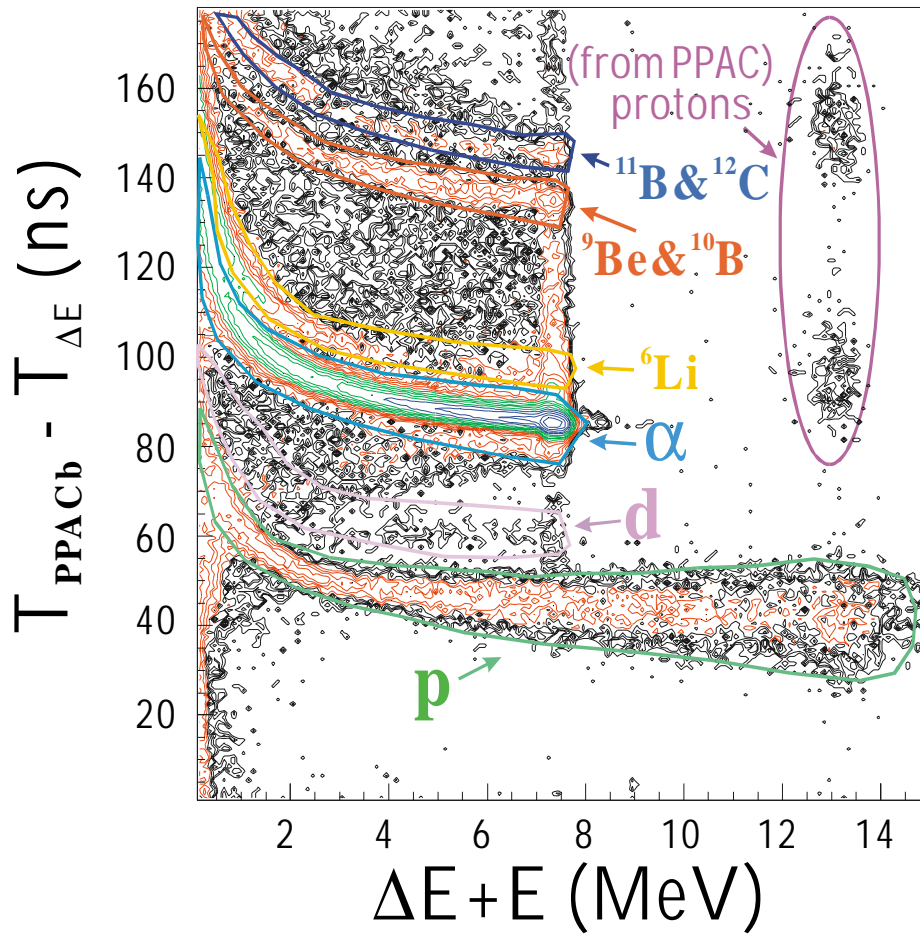
Target $(\text{CH}_2)_n$
 $\phi 3 \text{ cm}, 90 \mu\text{m}$
(C target)

$5 \times 5 \text{ cm}^2$
|--- ΔE (75 μm , PSD)
|--- E (1.5 cm, SSD)
|--- E_{Rej} (1.5 cm, SSD)

