

# Nuclear Structure of $^{44-46}\text{Ti}$ isotopes using shell model by OXBASH code

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## Abstract

Studying 44-46 Ti nuclei's nuclear composition in the shell model's framework was a key to understand the electronic transition at one orbital ( $1f_{7/2}$ ) and Even-even nucleus. Furthermore 40Ca shell is considered to be closed. The explanation nucleus's excited states of 44-46 Ti with neutrons ( $N=22,24$ ) with 4 to 6 nucleons outside of closed shell. The energy levels were calculated using the shell model and decrease in the probability of the electric quadruple transition  $B(E2)$  for the 44-46 Ti isotopes by utilizing the OXBASH code included within the f7 shell in conjunction with the F7MBZ& F742, The findings were analyzed and contrasted with certain previous experimental values. It was determined that the obtained theoretical results and the experimental data agree. There were a good agreement with theoretical and experimental results which paved the road for more applications. Ultimately, through this study, the best effective effort was reached, which is F7MBZ.

**Keywords:** *Energy levels; Shell model; OXBASH code,  $B(E2)$ .*

## 1. INTRODUCTION

The Shell model's calculations, implemented in a model space where the nucleons are constrained to occupy more than a few orbits without the need for scaling factors in order to replicate the recorded static moments or transition strengths are capable of reproducing observed static moments and transition strengths. Traditionally, Transition density models have undergone extensive testing by comparing computed and measured form factors of longitudinal electron scattering. Nuclear excitations have been studied using a variety of microscopic and macroscopic theories. [1]. The structure of neutron-rich nuclei has recently been the topic of substantial theoretical and experimental investigation. The current study is based on

the notion that nuclei with a large neutron excess can suffer major alterations to their fundamental shell structure [2]. With the nuclear shell model SM in mind, Talmi used SDI to investigate the surface delta relationship. to compute the properties of closed core nuclear scenarios involving few particles [3]. Computations in the well-known shell-model approach are carried out in a reduced Hilbert space, the so-called model space, and only particles outside a core made up of filled shells (valence particles) are considered active [4]. The purpose of this research is to look into the decreased transition probabilities and level schemes for 44Ti, 46Ti, and even isotopes, use the most recent OXBASH for Windows version the energy levels of some 44Ti states and 46Ti Compared

to the most recent data, the figures calculated in this paper show [2].

## 2. Theory

The force that created by collision of two nucleons is known as the residual interaction and the Hamiltonian operator perturbation that causes this interaction. The potential energy of two particles is equal to their sum then use the equation to symbolize the Hamiltonian operator in the state of perturbation [3].

$$H = H_0 + \sum_{i < j} V_{ij} \dots \dots \dots (1)$$

$H_0$ : is an unperturbed Hamiltonian,  $V_{ij}$ : remains of the two-body interaction [5]. In a D-dimensional Hilbert space [6], Schrödinger equation can be expressed in writing.

$$H |\psi_n\rangle = E_n |\psi_n\rangle \dots \dots \dots (2)$$

Where

$$H = H_0 + H_1 \dots \dots \dots (3)$$

$$H_0 = \sum_{i=1}^A (T_i + U_i) \dots \dots \dots (4)$$

$$H_0 = \sum_{i < j}^A (V_{ij}^{NN} - \sum_{i=1}^A U_i) \dots \dots \dots (5)$$

To disentangle [3] the nuclear Hamiltonian one-body potentials have been proposed as the integral of one-body terms. Which characterizes the nucleons' free-moving nature plus the interaction  $H_1$ . When solving the Schrodinger equation for a central potential, the energy of individual nucleons are given by the SPE. As shown in the states of a single particle, Whether it be a particle or a hole outside of a nucleus with a doubly-closed shell (DCS) in its neighbors (DCS  $\pm 1$ ) [7]. Lastly, by choosing the desired interaction the formation of the Hamiltonian, and after that, the calculations are carried out [8].

## 3. Calculations and Discussion

Titanium isotope structure is examined using the shell model employing the OXBASH system. For the calculations, different discussions are possible for the level of energy decreased likelihood of E2 transitions. The  $^{44-46}\text{Ti}$  isotopes have 22 and 22 (or 24) proton, and neutron respectively and The core is considered as  $^{40}\text{Ca}$  for  $^{44}\text{Ti}$  with 4 nucleons outside core and for  $^{46}\text{Ti}$  with 6 nucleons outside core.

The OXBASH code has been implemented by us in m-scheme as well as jj-coupling respectively. This study seeks consists of determining energy levels and lowered electric quadrupole transition probability B (E2).

### 3.1. Energy levels

The purpose of this investigation is to Determine the nuclei that are in close proximity to  $^{44}\text{Ti}$  because of the significant role that these nuclei play in recent developments in astrophysical applications. The determined energy levels and presented low-lying state experimental results for even-even nuclei. However, On the left, you can see the results of our calculations and right-hand experimental info for any band [9].

#### 3.1.1 Energy levels of $^{44}\text{Ti}$

For  $^{44}\text{Ti}$  isotope using (F7MBZ) interactions is shown in the table 1. When measured against the experimental data that is presented. Both of the parity and angular momentum are found to be identical to the ground state of level  $0^+$ . A good agreement was obtained for the values of the practical energies (1.08306, 2.45433, 4.01531, 7.670, 9.100) MeV corresponding to the angular momentum ( $2^+$ ,  $4^+$ ,  $6^+$ ,  $1,6^+$ ,  $2,4^+$ ) when compared with the calculated theoretical values. The total momentum and parity of the unconfirmed practical energies (6.5085, 6.810, 7.6711, 8.0399, 9.361, 9.713, 10.460) MeV was confirmed for the angular

