



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 01:25 PM UTC

PDB ID : 1TEC / pdb_00001tec
Title : CRYSTALLOGRAPHIC REFINEMENT BY INCORPORATION OF MOLECULAR DYNAMICS. THE THERMOSTABLE SERINE PROTEASE THERMITASE COMPLEXED WITH EGLIN-C
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Deposited on : 1989-05-24
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

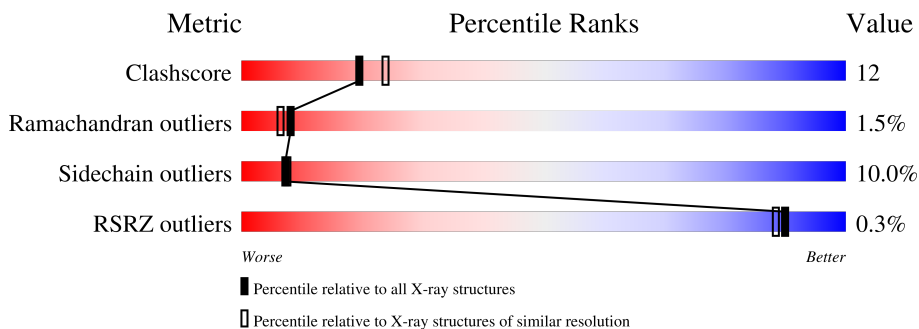
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

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

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2737 atoms, of which 0 are hydrogens and 0 are deuteriums.

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
			Total	C	N	O	S			
1	E	279	2004	1242	351	409	2	0	0	0

- Molecule 2 is a protein called EGLIN C.

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

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
3	E	2	2	2	0	0

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Na		
4	E	1	1	1	0	0

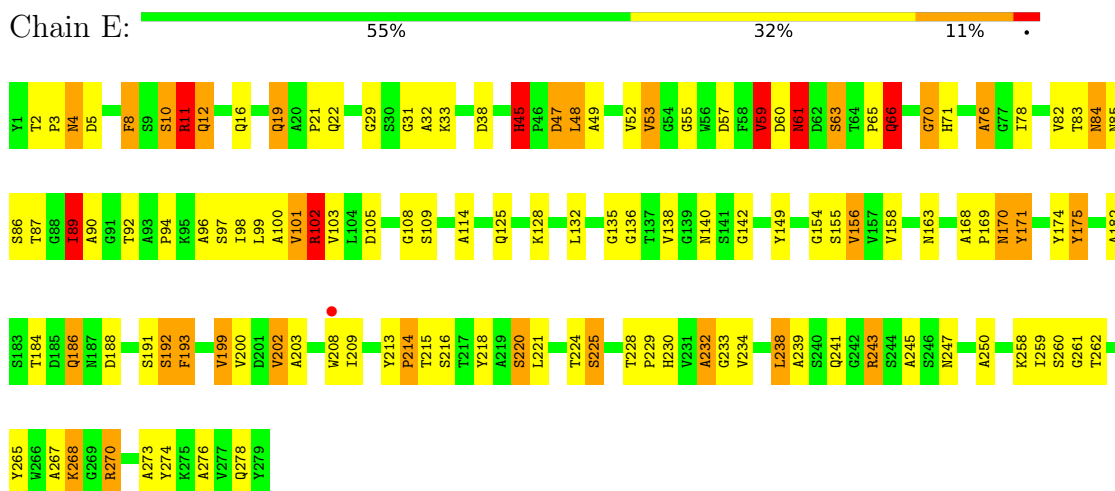
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	E	166	166	166	0	0
5	I	42	42	42	0	0