(4) Identification of RI Beam



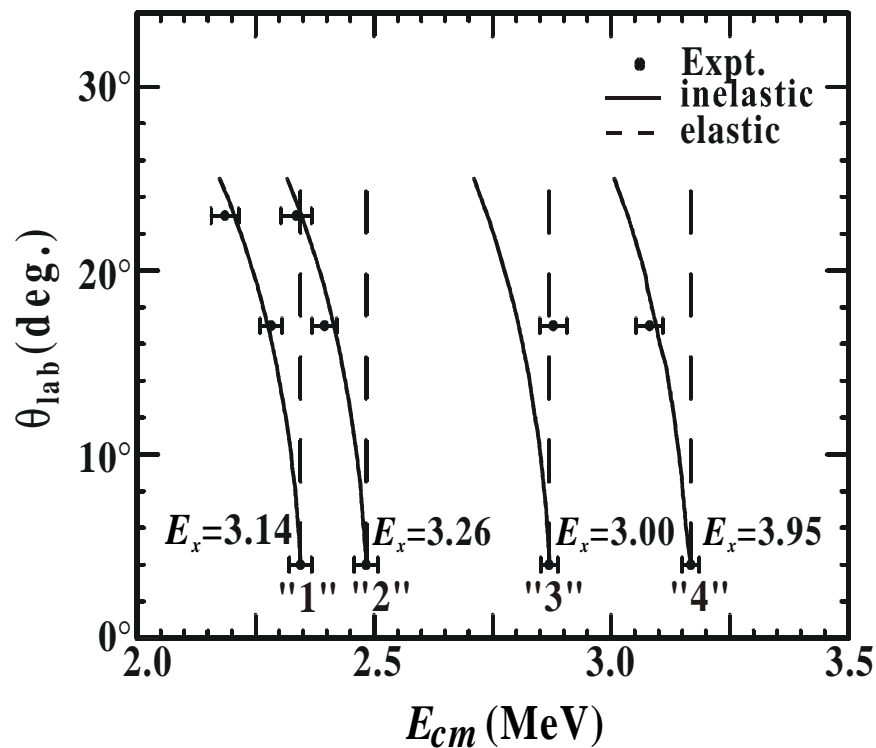
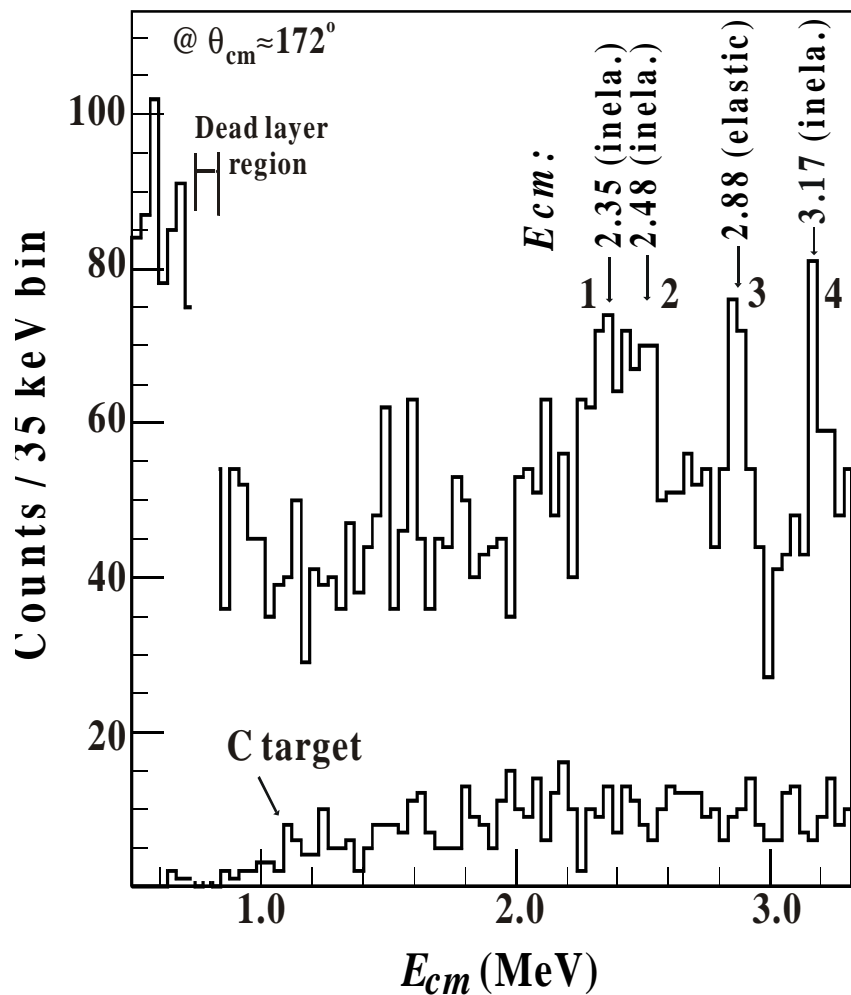
(5) Identification of Particles



