



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:24 PM UTC

PDB ID : 1Q87 / pdb_00001q87
Title : Crystal structure of the C-domain of the T.vaginalis Inr binding protein, IBP39 (tetragonal form)
Authors : Schumacher, M.A.; Johnson, P.J.
Deposited on : 2003-08-20
Resolution : 2.32 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
Xtriage (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

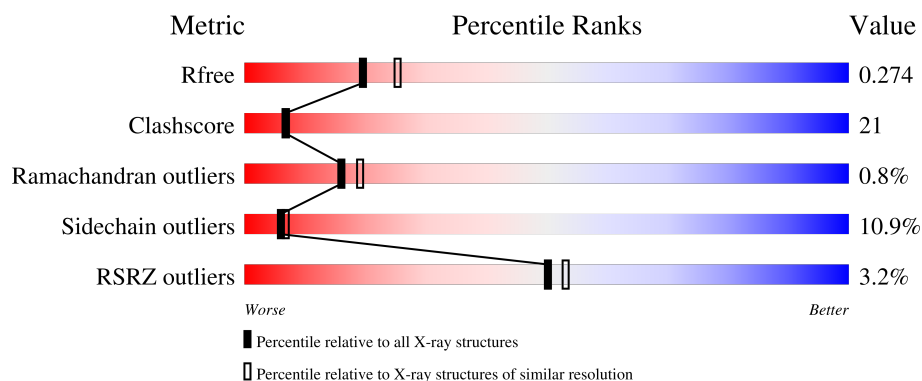
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.32 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))
R_{free}	180053	7754 (2.34-2.30)
Clashscore	190562	8383 (2.34-2.30)
Ramachandran outliers	187476	8303 (2.34-2.30)
Sidechain outliers	187428	8303 (2.34-2.30)
RSRZ outliers	180081	7760 (2.34-2.30)

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	A	221	
1	B	221	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3115 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 39 kDa initiator binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	185	Total	C	N	O	S	0	0	0
			1494	963	239	282	10			
1	B	193	Total	C	N	O	S	0	0	0
			1560	1005	249	296	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	342	HIS	-	expression tag	UNP Q95VR4
A	343	HIS	-	expression tag	UNP Q95VR4
A	344	HIS	-	expression tag	UNP Q95VR4
A	345	HIS	-	expression tag	UNP Q95VR4
A	346	HIS	-	expression tag	UNP Q95VR4
A	347	HIS	-	expression tag	UNP Q95VR4
B	342	HIS	-	expression tag	UNP Q95VR4
B	343	HIS	-	expression tag	UNP Q95VR4
B	344	HIS	-	expression tag	UNP Q95VR4
B	345	HIS	-	expression tag	UNP Q95VR4
B	346	HIS	-	expression tag	UNP Q95VR4
B	347	HIS	-	expression tag	UNP Q95VR4

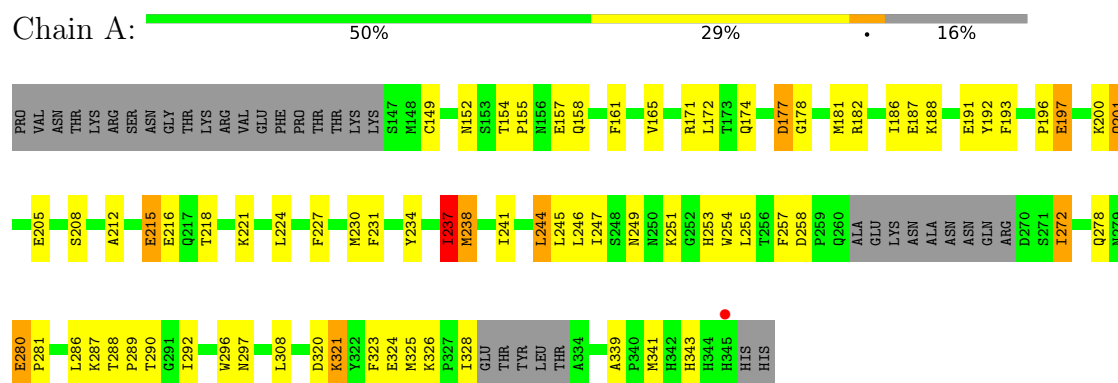
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	34	Total	O	0	0
			34	34		
2	B	27	Total	O	0	0
			27	27		