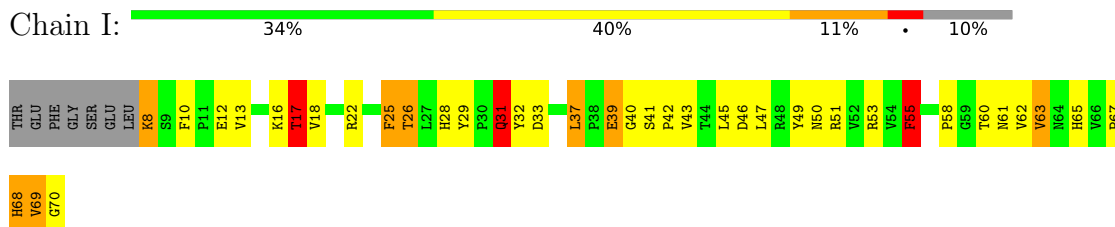
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

- Molecule 1: THERMITASE



- Molecule 2: EGLIN C



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.25Å 72.10Å 89.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.20 56.09 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.20) 76.2 (56.09-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.64 (at 2.20Å)	Xtrriage
Refinement program	GROMOS	Depositor
R, R_{free}	(Not available) , (Not available) 0.174 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	24.0	Xtrriage
Anisotropy	0.902	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 67.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2737	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: NA, 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	1.31	3/2049 (0.1%)	1.99	41/2805 (1.5%)
2	I	1.25	0/540	1.95	12/738 (1.6%)
All	All	1.30	3/2589 (0.1%)	1.99	53/3543 (1.5%)

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	43
2	I	0	19
All	All	0	62

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	71	HIS	CG-ND1	-7.05	1.30	1.38
1	E	70	GLY	CA-C	5.80	1.57	1.52
1	E	230	HIS	CG-ND1	-5.50	1.32	1.38

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	156	VAL	N-CA-CB	9.68	121.08	110.53
2	I	17	THR	N-CA-CB	-8.25	98.29	110.17
1	E	259	ILE	N-CA-C	-7.62	101.74	110.05
1	E	267	ALA	N-CA-C	7.62	119.66	111.36
1	E	89	ILE	CA-CB-CG2	7.61	123.44	110.50
2	I	37	LEU	N-CA-CB	-7.29	98.07	111.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	213	TYR	CA-C-N	6.95	126.71	119.76
1	E	213	TYR	C-N-CA	6.95	126.71	119.76
2	I	46	ASP	CA-CB-CG	6.94	119.54	112.60
1	E	278	GLN	OE1-CD-NE2	-6.86	115.74	122.60
1	E	170	ASN	CA-CB-CG	-6.77	105.83	112.60
1	E	214	PRO	CB-CA-C	-6.77	102.21	111.21
1	E	12	GLN	OE1-CD-NE2	-6.76	115.84	122.60
1	E	19	GLN	OE1-CD-NE2	-6.73	115.87	122.60
1	E	47	ASP	CA-CB-CG	-6.69	105.91	112.60
1	E	59	VAL	N-CA-CB	6.58	118.64	111.41
2	I	68	HIS	CB-CG-CD2	-6.45	122.82	131.20
1	E	5	ASP	CA-C-O	-6.24	113.32	119.49
2	I	12	GLU	N-CA-C	6.23	120.14	112.54
1	E	214	PRO	N-CA-CB	6.09	109.01	103.34
1	E	32	ALA	N-CA-C	-6.09	99.27	109.07
1	E	238	LEU	N-CA-CB	-6.00	101.26	110.20
2	I	50	ASN	OD1-CG-ND2	-5.97	116.63	122.60
1	E	168	ALA	CA-C-N	5.92	125.86	119.76
1	E	168	ALA	C-N-CA	5.92	125.86	119.76
1	E	202	VAL	CA-CB-CG2	5.91	120.44	110.40
1	E	241	GLN	OE1-CD-NE2	-5.88	116.72	122.60
1	E	188	ASP	CA-CB-CG	-5.86	106.74	112.60
1	E	78	ILE	N-CA-CB	-5.85	104.02	110.51
2	I	28	HIS	CA-CB-CG	-5.79	108.01	113.80
1	E	38	ASP	N-CA-CB	-5.78	103.61	109.51
1	E	61	ASN	CA-CB-CG	-5.71	106.89	112.60
1	E	238	LEU	CA-C-N	5.68	127.90	120.28
1	E	238	LEU	C-N-CA	5.68	127.90	120.28
2	I	13	VAL	CA-CB-CG2	5.65	120.00	110.40
1	E	169	PRO	N-CA-CB	5.64	108.33	103.31
1	E	33	LYS	N-CA-CB	-5.57	101.82	110.57
2	I	45	LEU	CB-CA-C	5.54	118.55	111.40
2	I	68	HIS	ND1-CG-CD2	5.53	111.63	106.10
1	E	175	TYR	N-CA-C	-5.52	99.60	108.55
1	E	192	SER	CA-C-O	-5.42	114.80	120.55
1	E	268	LYS	N-CA-C	5.38	119.97	113.41
2	I	31	GLN	CB-CA-C	-5.37	98.98	110.32
1	E	220	SER	CA-CB-OG	-5.34	100.41	111.10
1	E	65	PRO	N-CA-CB	5.27	106.09	102.35
1	E	84	ASN	CA-CB-CG	-5.21	107.39	112.60
1	E	225	SER	CA-C-N	5.18	127.22	120.28
1	E	225	SER	C-N-CA	5.18	127.22	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	102	ARG	NE-CZ-NH2	5.15	123.84	119.20
1	E	45	HIS	CB-CG-CD2	-5.08	124.60	131.20
1	E	11	ARG	CA-CB-CG	-5.04	104.03	114.10
1	E	261	GLY	O-C-N	-5.03	118.04	122.77
2	I	55	PHE	CA-CB-CG	-5.01	108.79	113.80