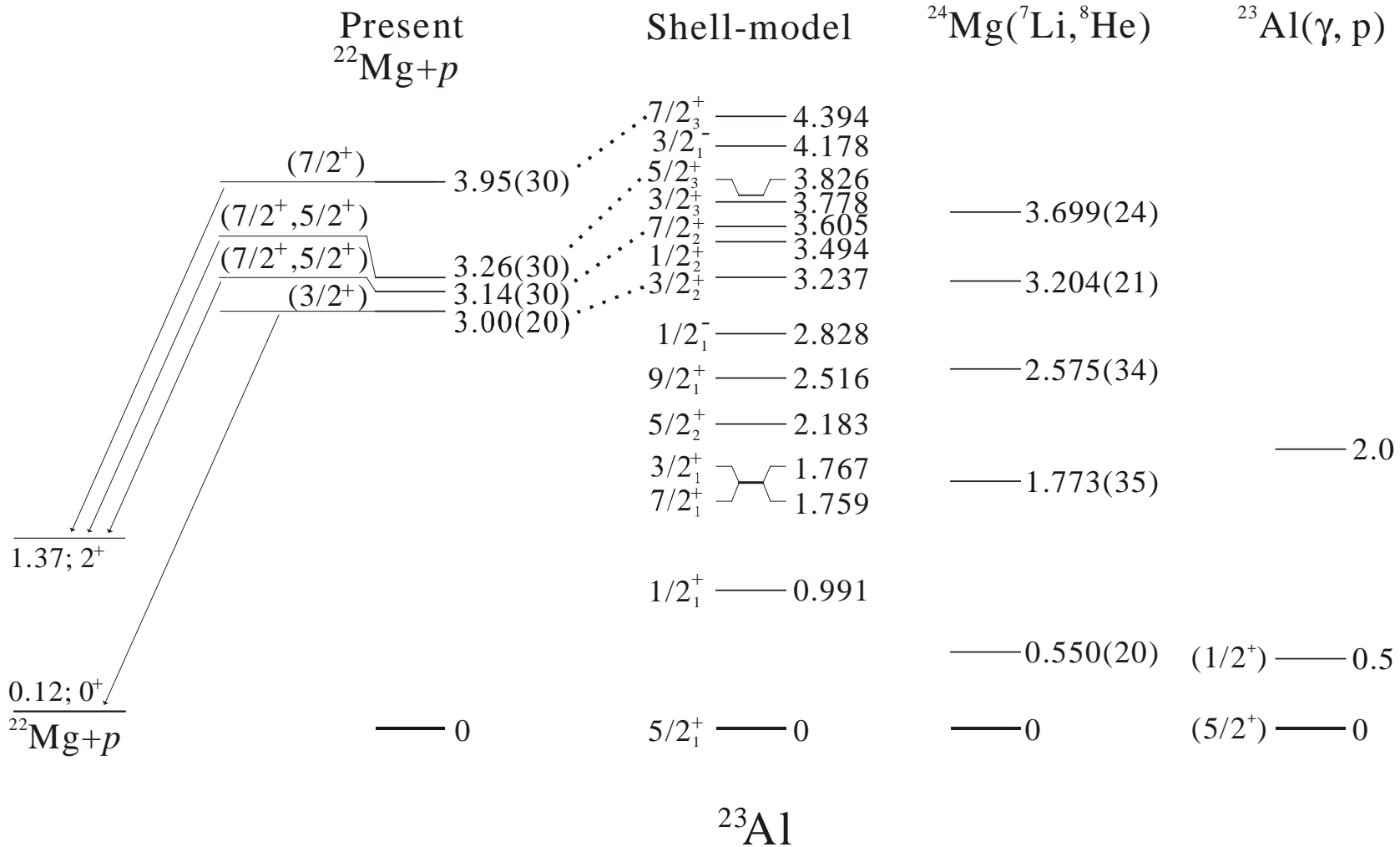
(Max energy for ΔE is 7.5 MeV)

Results

(1) Spectrum of $^{22}\text{Mg}+p$ Scattering



(2) Level Scheme in ^{23}Al



Model Calculations

(I) R-matrix code: SAMMY-M6-BETA

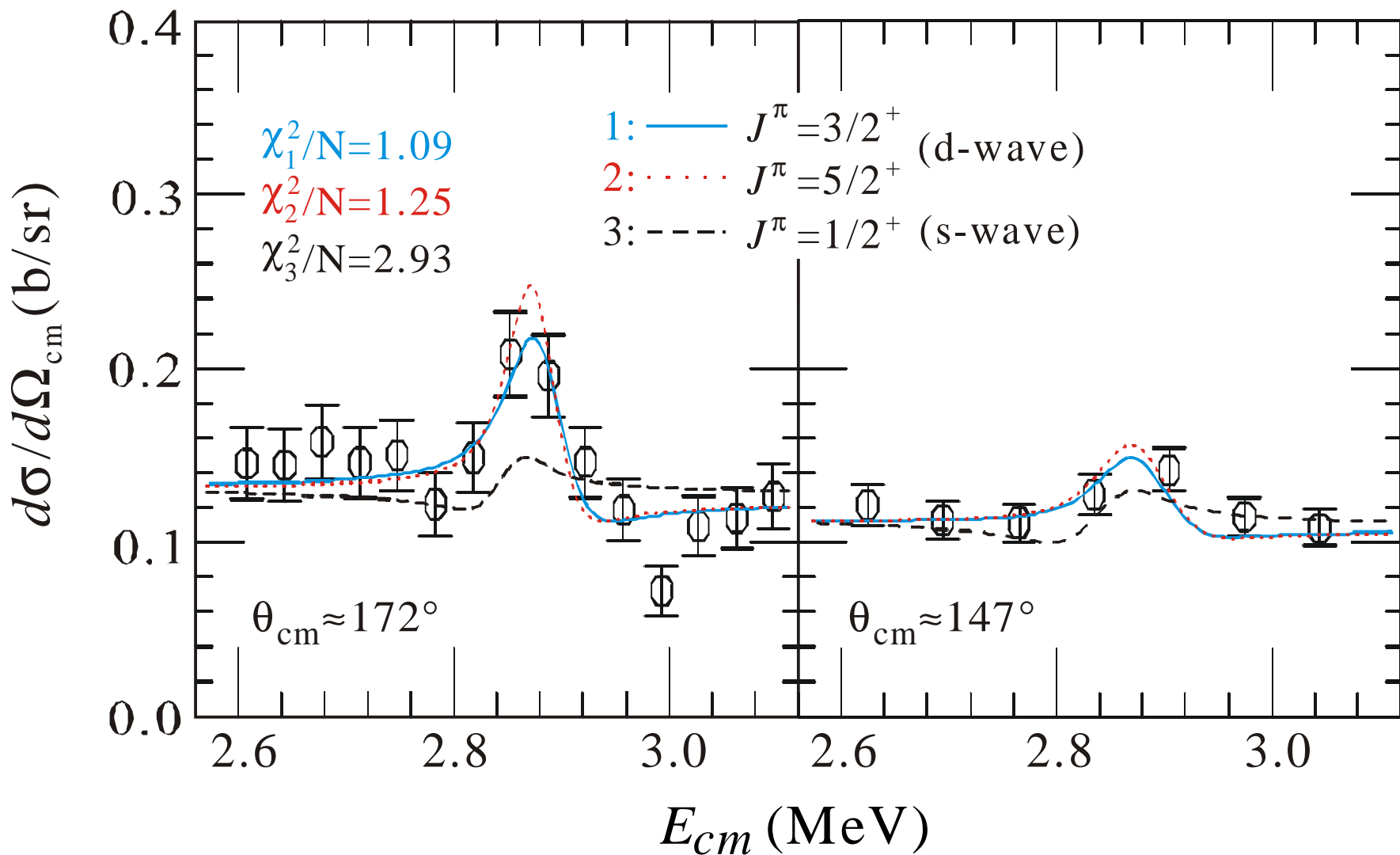
**A Code System for Multilevel R-Matrix Fits to
Neutron and Charged-Particle Cross-Section**