momentum (  $8+1$  ,  $0+2$ ,  $10+1$  , $12+1$  , $2+3$  , $4+4$ ,  $0+3$ ) are confirmed when compared with the calculated theoretical values , And The experimental energy value( $5.1511$  , $10.113$  ) MeV was confirmed for angular momentum ( $6-2$  , $3-2$ ) with positive parity In our calculations , We expect that practical values ( $6.959$  ,  $6.5718$  ,  $8.984$  ) MeV whose angular momentum is uncertain (  $4+$  ,  $8+$ ,  $10+$ ) have angular momentum (  $3+1$  ,  $7+1$ ,  $9+1$ ) , We found the values of energies for a specific angular momentum close to the values of the practical energies ( $7.634$  , $8.067$  ,  $8.318$  , $9.030$  , $9.280$  , $9.542$  , $9.668$  ,  $9.895$  , $10.520$ ) that have no specific angular momentum, and thus we expect that its angular momentum is the theoretically calculated momentum ( $4+3$  ,  $8+2$ ,  $5+2$  ,  $6+4$ ,  $8+3$ ,  $2+4$ ,  $7+2$  ,  $10+2$ ,  $6+5$ ). Through our calculations, we noticed that there are three levels with total angular momentum and parity that were not matched by any available practical value. For  $^{44}\text{Ti}$  isotope using ( F742) interactions is shown in the table2 . When measured against the experimental data that is presented , The parity and angular momentum are determined to be similar to the level  $0+$  ground state.. A good agreement was obtained for the values of the practical energies ( $1.08306$ ,  $2.45433$  ,  $4.01531$

, $7.458$  , $7.670$ ) MeV corresponding to the angular momentum (  $2+1$  ,  $4+1$ ,  $6+1$  , $8+2$  , $6+4$ ) when compared with the calculated theoretical values , confirmed The total momentum and parity of the unconfirmed practical energies (  $4.8031$ ,  $6.5718$  ,  $6.959$  ,  $6.84884$  , $7.6711$  , $8.0399$  ,  $8.180$  , $8.947$  , $9.143$  , $8.984$  ) MeV corresponding to the angular momentum (  $6+2$ ,  $8+1$ ,  $4+3$  , $6+3$  , $10+1$  ,  $12+1$ ,  $2+3$  ,  $4+5$  ,  $0+3$  ,  $10+2$ ) when compared with the calculated theoretical values , And The experimental energy value( $5.6706$  , $8.534$  ) MeV was confirmed for angular momentum ( $7-1$  , $3-2$ ) with positive parity In our calculations , A good agreement was obtained for the practical energy values ( $5.055$  , $5.250$  ) corresponding to the angular momentum ( $3-1$ , $5-1$  ) but with negative parity , We found the values of energies for a specific angular momentum close to the values of the practical energies ( $4.7922$  , $8.067$  ,  $8.318$  , $9.180$ ) that have no specific angular momentum, and thus we expect that its angular momentum is the theoretically calculated momentum ( $2+2$  ,  $4+4$ ,  $7+2$  ,  $6+5$ ). Through our calculations, we noticed that there are six levels with total angular momentum and parity that were not matched by any available practical value.

**Table (1): Excitation energy predicted by (F7MBZ) interactions and observed excitation energies[10] for the  $^{44}\text{Ti}$  nucleus are compared**

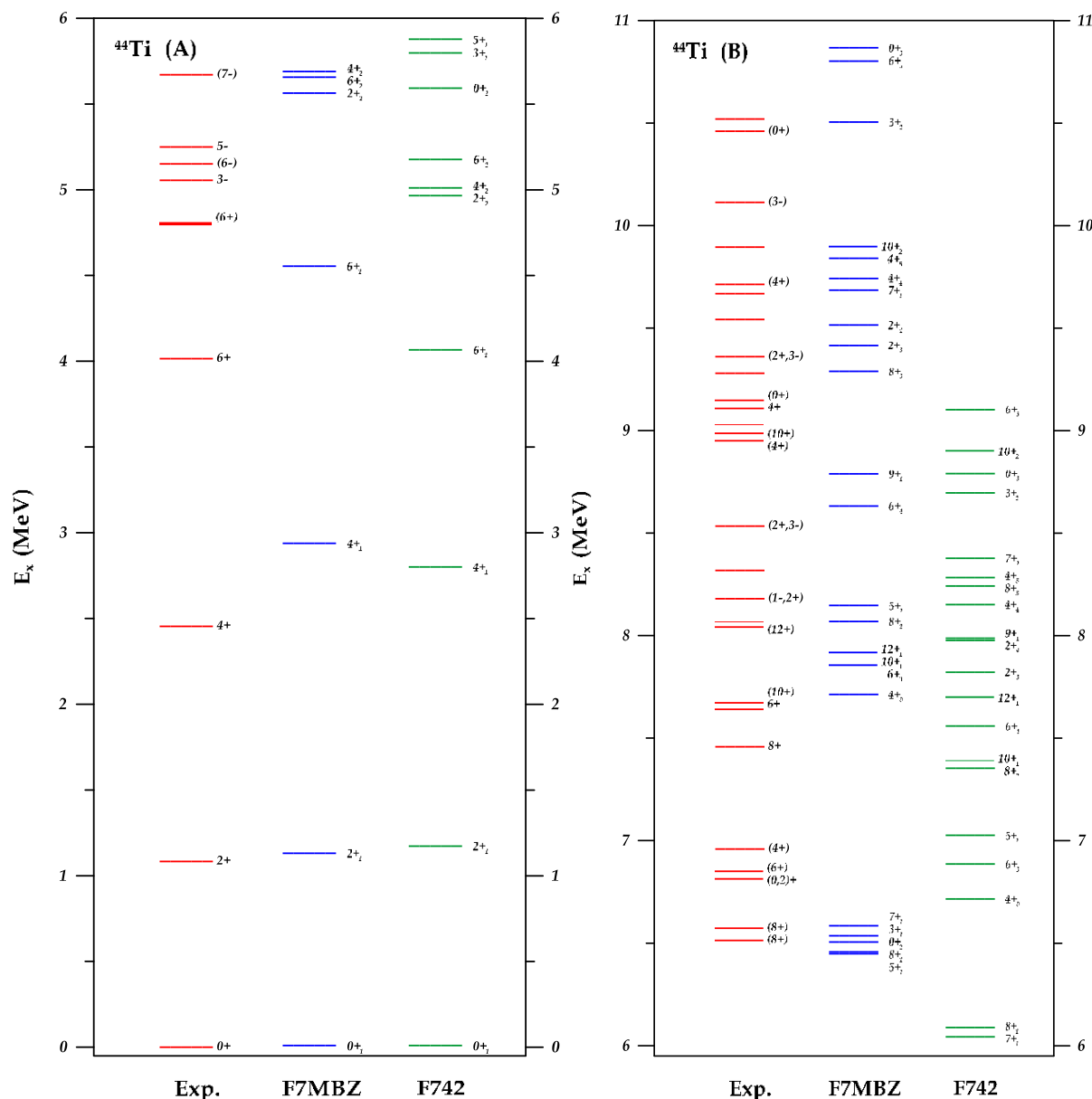
F7MBZ (Theoretical values)		Experimental values		Theoretical values for F7MBZ		Experimental values	
$J^{\pi+}$	E (MeV)	E (MeV)	$J^{\pi}$	$J^{\pi+}$	E (MeV)	E (MeV)	$J^{\pi}$
$0_1$	0	0	$0+$	$12_1$	7.905	8.0399	( $12+$ )
$2_1$	1.121	1.08306	$2+$	$8_2$	8.061	8.067	-----
$4_1$	2.929	2.45433	$4+$	$5_2$	8.139	8.318	-----
$6_1$	4.545	4.01531	$6+$	$6_4$	8.624	9.030	-----
$2_2$	5.554	4.7922	-----	$9_1$	8.781	8.984	( $10+$ )
$6_2$	5.647	5.1511	( $6-$ )	$8_3$	9.281	9.280	-----
$4_2$	5.680	-----	-----	$2_3$	9.407	9.361	( $2+,3-$ )
$5_1$	6.440	-----	-----	$2_4$	9.507	9.542	-----
$8_1$	6.449	6.5085	( $8+$ )	$7_2$	9.677	9.668	-----
$0_2$	6.497	6.810	( $0,2$ ) $+$	$4_4$	9.734	9.713	( $4+$ )
$3_1$	6.528	6.959	( $4+$ )	$4_5$	9.832	9.100	$4+$
$7_1$	6.577	6.5718	( $8+$ )	$10_2$	9.891	9.895	-----
$4_3$	7.705	7.634	-----	$3_2$	10.497	10.113	( $3-$ )

