



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:01 PM GMT

PDB ID : 2TEC  
Title : MOLECULAR DYNAMICS REFINEMENT OF A THERMITASE-EGLIN-C COMPLEX AT 1.98 ANGSTROMS RESOLUTION AND COMPARISON OF TWO CRYSTAL FORMS THAT DIFFER IN CALCIUM CONTENT  
Authors : Gros, P.; Betzel, C.; Dauter, Z.; Wilson, K.S.; Hol, W.G.J.  
Deposited on : 1990-10-26  
Resolution : 1.98 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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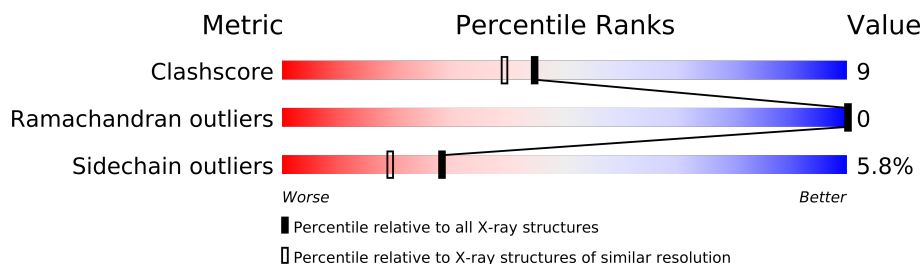
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.98 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	8091 (2.00-1.96)
Ramachandran outliers	78287	7989 (2.00-1.96)
Sidechain outliers	78261	7987 (2.00-1.96)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	E	279	
2	I	70	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2753 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called THERMITASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	279	Total	C	N	O	S	0	0	0
			2004	1242	351	409	2			

- Molecule 2 is a protein called EGLIN C.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	63	Total	C	N	O	0	0	0
			522	339	89	94			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	3	Total	Ca	0	0
			3	3		

- Molecule 4 is water.

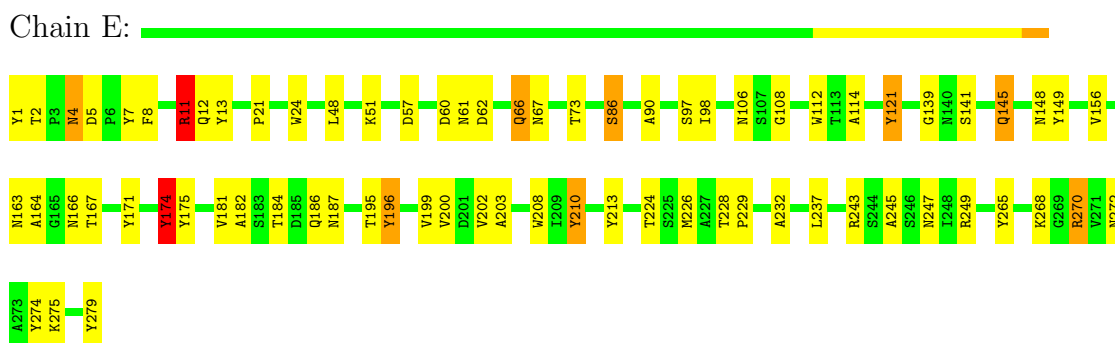
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	166	Total	O	0	0
			166	166		
4	I	58	Total	O	0	0
			58	58		

### 3 Residue-property plots

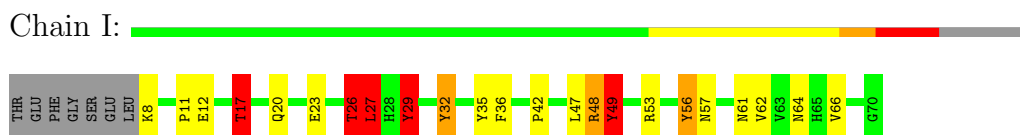
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

#### • Molecule 1: THERMITASE



#### • Molecule 2: EGLIN C



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	49.06Å 67.23Å 90.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.98	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.98)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	GROMOS	Depositor
R, $R_{free}$	0.165 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2753	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	11.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	E	0.95	0/2049	1.50	26/2805 (0.9%)
2	I	0.94	0/540	1.64	12/738 (1.6%)
All	All	0.95	0/2589	1.53	38/3543 (1.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	E	0	17
2	I	0	7
All	All	0	24