Data Using Bayes' Equations (By [Nancy M. Larson @ORNL](#))

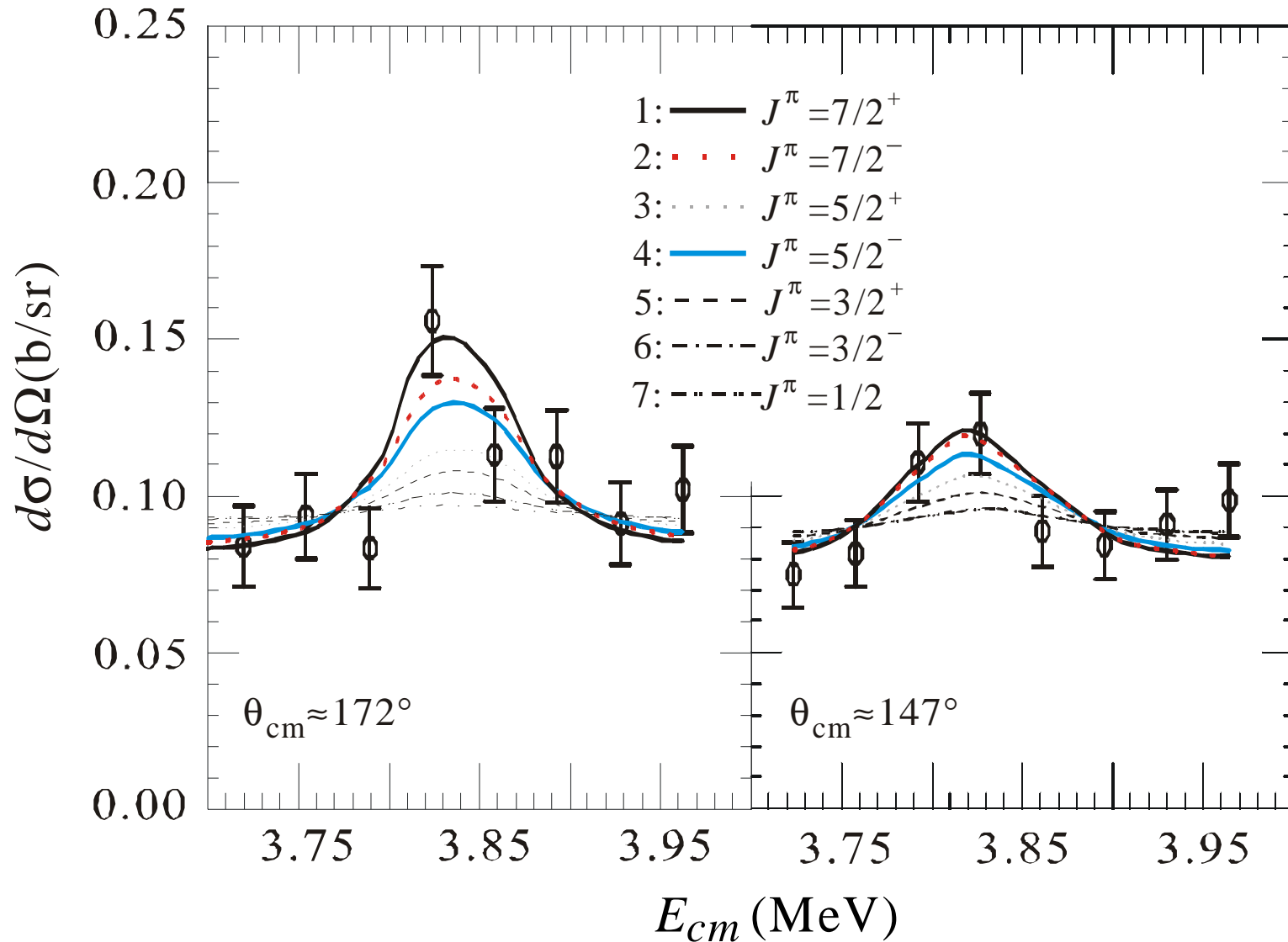
Data Type of Cross Sections

- 1. Elastic (both Angle-Integrated and Differential);**
- 2. Inelastic scattering;**
- 3. Angle-Differential Reaction Cross Sections;**
Total; Transmission; Fission (or Reaction); Capture;
Absorption, etc.