6 <sub>3</sub>	7.842	7.670	6+	6 <sub>5</sub>	10.794	10.520	-----
10 <sub>1</sub>	7.845	7.6711	(10+)	0 <sub>3</sub>	10.859	10.460	(0+)

**Table (2): Excitation energy predicted by (F742) interactions and observed excitation energies[10] for the  $^{44}\text{Ti}$  nucleus are compared.**

F742(Theoretical values)		Experimental values		F742(Theoretical values)		Experimental values	
$J^{\pi+}$	E (MeV)	E (MeV)	$J^{\pi}$	$J^{\pi+}$	E (MeV)	E (MeV)	$J^{\pi}$
0 <sub>1</sub>	0	0	0+	8 <sub>2</sub>	7.345	7.458	8+
2 <sub>1</sub>	1.163	1.08306	2+	10 <sub>1</sub>	7.380	7.6711	(10+)
4 <sub>1</sub>	2.791	2.45433	4+	6 <sub>4</sub>	7.551	7.670	6+
6 <sub>1</sub>	4.057	4.01531	6+	12 <sub>1</sub>	7.689	8.0399	(12+)
2 <sub>2</sub>	4.956	4.7922	-----	2 <sub>3</sub>	7.813	8.180	(1-,2+)
4 <sub>2</sub>	5.001	-----	-----	2 <sub>4</sub>	7.968	-----	-----
6 <sub>2</sub>	5.167	4.8031	(6+)	9 <sub>1</sub>	7.979	--	-----
0 <sub>2</sub>	5.583	-----	-----	4 <sub>4</sub>	8.143	8.067	-----
3 <sub>1</sub>	5.788	5.055	3-	8 <sub>3</sub>	8.234	-----	-----
5 <sub>1</sub>	5.868	5.250	5-	4 <sub>5</sub>	8.275	8.947	(4+)
7 <sub>1</sub>	6.035	5.6706	(7-)	7 <sub>2</sub>	8.369	8.318	-----
8 <sub>1</sub>	6.080	6.5718	(8+)	3 <sub>2</sub>	8.688	8.534	(2+,3-)
4 <sub>3</sub>	6.707	6.959	(4+)	0 <sub>3</sub>	8.782	9.143	(0+)
6 <sub>3</sub>	6.878	6.84884	(6+)	10 <sub>2</sub>	8.891	8.984	(10+)
5 <sub>2</sub>	7.018	-----	-----	6 <sub>5</sub>	9.094	9.180	-----

**Fig.1. A comparison between theoretical energy levels for two interaction and the experimental data for  $^{44}\text{Ti}$ .**



when compared with the calculated theoretical values , And The experimental energy value(5.409,5.530 ,6.8303,7.019 ) MeV was confirmed for angular momentum (3-2 , 3-3,9-1 , 3-4 ) with positive parity In our calculations , We expect that practical values (9.168, 9.205) MeV whose angular momentum is uncertain ( 4+ , 6+) have angular momentum ( 3+8 , 7+7) , We found the values of energies for a specific angular momentum close to the values of the practical energies (3.213 , 3.7715 , 4.003 , 5.117 ,5.154 , 5.811 ,6.025 ,6.021, 6.251 , 6.266 , 6.305 ,6.574 ,6.616 , 6.794 , 6.851 ,6.974 , 7.101 ,7.147 , 7.172 ,7.201 , 7.288 ,7.350 , 7.472 , 7.558 , 7.584 , 7.608 ,7.735 ,7.788 ,7.849 ,7.874 , 7.937 ,8.013 ,8.040 , 8.088 ,8.134 ,8.230 ,8.2939 ,8.384,8.467 ,8.530 ,8.621, 8.701 ,8.761 ,8.808 ,8.940 ,8.984 ,9.070, 9.141 ,9.253 , 9.345 , 9.572 ,9.649 ,9.761 ,9.790,