There are no chirality outliers.

All (62) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	101	VAL	Mainchain
1	E	102	ARG	Mainchain
1	E	105	ASP	Mainchain
1	E	108	GLY	Mainchain
1	E	11	ARG	Sidechain
1	E	114	ALA	Mainchain
1	E	132	LEU	Mainchain
1	E	136	GLY	Mainchain
1	E	142	GLY	Mainchain
1	E	149	TYR	Sidechain
1	E	155	SER	Mainchain
1	E	16	GLN	Mainchain
1	E	171	TYR	Sidechain
1	E	174	TYR	Mainchain,Sidechain
1	E	175	TYR	Sidechain
1	E	191	SER	Mainchain
1	E	192	SER	Mainchain
1	E	193	PHE	Sidechain
1	E	209	ILE	Mainchain
1	E	21	PRO	Mainchain
1	E	216	SER	Mainchain
1	E	218	TYR	Sidechain
1	E	232	ALA	Mainchain
1	E	243	ARG	Sidechain
1	E	250	ALA	Mainchain
1	E	258	LYS	Mainchain
1	E	262	THR	Mainchain
1	E	265	TYR	Sidechain
1	E	270	ARG	Mainchain
1	E	3	PRO	Mainchain
1	E	31	GLY	Mainchain
1	E	45	HIS	Sidechain

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Mol	Chain	Res	Type	Group
1	E	52	VAL	Mainchain
1	E	60	ASP	Mainchain
1	E	61	ASN	Mainchain
1	E	63	SER	Mainchain
1	E	66	GLN	Mainchain
1	E	70	GLY	Mainchain
1	E	76	ALA	Mainchain
1	E	8	PHE	Sidechain
1	E	84	ASN	Mainchain
1	E	87	THR	Mainchain
2	I	16	LYS	Mainchain
2	I	18	VAL	Mainchain
2	I	25	PHE	Sidechain
2	I	26	THR	Mainchain
2	I	29	TYR	Sidechain
2	I	31	GLN	Mainchain
2	I	32	TYR	Sidechain
2	I	39	GLU	Mainchain
2	I	42	PRO	Mainchain
2	I	43	VAL	Mainchain
2	I	49	TYR	Sidechain
2	I	53	ARG	Sidechain
2	I	55	PHE	Mainchain
2	I	58	PRO	Mainchain
2	I	63	VAL	Mainchain
2	I	65	HIS	Mainchain,Sidechain
2	I	69	VAL	Mainchain
2	I	8	LYS	Mainchain