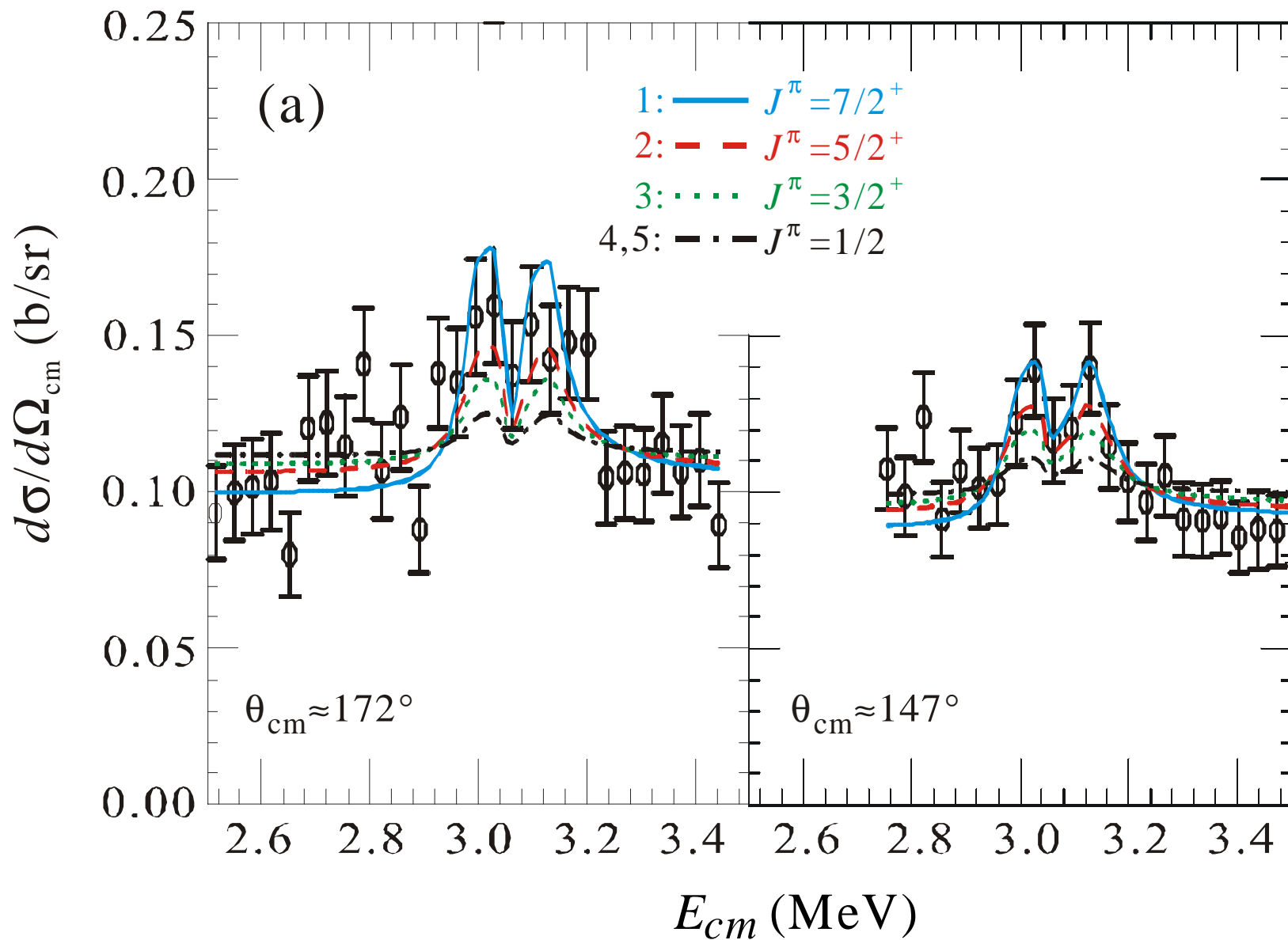
(I-1) *R*-matrix fits for 3.00-MeV state