9.852 , 9.864 ,10.038 , 10.212 ,10.441 , 10.602 ,10.730 ,10.866 , 10.938 , 10.980 , 10.051 ,11.299 , 11.426 , 12.974 ) that have no specific angular momentum and thus we expect that its angular momentum is the theoretically calculated momentum (4+2 , 3+1, 5+1 , 0+2, 5+2, , 7+1, 7+2, 4+6 , 5+ 3, 6+4, 6+5 , 7+3, 5+4, 1+2, 6+6 , 8+3, 0+ 3, 2+6, 4+7,6+7 , 4+8, 5+5, 10+2, 8+4 , 4+9, 7+4, 8+5 , 5+6, 9+2 , 6+8, 2+7, 7+5, 9+3, 3+6, 4+10,6+9, 5+7, 3+7, 8+6, 2+8, 2+9 , 5+8, 8+7 , 7+6, 6+10, 5+9, 9+4 , 5+10 ,0+4 , 8+8, 8+9, 9+5, 7+8 , 11+2, 3+9 , 10+5, 10+6, 7+9, 8+10 , 12+2, 13+1,3+10 , 14+1, 9+6, 7+10, 11+3, 0+5 ,10+7). Through our calculations, we noticed that there are three levels with total angular momentum and parity that were not matched by any available practical value.

**Table (3): Excitation energy predicted by (F7MBZ) interactions and observed excitation energies[10] for the  $^{46}\text{Ti}$  nucleus are compared.**

F7MBZ (Theoretical values)		Experimental values		Theoretical values F7MBZ		Experimental values	
$J^{\pi+}$	E (MeV)	E (MeV)	$J^{\pi}$	$J^{\pi+}$	E (MeV)	E (MeV)	$J^{\pi}$
0 <sub>1</sub>	0	0	0+	9 <sub>2</sub>	7.863	7.849	-----
2 <sub>1</sub>	1.101	0.889286	2+	6 <sub>8</sub>	7.884	7.874	-----
4 <sub>1</sub>	2.174	2.009846	4+	2 <sub>7</sub>	7.932	7.937	-----
2 <sub>2</sub>	2.769	2.9618	2+	7 <sub>5</sub>	8.028	8.013	-----
4 <sub>2</sub>	3.218	3.213	-----	9 <sub>3</sub>	8.051	8.040	-----
6 <sub>1</sub>	3.267	3.29886	6+	3 <sub>6</sub>	8.064	8.088	-----
2 <sub>3</sub>	3.690	3.2357	2+	4 <sub>10</sub>	8.190	8.134	-----
3 <sub>1</sub>	3.747	3.7715	-----	6 <sub>9</sub>	8.226	8.230	-----
4 <sub>3</sub>	3.901	3.849	(4+)	10 <sub>3</sub>	8.228	8.2839	10+,11+,12+
1 <sub>1</sub>	3.996	3.731	1+	5 <sub>7</sub>	8.277	8.293	-----
5 <sub>1</sub>	4.022	4.003	-----	12 <sub>1</sub>	8.292	8.2175	12+
6 <sub>2</sub>	4.190	4.398	(5-,6+)	3 <sub>7</sub>	8.431	8.384	-----
6 <sub>3</sub>	4.888	4.527	(6+)	8 <sub>6</sub>	8.452	8.467	-----
8 <sub>1</sub>	4.986	4.8969	8+	2 <sub>8</sub>	8.499	8.530	-----
0 <sub>2</sub>	5.193	5.117	-----	2 <sub>9</sub>	8.674	8.621	-----
5 <sub>2</sub>	5.202	5.154	-----	5 <sub>8</sub>	8.721	8.701	-----
4 <sub>4</sub>	5.208	5.794	4+	8 <sub>7</sub>	8.778	8.761	-----
3 <sub>2</sub>	5.470	5.409	3-	7 <sub>6</sub>	8.915	8.808	-----
2 <sub>4</sub>	5.493	5.604	(2+)	6 <sub>10</sub>	8.939	8.940	-----
4 <sub>5</sub>	5.633	5.992	(4+)	5 <sub>9</sub>	8.953	8.984	-----
3 <sub>3</sub>	5.672	5.530	3-	9 <sub>4</sub>	9.067	9.070	-----
7 <sub>1</sub>	5.700	5.811	-----	5 <sub>10</sub>	9.156	9.141	-----
2 <sub>5</sub>	5.794	5.872	(2+)	10 <sub>4</sub>	9.173	-----	-----
8 <sub>2</sub>	5.938	6.2004	8+	3 <sub>8</sub>	9.226	9.168	4+

7 <sub>2</sub>	5.945	6.025	-----	7 <sub>7</sub>	9.235	9.205	6+
4 <sub>6</sub>	6.094	6.021	-----	0 <sub>4</sub>	9.241	9.253	-----
5 <sub>3</sub>	6.132	6.251	-----	1 <sub>3</sub>	9.319	9.170	1+
6 <sub>4</sub>	6.266	6.266	-----	8 <sub>8</sub>	9.348	9.345	-----
10 <sub>1</sub>	6.438	6.2419	10+	8 <sub>9</sub>	9.534	9.572	-----
6 <sub>5</sub>	6.443	6.305	-----	9 <sub>5</sub>	9.644	9.649	-----
7 <sub>3</sub>	6.588	6.674	-----	2 <sub>10</sub>	9.740	9.615	2+
5 <sub>4</sub>	6.657	6.616	-----	7 <sub>9</sub>	9.747	9.761	-----
1 <sub>2</sub>	6.860	6.794	-----	11 <sub>2</sub>	9.822	9.790	-----
6 <sub>6</sub>	6.912	6.851	-----	3 <sub>9</sub>	9.866	9.852	-----
9 <sub>1</sub>	6.921	6.8303	9-	10 <sub>5</sub>	9.883	9.864	-----
8 <sub>3</sub>	6.983	6.974	-----	10 <sub>6</sub>	10.140	10.038	-----
3 <sub>4</sub>	7.051	7.019	(3-,4+)	7 <sub>9</sub>	10.208	10.212	-----
0 <sub>3</sub>	7.111	7.101	-----	1 <sub>4</sub>	10.327	10.180	1+
2 <sub>6</sub>	7.117	7.147	-----	8 <sub>10</sub>	10.445	10.441	-----
4 <sub>7</sub>	7.186	7.172	-----	12 <sub>2</sub>	10.638	10.602	-----
6 <sub>7</sub>	7.255	7.201	-----	13 <sub>1</sub>	10.749	10.730	-----
4 <sub>8</sub>	7.406	7.288	-----	3 <sub>10</sub>	10.840	10.866	-----
5 <sub>5</sub>	7.433	7.350	-----	14 <sub>1</sub>	10.900	10.938	-----
10 <sub>2</sub>	7.459	7.472	-----	9 <sub>6</sub>	10.912	10.980	-----
3 <sub>5</sub>	7.576	7.534	(3-)	7 <sub>10</sub>	10.954	11.051	-----
8 <sub>4</sub>	7.593	7.558	-----	11 <sub>3</sub>	11.271	11.299	-----
4 <sub>9</sub>	7.594	7.584	-----	0 <sub>5</sub>	11.438	11.426	-----
7 <sub>4</sub>	7.619	7.608	-----	12 <sub>3</sub>	11.841	-----	-----
8 <sub>5</sub>	7.771	7.735	-----	9 <sub>7</sub>	12.200	-----	-----
11 <sub>1</sub>	7.774	7.9418	11+	10 <sub>7</sub>	12.351	12.974	-----
5 <sub>6</sub>	7.804	7.788	-----				