5.2 Too-close contacts [\(i\)](#)

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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2004	0	1909	48	0
2	I	522	0	498	12	1
3	E	2	0	0	0	0
4	E	1	0	0	0	0
5	E	166	0	0	2	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	I	42	0	0	2	0
All	All	2737	0	2407	57	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (57) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:186:GLN:NE2	1:E:186:GLN:H	1.63	0.96
1:E:243:ARG:HD3	1:E:247:ASN:HD22	1.52	0.75
1:E:243:ARG:HD3	1:E:247:ASN:ND2	2.01	0.75
1:E:4:ASN:ND2	1:E:86:SER:H	1.86	0.73
1:E:163:ASN:HD21	1:E:224:THR:H	1.38	0.72
1:E:59:VAL:HG22	1:E:102:ARG:NH1	2.04	0.71
1:E:19:GLN:HE21	1:E:22:GLN:NE2	1.89	0.70
1:E:66:GLN:HE21	1:E:66:GLN:H	1.39	0.70
1:E:138:VAL:HG12	2:I:40:GLY:HA3	1.73	0.70
1:E:186:GLN:H	1:E:186:GLN:HE21	1.44	0.65
2:I:10:PHE:HB2	2:I:67:PRO:HG2	1.79	0.65
1:E:182:ALA:HB2	1:E:200:VAL:HG11	1.79	0.65
1:E:186:GLN:H	1:E:186:GLN:CD	2.07	0.63
1:E:8:PHE:HA	1:E:12:GLN:HE21	1.62	0.63
1:E:128:LYS:HB2	5:E:473:HOH:O	2.01	0.60
1:E:184:THR:HG23	1:E:203:ALA:HB1	1.84	0.60
2:I:25:PHE:HB3	5:I:78:HOH:O	2.02	0.58
1:E:225:SER:O	1:E:229:PRO:HD3	2.04	0.58
1:E:229:PRO:HA	1:E:232:ALA:HB3	1.85	0.58
1:E:138:VAL:HG12	2:I:40:GLY:CA	2.34	0.56
1:E:220:SER:O	1:E:221:LEU:HD23	2.05	0.56
1:E:76:ALA:HB1	1:E:98:ILE:HD13	1.91	0.53
2:I:60:THR:OG1	2:I:62:VAL:HG22	2.09	0.52
2:I:8:LYS:O	2:I:68:HIS:HB2	2.09	0.52
1:E:47:ASP:O	1:E:82:VAL:HG23	2.09	0.52
1:E:101:VAL:HG12	1:E:103:VAL:HG13	1.92	0.50
1:E:158:VAL:HG12	1:E:228:THR:HG23	1.94	0.50
2:I:22:ARG:O	2:I:26:THR:HG23	2.11	0.50
1:E:199:VAL:HG11	5:E:430:HOH:O	2.14	0.48
1:E:45:HIS:HB3	1:E:48:LEU:HB2	1.96	0.48
1:E:29:GLY:O	1:E:96:ALA:HB2	2.13	0.48
2:I:70:GLY:HA3	5:I:107:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:83:THR:HB	1:E:94:PRO:HB3	1.96	0.47
1:E:186:GLN:NE2	1:E:186:GLN:N	2.47	0.47
1:E:208:TRP:HE3	1:E:220:SER:OG	1.97	0.47
1:E:234:VAL:HG22	1:E:273:ALA:HB2	1.97	0.47
1:E:66:GLN:H	1:E:66:GLN:NE2	2.09	0.46
1:E:238:LEU:HD11	1:E:276:ALA:HB1	1.99	0.45
1:E:55:GLY:HA3	1:E:100:ALA:O	2.17	0.45
1:E:4:ASN:C	1:E:4:ASN:HD22	2.24	0.45
1:E:19:GLN:O	1:E:274:TYR:HB2	2.16	0.45
1:E:57:ASP:OD1	1:E:57:ASP:C	2.60	0.45
1:E:186:GLN:HE21	1:E:186:GLN:N	2.12	0.44
1:E:128:LYS:HG2	1:E:239:ALA:HB1	2.00	0.44
1:E:10:SER:C	1:E:11:ARG:HG3	2.43	0.44
2:I:17:THR:HA	2:I:61:ASN:O	2.19	0.43
1:E:92:THR:HB	1:E:233:GLY:HA3	1.99	0.43
1:E:89:ILE:HG23	1:E:90:ALA:O	2.19	0.43
2:I:37:LEU:O	2:I:55:PHE:HA	2.19	0.42
1:E:135:GLY:HA2	1:E:171:TYR:O	2.19	0.42
1:E:154:GLY:O	1:E:245:ALA:HB2	2.19	0.42
1:E:193:PHE:CD2	1:E:193:PHE:C	2.98	0.42
2:I:51:ARG:O	2:I:69:VAL:HA	2.20	0.41
1:E:10:SER:O	1:E:11:ARG:HG3	2.20	0.41
1:E:214:PRO:HA	1:E:215:THR:HA	1.75	0.41
1:E:53:VAL:HG13	1:E:97:SER:HB2	2.03	0.40
1:E:138:VAL:CG1	2:I:40:GLY:HA3	2.48	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:62:VAL:CG2	5:E:399:HOH:O[2_455]	2.15	0.05
5:E:429:HOH:O	5:E:459:HOH:O[3_544]	2.19	0.01