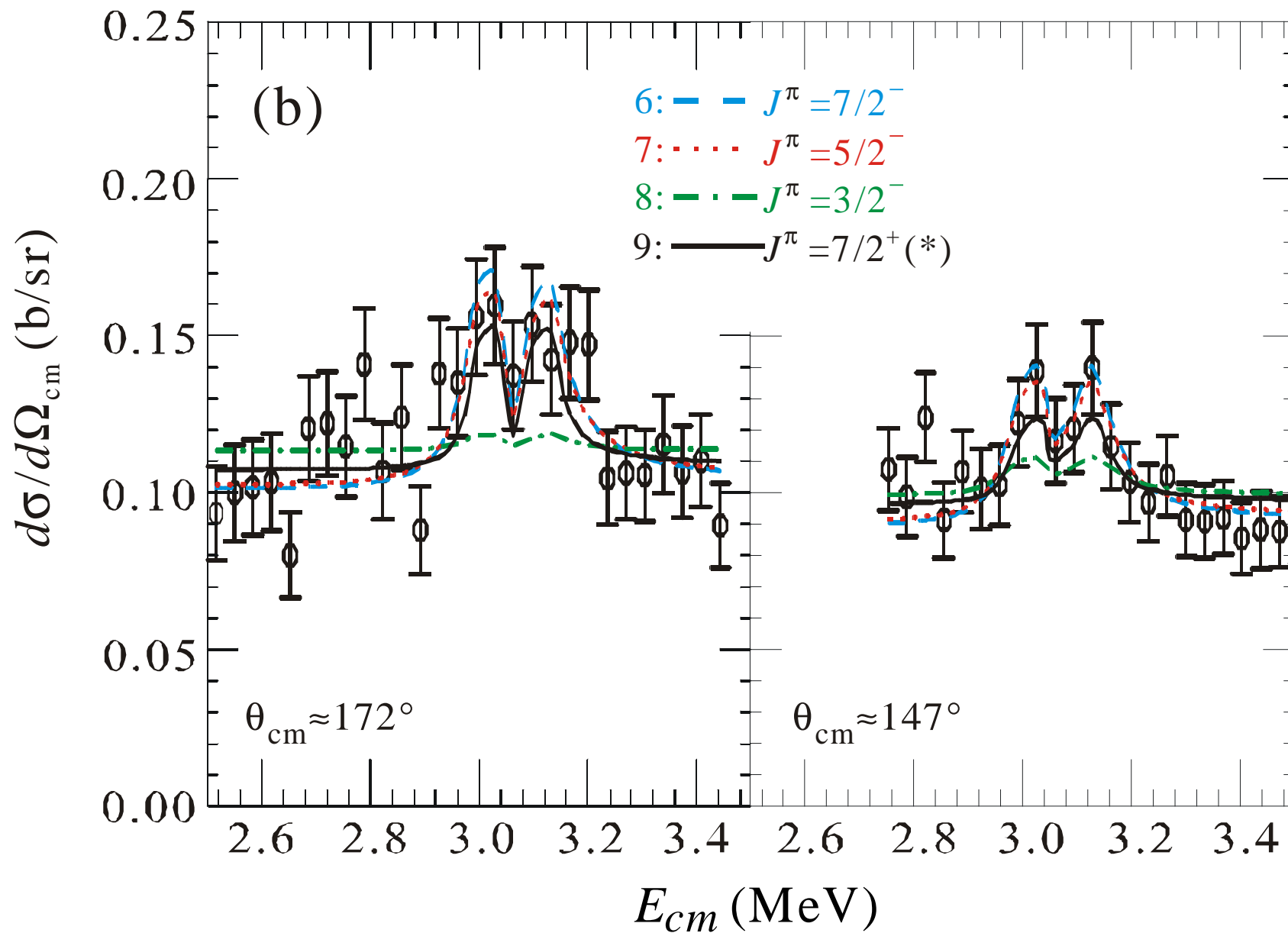
(I-2) *R*-matrix fits for 3.95-MeV state



(I-3a) Fits for 3.14, 3.26 MeV states



(I-3b) Fits for 3.14, 3.26 MeV states



(I-4) *R*-matrix fitting results

E_x (MeV)	J^π	Γ_ρ (keV)	$\Gamma_{\rho'}$ (keV)
3.00 (0.02)	(5/2 ⁺ , 3/2 ⁺)	17 (3), 32 (5)	
3.14 (0.03)	(7/2, 5/2)	$2 \leq \Gamma_\rho \leq 5$	30 (20)
3.26 (0.03)	(7/2, 5/2)	$2 \leq \Gamma_\rho \leq 5$	30 (20)
3.95 (0.03)	(7/2, 5/2 ⁻)	20 (10)	30 (20)

(II) Shell model code: OXBASH (B. A. Brown, MSU)

MODEL SPACE: full $(0 + 1) \hbar\omega$ (named as *spsdpf*)

INTERACTION: Isospin-conserving *WBT* interaction of
Warburton and Brown

ADVANTAGE: Positive $(0\hbar\omega)$ and negative $(1\hbar\omega)$ states
can be treated within the same procedure.