For <sup>46</sup>Ti isotope using ( F742) interactions is shown in the table4 . When measured against the experimental data that is presented, Both the parity and angular momentum are found to be identical to the ground state of level 0+. A good agreement was obtained for the values of the practical energies (0.889286, 2.009846 , 2.9618, 3.29886,3.2357, 3.731,4.675 ,4.8969 ,4.950 ,6.2004,6.118 ,6.338 ,6.2419,6.395,6.685 ,7.9418 , 8.2175) MeV corresponding to the angular momentum ( 2+1 , 4+1,2+2 , 6+1 , 2+3,1+1, 0+2,8+1 ,2+4 , 8+2 , 2+6, 4+7, 10+1 ,4+8 , 4+9,11+1,12+1) , when compared with the calculated theoretical values , confirmed The total momentum and parity of the unconfirmed practical energies (4.527, 5.079, 6.361, 5.965, 7.019 ,8.020 ) corresponding to the angular momentum (6+3, 4+5 , 6+4, 6+5, 4+10, 0+4) when compared with the calculated theoretical values , And The experimental energy

value(3.82643 ,6.958 ) MeV was confirmed for angular momentum (5-1 ,3-6) with positive parity In our calculations , A good agreement was obtained for the practical energy values (3.44139 ,3.826, 5.1976) corresponding to the angular momentum (4-3,5-1,7-4) but with negative parity, We found the values of energies for a specific angular momentum close to the values of the practical energies (3.213 , 3.551 , 3.7715 , 4.617, 4.845 ,5.000 , 5.094 , 5.117 , 5.154 , 5.180 , 5.811 , 5.903 , 6.025 , 6.021 , 6.251 , 6.266 , 6.305 ,6.424, 6.513 , 6.574 , 6.616, 6.794 , 6.851 , 6.974 , 6.041 , 7.101 ,7.147 , 7.172 ,7.201 , 7.238 ,7.288 , 7.350 , 7.472 ,7.558 , 7.660 , 7.710 , 7.730 , 7.735 , 7.788 , 7.849 ,7.874 ,7.917, 7.979 , 8.013 , 8.182 , 8.230 ,8.346 , 8.384 , 8.467 , 8.530 ,8.574 ,8.621 , 8.662 , 8.701 ,8.984 , 9.070,9.111 ,9.141 ,9.253 , 9.304 , 9.345 , 9.474 , 9.761 , 9.864 , 10.038 , 10.212 , 10.441 , 10.602 , 10.866 , 10.980 )

that have no specific angular momentum and thus we expect that its angular momentum is the theoretically calculated momentum (4+2 , 3+1 , 6+2 , 4+4 , 5+2 , 3+2 , 3+3 , 2+5 , 4+6 , 7+1 , 5+3 , 5+4 , 1+2 , 7+3 , 3+4 , 0+3 , 6+6 , , 8+3 , 5+5 , 6+7 , 3+5 , 2+7 , 5+6 , 8+4 , 7+4 , 6+8,10+2 , 8+5 , 5+7 , 6+9 , 7+5 , 3+7 , 9+3 , 2+9 , 5+8 , 8+6 , 10+3 , 8+7 ,1+3 , 5+9 , 3+8 ,

6+10 , 7+6 , 5+10 , 7+7 , 2+10 , 9+4 , 3+9 , 10+4 , 8+8 , 7+8 ,8+9 , 1+4 , 9+5 , 7+9 , 3+10,10+5 , 11+2 ,8+10 , 10+6 , 0+5 , 7+10 , 9+6 , 122 , 13+1 , 11+3 , 14+1 , 9+7 , 12+3 , 10+7 ). Through our calculations, we noticed that there are three levels with total angular momentum and parity that were not matched by any available practical value.