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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%)	254 (92%)	18 (6%)	5 (2%)	6	4
2	I	61/70 (87%)	58 (95%)	3 (5%)	0	100	100
All	All	338/349 (97%)	312 (92%)	21 (6%)	5 (2%)	8	6

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	49	ALA
1	E	53	VAL
1	E	61	ASN
1	E	125	GLN
1	E	140	ASN

5.3.2 Protein sidechains [i](#)

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%)	183 (91%)	19 (9%)	8	9
2	I	58/64 (91%)	51 (88%)	7 (12%)	5	4
All	All	260/266 (98%)	234 (90%)	26 (10%)	7	7

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	2	THR
1	E	4	ASN
1	E	10	SER
1	E	48	LEU
1	E	59	VAL
1	E	63	SER
1	E	66	GLN
1	E	85	ASN

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Mol	Chain	Res	Type
1	E	89	ILE
1	E	99	LEU
1	E	109	SER
1	E	156	VAL
1	E	170	ASN
1	E	186	GLN
1	E	199	VAL
1	E	202	VAL
1	E	260	SER
1	E	268	LYS
1	E	270	ARG
2	I	17	THR
2	I	31	GLN
2	I	33	ASP
2	I	39	GLU
2	I	41	SER
2	I	47	LEU
2	I	63	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	4	ASN
1	E	12	GLN
1	E	22	GLN
1	E	66	GLN
1	E	117	ASN
1	E	148	ASN
1	E	163	ASN
1	E	177	ASN
1	E	186	GLN
1	E	247	ASN
2	I	61	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	E	279/279 (100%)	-0.20	1 (0%) 88 87	2, 17, 30, 46	0
2	I	63/70 (90%)	-0.20	0 100 100	11, 20, 31, 38	0
All	All	342/349 (97%)	-0.20	1 (0%) 90 88	2, 18, 31, 46	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	208	TRP	3.7

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	CA	E	343	1/1	0.96	0.21	0,0,0,0	0
3	CA	E	344	1/1	0.96	0.15	11,11,11,11	0
4	NA	E	345	1/1	0.96	0.05	9,9,9,9	0

6.5 Other polymers [i](#)

There are no such residues in this entry.