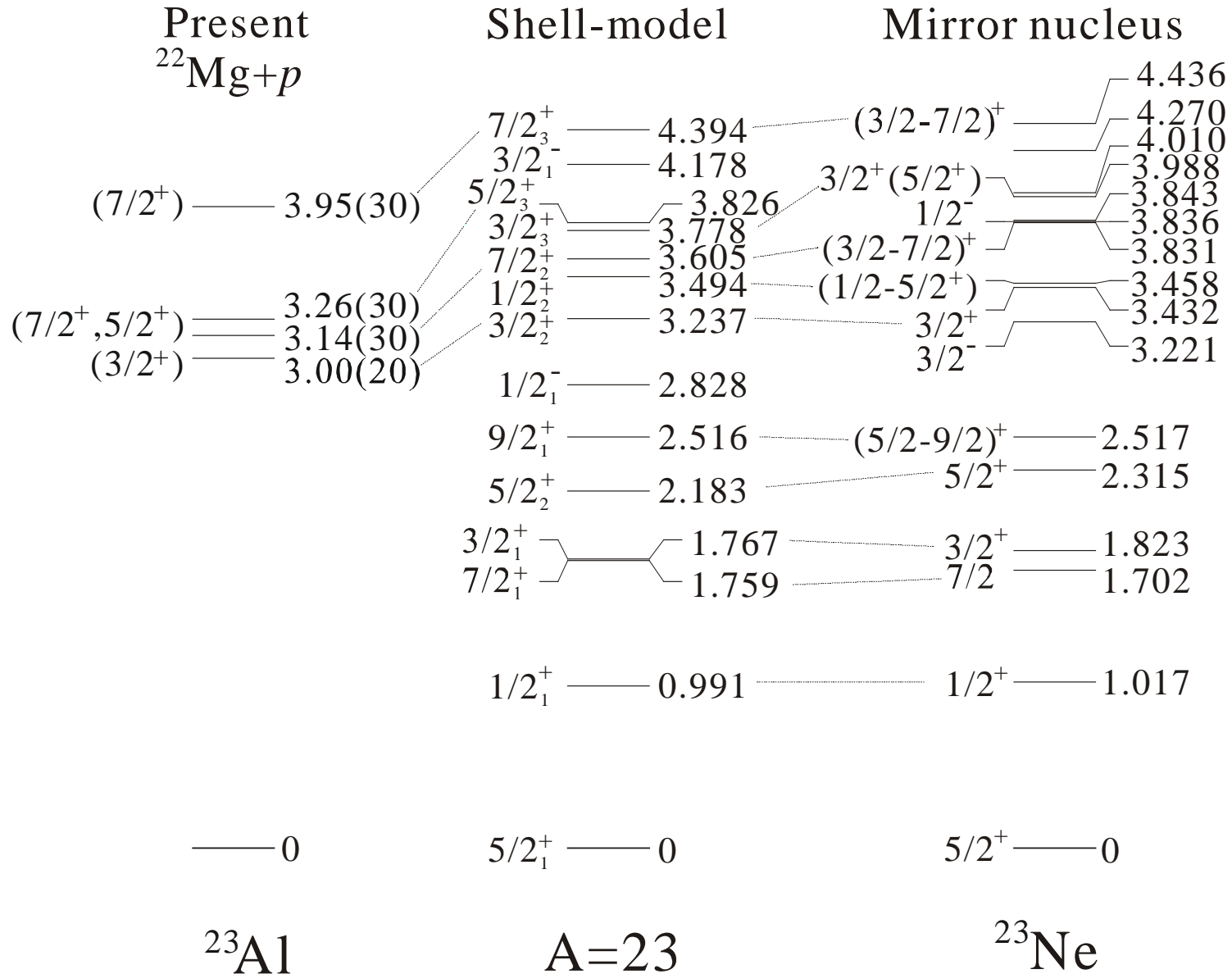
(II-1) Spectroscopic factor C^2S

E_x ^{23}Ne	E_x (cal.)	ΔE_x	J^π (expt.)	J^π (calc.)	l	C^2S	
						(d, p)	calc.
0	0	0	$5/2^+$	$5/2_1^+$	2	0.22	0.34
1.017	0.991	0.03	$1/2^+$	$1/2_1^+$	0	0.70	0.66
1.823	1.767	0.06	$3/2^+$	$3/2_1^+$	2	0.02	0.08
2.315	2.183	0.13	$5/2^+$	$5/2_2^+$	2	0.06	0.09
3.221	4.178	-0.96	$3/2^-$	$3/2_1^-$	1	0.30	0.28
3.432	3.237	0.20	$3/2^+$	$3/2_2^+$	2	0.30	0.28
3.836	2.828	1.01	$1/2^-$	$1/2_1^-$	1	0.11	0.01
3.988	3.778	0.21	$3/2^+ (5/2^+)$	$3/2_3^+$	2	0.30	0.22