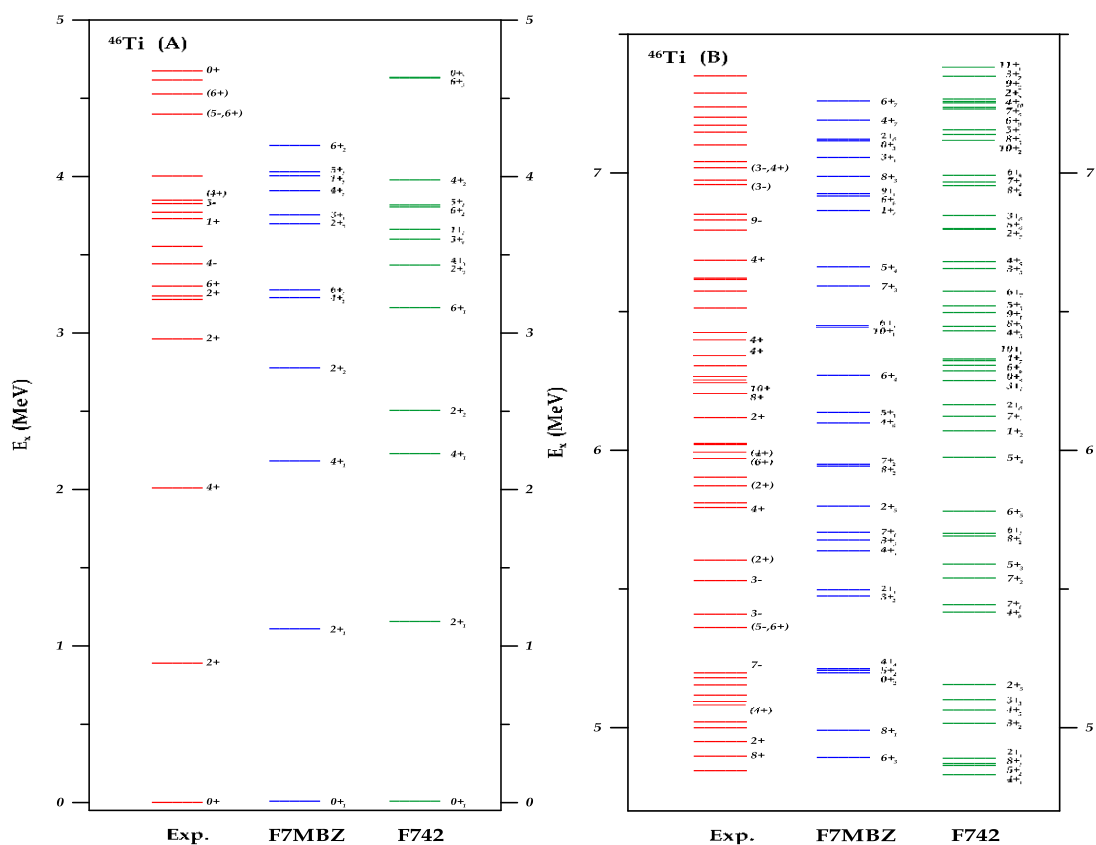
**Table (4): Excitation energy predicted by (F742) interactions and observed excitation energies[10] for the  $^{46}\text{Ti}$  nucleus are compared.**