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	270	ARG	NE-CZ-NH1	12.48	126.54	120.30
1	E	265	TYR	CB-CG-CD2	-12.42	113.55	121.00
1	E	274	TYR	CB-CG-CD1	-11.96	113.82	121.00
1	E	270	ARG	NE-CZ-NH2	-11.72	114.44	120.30
1	E	279	TYR	CB-CG-CD2	-9.30	115.42	121.00
1	E	174	TYR	CB-CG-CD2	-8.68	115.79	121.00
1	E	175	TYR	CB-CG-CD2	-8.61	115.83	121.00
1	E	145	GLN	CB-CA-C	-7.34	95.72	110.40
1	E	11	ARG	NH1-CZ-NH2	-7.29	111.38	119.40
2	I	53	ARG	NE-CZ-NH1	7.18	123.89	120.30
1	E	121	TYR	CB-CG-CD1	-7.07	116.76	121.00
1	E	13	TYR	CB-CG-CD2	-6.93	116.84	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	48	ARG	NE-CZ-NH1	6.92	123.76	120.30
2	I	29	TYR	CB-CG-CD1	-6.69	116.98	121.00
1	E	270	ARG	CD-NE-CZ	6.54	132.76	123.60
1	E	1	TYR	CB-CG-CD2	-6.52	117.09	121.00
2	I	27	LEU	CB-CG-CD1	6.49	122.03	111.00
1	E	5	ASP	CB-CG-OD1	-6.05	112.85	118.30
2	I	35	TYR	CB-CG-CD2	-6.04	117.38	121.00
2	I	17	THR	CA-CB-CG2	6.01	120.81	112.40
2	I	48	ARG	NH1-CZ-NH2	-5.99	112.82	119.40
2	I	17	THR	N-CA-CB	-5.95	99.00	110.30
2	I	56	TYR	CB-CG-CD1	-5.76	117.54	121.00
1	E	274	TYR	CG-CD2-CE2	-5.58	116.83	121.30
1	E	210	TYR	CB-CG-CD2	-5.54	117.67	121.00
1	E	274	TYR	CD1-CG-CD2	5.53	123.99	117.90
1	E	114	ALA	CB-CA-C	5.44	118.26	110.10
1	E	171	TYR	CB-CG-CD2	-5.43	117.74	121.00
1	E	279	TYR	CG-CD1-CE1	-5.43	116.96	121.30
1	E	181	VAL	CG1-CB-CG2	-5.41	102.25	110.90
1	E	11	ARG	NE-CZ-NH2	-5.35	117.63	120.30
1	E	97	SER	N-CA-CB	-5.28	102.57	110.50
1	E	249	ARG	NE-CZ-NH1	5.23	122.92	120.30
2	I	29	TYR	CG-CD1-CE1	-5.23	117.12	121.30
2	I	49	TYR	CB-CG-CD2	-5.22	117.87	121.00
2	I	32	TYR	CB-CG-CD1	-5.21	117.88	121.00
1	E	202	VAL	CA-CB-CG1	5.15	118.62	110.90
1	E	60	ASP	CB-CG-OD2	-5.12	113.69	118.30

There are no chirality outliers.

All (24) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	108	GLY	Mainchain
1	E	11	ARG	Sidechain
1	E	121	TYR	Sidechain
1	E	139	GLY	Mainchain
1	E	148	ASN	Mainchain
1	E	149	TYR	Sidechain
1	E	174	TYR	Sidechain
1	E	187	ASN	Mainchain
1	E	195	THR	Mainchain
1	E	196	TYR	Sidechain
1	E	210	TYR	Sidechain

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Mol	Chain	Res	Type	Group
1	E	213	TYR	Sidechain
1	E	232	ALA	Mainchain
1	E	270	ARG	Sidechain
1	E	272	ASN	Sidechain
1	E	57	ASP	Mainchain
1	E	7	TYR	Sidechain
2	I	26	THR	Mainchain
2	I	29	TYR	Sidechain
2	I	36	PHE	Sidechain
2	I	47	LEU	Mainchain
2	I	48	ARG	Sidechain
2	I	49	TYR	Sidechain
2	I	56	TYR	Sidechain

## 5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2004	0	1908	34	0
2	I	522	0	498	11	0
3	E	3	0	0	0	0
4	E	166	0	0	10	1
4	I	58	0	0	2	0
All	All	2753	0	2406	42	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (42) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:11:ARG:HB3	1:E:11:ARG:HH11	1.48	0.78
2:I:17:THR:HG22	2:I:20:GLN:H	1.47	0.78
1:E:66:GLN:HE21	1:E:66:GLN:H	1.34	0.76
1:E:12:GLN:HE22	1:E:90:ALA:H	1.32	0.76
1:E:156:VAL:HB	4:E:497:HOH:O	1.87	0.74