(II-2) Spectroscopic factor C²S

E _x exp.	E _x calc.	J ^π calc.	to Ground State of ²² Mg (0 ⁺)			to Excited State of ²² Mg (2 ⁺)		
			1d _{5/2}	1d _{3/2}	2s _{1/2}	1d _{5/2}	1d _{3/2}	2s _{1/2}
	0.0	5/2 ₁ ⁺	0.335			0.82	0.005	0.002
0.55	0.991	1/2 ₁ ⁺			0.66	0.40	0.026	
1.77	1.767	3/2 ₁ ⁺		0.076		0.07	0.05	0.40
	2.183	5/2 ₂ ⁺	0.086			0.15	0.01	0.62
3.00	3.237	3/2 ₂ ⁺		0.284		0.005	0.06	0.02
3.14	3.605	7/2 ₂ ⁺				0.024	0.323	
	3.778	3/2 ₃ ⁺		0.22		0.005	0.11	0.17
3.26	3.826	5/2 ₃ ⁺	0.007			0.01	0.033	0.023
3.95	4.394	7/2 ₃ ⁺				0.002	0.18	

(II-3) States in ^{23}Al & ^{23}Ne



(II-3) Calc. of Partial Width Γ_p

By C. Iliadis [Nucl. Phys. **A618**, (1997) 166]

$$\Gamma_p = \frac{2\hbar^2}{\mu R^2} P_\ell C^2 S \theta_{sp}^2$$

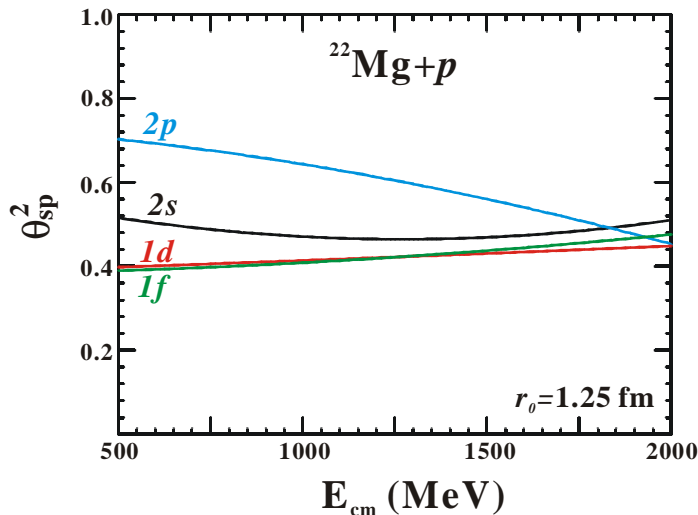
μ is the reduced mass in *amu*.

P_ℓ the Coulomb penetrability calculated with standard program.

C^2 is the isospin Clebsch-Gordan coefficient, S is the single-particle Spectroscopic factor. $C^2 S$ can be calculated within shell model.

θ_{sp}^2 dimensionless single-particle reduced width.

$R = r_0(A_t^{1/3} + A_p^{1/3})$ is the channel (or interaction) radius in *fm*.



For radius range of $r_0 = 1.2 \sim 1.5$ fm

Uncertainty of θ_{sp}^2 : 15 ~ 25 %

Uncertainty of P_ℓ : 40 ~ 60 %

Uncertainty of Γ_p : 2 ~ 8 %

Overall uncertainty of Γ_p : ≤ 15 %

Comparisons

E_x	J^π_{expt}	J^π_{calc}	ℓ	$\Gamma_{p,(p')}^{\text{exp}}$	$\Gamma_{p,(p')}^{\text{cal}}$
0.55	(1/2 ⁺)	1/2 ₁ ⁺	0		0.074
1.77	(3/2 ⁺)	3/2 ₁ ⁺	2		1.0
3.00	(3/2 ⁺ , 5/2 ⁺)	3/2 ₂ ⁺	2	32(5), 17(3)	44(7)
3.14*	(7/2 ⁺ , 5/2 ⁺)	7/2 ₂ ⁺	2	30(20)	6(1)
3.26*	(7/2 ⁺ , 5/2 ⁺)	5/2 ₃ ⁺	0	30(20)	10(2)
3.95*	(7/2 ⁺)	7/2 ₃ ⁺	2	30(20)	18(3)

* Their widths are for the decay branches to the 1st states in ²²Mg.

(Energy in MeV and width in keV)

Conclusion

(1) New observations:

$E_x = 3.00$ MeV, $J^\pi = (3/2^+)$ with $\Gamma_p = 28(5)$ keV (elastic)

$E_x = 3.95$ MeV, $J^\pi = (7/2^+)$ with $\Gamma_p = 30(15)$ keV (inelastic)

[2] $E_x = 3.14$ & 3.26 MeV states:

inelastic p-decay to the first excited state in ^{22}Mg

Both most likely with $J^\pi = (7/2^+, 5/2^+)$.

[3] Weak-coupling picture

E_x	3.00	3.14	3.26	3.95
Coupling	$0_1^+ \otimes d_{3/2}$	$2_1^+ \otimes d_{3/2}$	$2_1^+ \otimes s_{1/2}$ & $2_1^+ \otimes d_{3/2}$	$2_1^+ \otimes d_{3/2}$