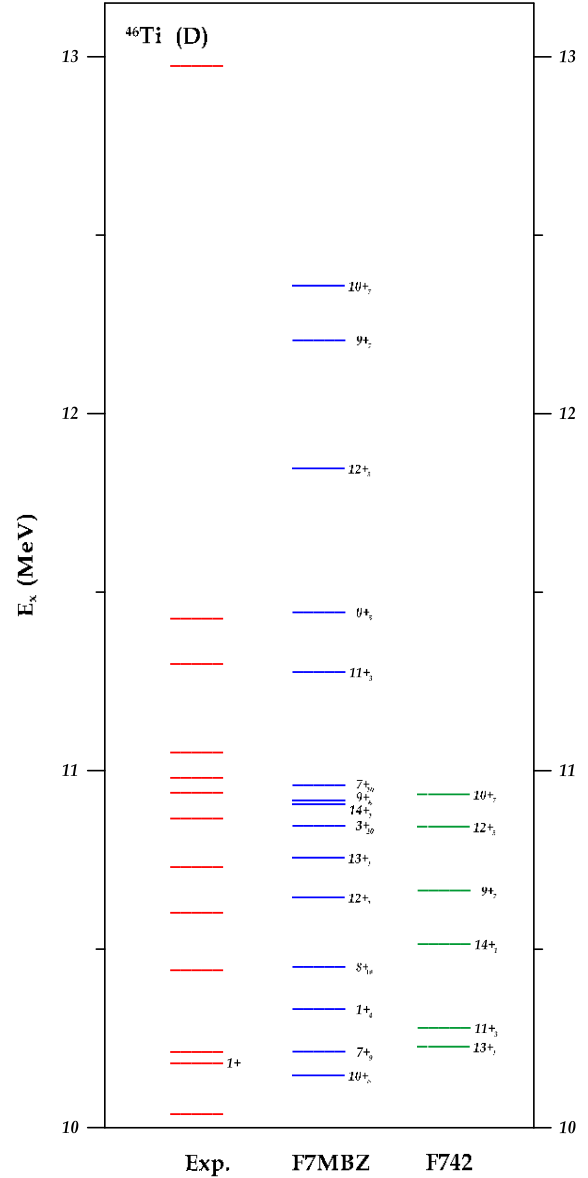
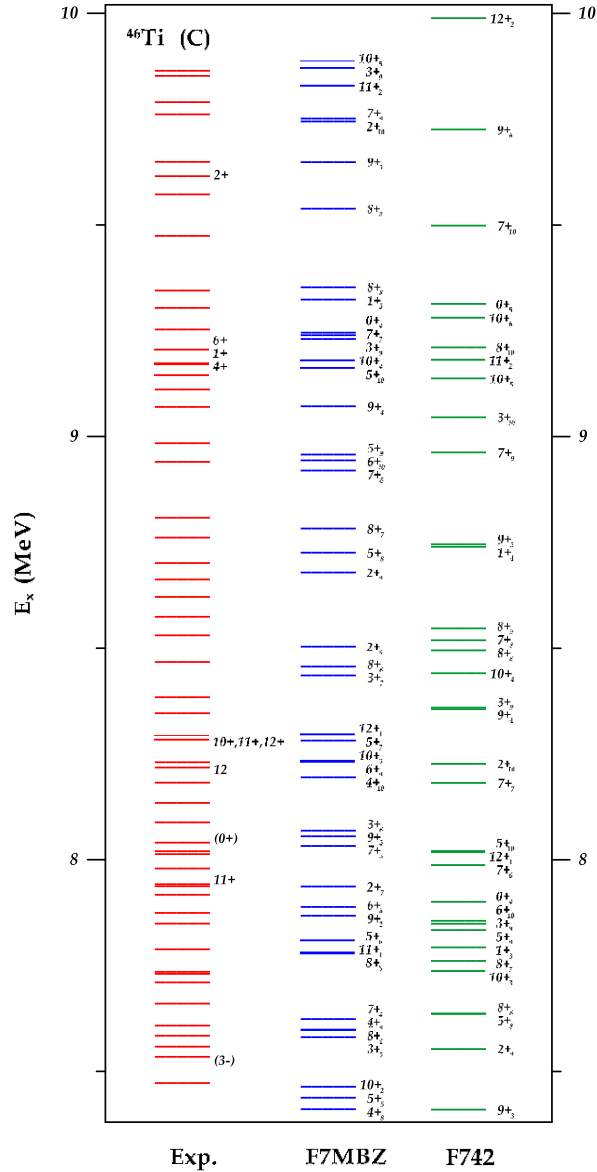
F742(Theoretical values)		Experimental values		F742Theoretical values		Experimental values	
$J^{\pi+}$	E (MeV)	E (MeV)	$J^{\pi}$	$J^{\pi+}$	E (MeV)	E (MeV)	$J^{\pi}$
0 <sub>1</sub>	0	0	0+	8 <sub>5</sub>	7.134	7.147	-----
2 <sub>1</sub>	1.148	0.889286	2+	5 <sub>7</sub>	7.151	7.172	-----
4 <sub>1</sub>	2.221	2.009846	4+	6 <sub>9</sub>	7.226	7.238	-----
2 <sub>2</sub>	2.497	2.9618	2+	7 <sub>5</sub>	7.232	7.288	-----
4 <sub>2</sub>	3.970	3.213	-----	4 <sub>10</sub>	7.248	7.019	(3-,4+)
6 <sub>1</sub>	3.154	3.29886	6+	2 <sub>8</sub>	7.253	-----9 <sub>2</sub>	-----
4 <sub>3</sub>	3.426	3.44139	4-2+		7.262	-----10 <sub>2</sub>	-----
2 <sub>3</sub>	3.426	3.2357	-----		7.110	7.172 3 <sub>7</sub>	--11+
3 <sub>1</sub>	3.591	3.5531	1+		7.344	7.350 11 <sub>1</sub>	-----
1 <sub>1</sub>	3.654	3.731	-----		7.377	7.9418 9 <sub>3</sub>	-----
6 <sub>2</sub>	3.797	3.7715	5-		7.405	7.472 2 <sub>9</sub>	-----
5 <sub>1</sub>	3.810	3.82643	(6+)		7.548	7.558 5 <sub>8</sub>	-----
6 <sub>3</sub>	4.620	4.527	0+		7.631	7.660 8 <sub>6</sub>	-----
0 <sub>2</sub>	4.625	4.675	-----		7.632	7.710 10 <sub>3</sub>	-----
4 <sub>4</sub>	4.826	4.617	-----		7.729	7.730 8 <sub>7</sub>	-----
5 <sub>2</sub>	4.859	4.845	8+		7.756	7.735 1 <sub>3</sub>	(0+)
8 <sub>1</sub>	4.866	4.8969	2+		7.788	7.788 5 <sub>9</sub>	-----12
2 <sub>4</sub>	4.885	4.950	-----		7.829	7.849 3 <sub>8</sub>	-----
3 <sub>2</sub>	5.011	5.000	(4+)		7.844	7.874 6 <sub>10</sub>	-----
4 <sub>5</sub>	5.059	5.079	-----		7.851	7.917 0 <sub>4</sub>	-----
3 <sub>3</sub>	5.096	5.094	-----		7.896	8.020 7 <sub>6</sub>	-----
2 <sub>5</sub>	5.151	5.117	-----		7.980	7.979 12 <sub>1</sub>	-----
4 <sub>6</sub>	5.412	5.154	-----		8.013	8.2175 5 <sub>10</sub>	-----
7 <sub>1</sub>	5.439	5.180	7-		8.015	8.013 7 <sub>7</sub>	-----
7 <sub>2</sub>	5.535	5.1976	-----		8.177	8.182 2 <sub>10</sub>	-----
5 <sub>3</sub>	5.585	5.811	8+		8.222	8.230 9 <sub>4</sub>	-----
8 <sub>2</sub>	5.686	6.2004	(5-,6+)		8.351	8.346 3 <sub>9</sub>	-----
6 <sub>4</sub>	5.696	5.361	(6+)		8.355	8.384 10 <sub>4</sub>	-----
6 <sub>5</sub>	5.776	5.965	-----		8.436	8.467 8 <sub>8</sub>	-----
5 <sub>4</sub>	5.970	5.903	-----2+		8.490	8.530 7 <sub>8</sub>	-----
1 <sub>2</sub>	6.066	6.025	-----		8.514	8.574 8 <sub>9</sub>	-----
7 <sub>3</sub>	6.118	5.021	-----		8.542	6.621 1 <sub>4</sub>	-----
2 <sub>6</sub>	6.160	6.118	-----		8.735	8.662 9 <sub>5</sub>	-----
3 <sub>4</sub>	6.247	6.251	4+		8.741	8.701 7 <sub>9</sub>	-----
0 <sub>3</sub>	6.282	6.266	-----		8.958	8.984 3 <sub>10</sub>	-----
6 <sub>6</sub>	6.302	6.305	-----		9.041	9.070 10 <sub>5</sub>	-----
4 <sub>7</sub>	6.319	6.338	-----		9.131	9.111	-----



10 <sub>1</sub>	6.325	6.2419	10+		11 <sub>2</sub>	9.175	9.141	----
4 <sub>8</sub>	6.426	6.395	4+		8 <sub>10</sub>	9.206	9.253	----
8 <sub>3</sub>	6.442	6.424	-----		10 <sub>6</sub>	9.275	9.304	----
9 <sub>1</sub>	6.492	-----	-----		0 <sub>5</sub>	9.309	9.345	----
5 <sub>5</sub>	6.516	6.513	-----		7 <sub>10</sub>	9.494	9.474	----
6 <sub>7</sub>	6.569	6.574	-----		9 <sub>6</sub>	9.721	9.761	----
3 <sub>5</sub>	6.651	6.616	-----		12 <sub>2</sub>	9.984	9.864	----
4 <sub>9</sub>	6.676	6.685	4+		13 <sub>1</sub>	10.220	10.038	----
2 <sub>7</sub>	6.792	6.794	-----		11 <sub>3</sub>	10.274	10.212	----
5 <sub>6</sub>	6.795	6.851	-----		14 <sub>1</sub>	10.509	10.441	----
3 <sub>6</sub>	6.842	6.958	(3-)		9 <sub>7</sub>	10.659	10.602	----
8 <sub>4</sub>	6.950	6.974	-----		12 <sub>3</sub>	10.836	10.866	----
7 <sub>4</sub>	6.963	7.041	-----		10 <sub>7</sub>	10.927	10.980	----
6 <sub>8</sub>	6.987	7.101	-----					