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:163:ASN:HD21	1:E:224:THR:H	1.36	0.72
1:E:186:GLN:HG2	4:E:433:HOH:O	1.96	0.64
1:E:141:SER:O	1:E:145:GLN:HG2	1.98	0.62
1:E:61:ASN:HB3	4:E:347:HOH:O	2.01	0.61
1:E:174:TYR:HE2	1:E:199:VAL:HG11	1.65	0.61
1:E:4:ASN:HD21	1:E:86:SER:H	1.48	0.61
1:E:243:ARG:HH21	1:E:247:ASN:HD21	1.48	0.60
1:E:174:TYR:CE2	1:E:199:VAL:HG11	2.41	0.55
2:I:12:GLU:HB2	4:I:108:HOH:O	2.08	0.53
1:E:164:ALA:HB3	4:E:449:HOH:O	2.09	0.52
1:E:12:GLN:NE2	1:E:90:ALA:H	2.07	0.48
2:I:57:ASN:HB2	2:I:64:ASN:ND2	2.28	0.48
1:E:112:TRP:HD1	4:E:493:HOH:O	1.97	0.48
1:E:8:PHE:HA	1:E:12:GLN:HE21	1.79	0.48
1:E:243:ARG:HH21	1:E:247:ASN:ND2	2.11	0.48
1:E:66:GLN:NE2	1:E:66:GLN:H	2.06	0.47
1:E:51:LYS:HB3	1:E:98:ILE:HD12	1.96	0.47
1:E:106:ASN:HB2	4:E:364:HOH:O	2.14	0.46
2:I:23:GLU:HG2	2:I:27:LEU:HD22	1.96	0.46
1:E:182:ALA:HB2	1:E:200:VAL:HG11	1.98	0.46
1:E:275:LYS:HG2	4:E:444:HOH:O	2.17	0.45
1:E:112:TRP:CZ3	2:I:42:PRO:HB3	2.52	0.45
1:E:21:PRO:HA	1:E:24:TRP:CE3	2.51	0.44
1:E:166:ASN:HB2	4:E:473:HOH:O	2.17	0.44
1:E:4:ASN:ND2	1:E:86:SER:H	2.15	0.43
1:E:245:ALA:HA	4:E:497:HOH:O	2.18	0.43
1:E:184:THR:HG23	1:E:203:ALA:HB1	2.01	0.42
1:E:62:ASP:HB2	4:E:346:HOH:O	2.19	0.42
1:E:208:TRP:CH2	2:I:49:TYR:HE1	2.37	0.42
1:E:208:TRP:HH2	2:I:49:TYR:HE1	1.66	0.42
2:I:29:TYR:HB3	2:I:32:TYR:HD1	1.85	0.42
1:E:67:ASN:ND2	1:E:73:THR:H	2.18	0.41
2:I:11:PRO:HA	2:I:66:VAL:HG13	2.03	0.41
2:I:23:GLU:O	2:I:27:LEU:HB2	2.21	0.41
1:E:228:THR:HB	1:E:229:PRO:HD3	2.02	0.41
1:E:167:THR:HA	1:E:196:TYR:O	2.21	0.40
2:I:26:THR:HG22	4:I:78:HOH:O	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:E:416:HOH:O	4:E:424:HOH:O[3.655]	2.02	0.18

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	277/279 (99%)	268 (97%)	9 (3%)	0	100	100
2	I	61/70 (87%)	60 (98%)	1 (2%)	0	100	100
All	All	338/349 (97%)	328 (97%)	10 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	202/202 (100%)	193 (96%)	9 (4%)	38	29
2	I	58/64 (91%)	52 (90%)	6 (10%)	10	5
All	All	260/266 (98%)	245 (94%)	15 (6%)	28	19

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	2	THR
1	E	4	ASN
1	E	11	ARG
1	E	48	LEU

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Mol	Chain	Res	Type
1	E	66	GLN
1	E	86	SER
1	E	226	MET
1	E	237	LEU
1	E	268	LYS
2	I	8	LYS
2	I	17	THR
2	I	26	THR
2	I	27	LEU
2	I	61	ASN
2	I	62	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	4	ASN
1	E	12	GLN
1	E	117	ASN
1	E	148	ASN
1	E	163	ASN
1	E	177	ASN
1	E	247	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.