**Fig.2. A comparison between theoretical energy levels for two interaction and the experimental data for  $^{46}\text{Ti}$ .**





### 3.2 B(E2) Calculations :

#### 3.2.1 B(E2) for $^{44}\text{Ti}$

Within the nuclear shell model, (F7MBZ & F742) projected that the chance of an electric quadrupole transition B (E2) for  $^{44}\text{Ti}$  would be lower. The transition probability was determined for each in-band transition assuming a pure E2 transition by using the harmonic oscillator potential (HO, b), Where

$b < 0$ . Core polarization effects have been taken into account by selecting the effective charges for the proton ( $e_p=1.46e$ ) and the neutron ( $e_n=0.5e$ ). Table 5  $^{44}\text{Ti}$  was calculated with the help of the efficient interactions of F7MBZ and F742. Over all, there appears to be a fair amount of concordance between the computed results and the available experimental data.

**Table 5: the B (E2) values for 44Ti ground-state band. They use e2fm4 units, which match the experimental results.**

$J_i^+$	$\rightarrow$	$J_f^+$	Theory B(E2) ( $e^2 \text{ fm}^4$ )		Exp. B(E2) ( $e^2 \text{ fm}^4$ ) $e_p = 1.46e$ , $e_n = 0.5e$ $b = 1.888 \text{ fm}$
			F7MBZ	F742	
2 <sub>1</sub>	$\rightarrow$	0 <sub>1</sub>	119.9	119.8	119.95
0 <sub>2</sub>	$\rightarrow$	2 <sub>1</sub>	7.257	14.03	3137.152
2 <sub>2</sub>	$\rightarrow$	2 <sub>1</sub>	21.55	27.96	64.588
2 <sub>2</sub>	$\rightarrow$	0 <sub>2</sub>	12.77	12.64	212.219
2 <sub>2</sub>	$\rightarrow$	0 <sub>1</sub>	1.285	1.029	1.384
2 <sub>3</sub>	$\rightarrow$	0 <sub>1</sub>	0.03518	0.0005722	5.259
2 <sub>3</sub>	$\rightarrow$	0 <sub>2</sub>	0.02539	13.38	55.362
2 <sub>4</sub>	$\rightarrow$	0 <sub>1</sub>	0.06678	0.03225	1.199
4 <sub>1</sub>	$\rightarrow$	2 <sub>1</sub>	154.7	150.5	276.808
4 <sub>2</sub>	$\rightarrow$	2 <sub>1</sub>	2.295	6.207	24.913
4 <sub>2</sub>	$\rightarrow$	2 <sub>2</sub>	28.96	28.03	193.765
4 <sub>3</sub>	$\rightarrow$	2 <sub>1</sub>	0.007912	0.09484	4.152
4 <sub>3</sub>	$\rightarrow$	2 <sub>3</sub>	0.02051	5.635	258.354
6 <sub>1</sub>	$\rightarrow$	4 <sub>1</sub>	89.09	65.62	156.858

### 3.2.2 B(E2) for 46Ti

Within the nuclear shell model, (F7MBZ & F742) projected that the chance of an electric quadruple transition B (E2) for 46Ti would be lower. The transition probability was determined for each in-band transition assuming a pure E2 transition by using the harmonic oscillator potential (HO, b), Where

$b < 0$ . Core polarization effects have been taken into account by selecting the effective charges for the proton ( $e_p = 1.157e$ ) and the neutron ( $e_n = 0.771e$ ). Table 6 46Ti was calculated with the help of the efficient interactions of F7MBZ and F742. Over all, there appears to be a fair amount of concordance between the computed results and the available experimental data.

**Table 6: the B (E2) values for 46Ti ground-state band. They use e2fm4 units, which match the experimental results**

$J_i^+$	$\rightarrow$	$J_f^+$	Theory B(E2) ( $e^2 \text{ fm}^4$ )		Exp. B(E2) ( $e^2 \text{ fm}^4$ ) $e_p = 1.57e$ , $e_n = 0.771e$ $b = 1.902 \text{ fm}$
			F7MBZ	F742	
2 <sub>1</sub>	$\rightarrow$	0 <sub>1</sub>	120.1	119.6	-----
4 <sub>1</sub>	$\rightarrow$	2 <sub>1</sub>	125.5	121.9	489.516
0 <sub>2</sub>	$\rightarrow$	2 <sub>1</sub>	13.69	20.78	50.910
2 <sub>2</sub>	$\rightarrow$	2 <sub>1</sub>	50.3	41.96	

2 <sub>2</sub>	→	0 <sub>1</sub>	11.05	12.49	0.627
2 <sub>3</sub>	→	0 <sub>1</sub>	4.476	3.012	8.713
6 <sub>1</sub>	→	4 <sub>1</sub>	125	108.2	160.561
0 <sub>4</sub>	→	2 <sub>1</sub>	0.08582	0.00218	21.245
8 <sub>1</sub>	→	6 <sub>1</sub>	133.2	124.1	110.631
10 <sub>2</sub>	→	8 <sub>1</sub>	66.48	69.85	152.729

#### 4. Conclusions

All figures show, success in reaching an agreement nearly all energy levels of the isotopes of Titanium, and a proper reproduction of the level arrangement is made. We can evaluate practically any calculations in relation to (F7MBZ & F742) data. Met with some success in replicating the level structure exhibited. Generally speaking, the greatest and most thorough results possess the biggest model space while performing calculations in the infinite sphere. In OXBASH, the model space is described based on the nucleon valence orbits that are now excited together with the outcomes of our calculations are often in agreement with experimental findings

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