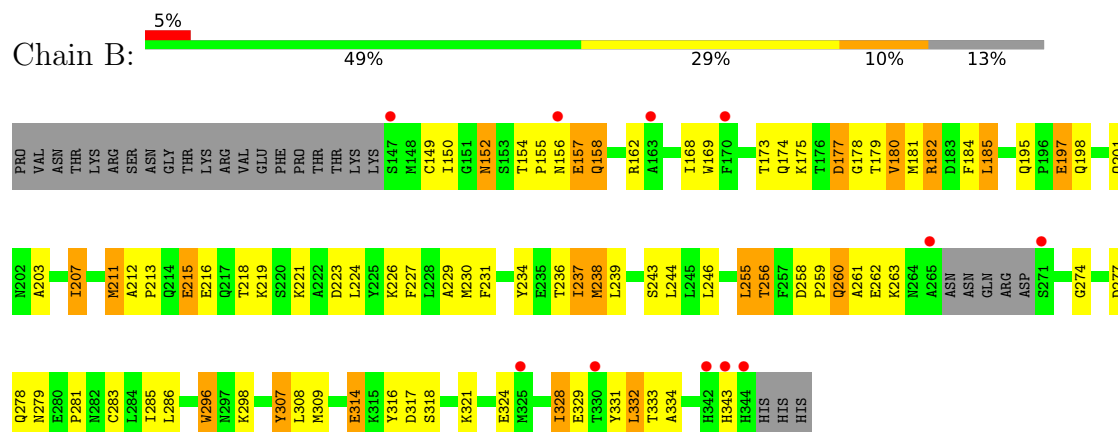
3 Residue-property plots

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: 39 kDa initiator binding protein



- Molecule 1: 39 kDa initiator binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	76.42Å 76.42Å 157.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	65.00 – 2.32 65.00 – 2.32	Depositor EDS
% Data completeness (in resolution range)	91.0 (65.00-2.32) 91.1 (65.00-2.32)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.32Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.215 , 0.274 0.217 , 0.274	Depositor DCC
R_{free} test set	1891 reflections (9.89%)	wwPDB-VP
Wilson B-factor (Å ²)	61.2	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 66.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3115	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% 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

5.1 Standard geometry

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	A	0.76	1/1531 (0.1%)	1.07	8/2070 (0.4%)
1	B	0.69	0/1599	1.03	9/2164 (0.4%)
All	All	0.73	1/3130 (0.0%)	1.05	17/4234 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	238	MET	SD-CE	-5.50	1.65	1.79

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	296	TRP	N-CA-C	7.83	121.66	108.90
1	A	339	ALA	CA-C-N	7.75	127.80	119.90
1	A	339	ALA	C-N-CA	7.75	127.80	119.90
1	B	298	LYS	CA-C-N	7.21	127.84	119.47
1	B	298	LYS	C-N-CA	7.21	127.84	119.47
1	A	296	TRP	N-CA-C	6.94	120.55	109.24
1	B	331	TYR	N-CA-C	6.79	119.32	110.43
1	A	343	HIS	N-CA-C	-6.47	104.87	112.89
1	B	260	GLN	N-CA-C	-6.16	106.44	112.97
1	A	324	GLU	N-CA-C	-6.14	103.91	111.40
1	B	343	HIS	N-CA-C	-5.94	105.99	113.18
1	A	237	ILE	N-CA-C	5.38	116.76	110.62
1	B	328	ILE	N-CA-C	-5.37	105.48	110.53
1	A	280	GLU	N-CA-C	-5.31	99.72	108.75
1	A	281	PRO	N-CA-C	5.30	123.39	112.47
1	B	307	TYR	N-CA-C	5.29	119.37	113.02
1	B	324	GLU	N-CA-C	-5.17	105.27	112.45

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	A	1494	0	1453	45	0
1	B	1560	0	1523	84	0
2	A	34	0	0	4	0
2	B	27	0	0	3	0
All	All	3115	0	2976	129	0

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 21.

All (129) 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:A:177:ASP:CG	1:A:178:GLY:H	1.74	0.94
1:B:317:ASP:H	1:B:321:LYS:HD2	1.48	0.78
1:B:177:ASP:OD1	1:B:178:GLY:N	2.19	0.76
1:A:177:ASP:CG	1:A:178:GLY:N	2.44	0.75
1:B:203:ALA:O	1:B:207:ILE:HG12	1.86	0.75
1:B:197:GLU:H	1:B:197:GLU:CD	1.91	0.75
1:A:341:MET:HA	1:A:341:MET:HE2	1.69	0.74
1:B:332:LEU:HD13	1:B:334:ALA:H	1.52	0.73
1:A:191:GLU:HG2	2:A:367:HOH:O	1.89	0.72
1:A:287:LYS:HD2	1:A:292:ILE:HD11	1.72	0.72
1:A:321:LYS:HE3	2:A:352:HOH:O	1.89	0.72
1:B:226:LYS:NZ	1:B:278:GLN:HE22	1.89	0.70
1:B:226:LYS:HG2	1:B:230:MET:HE2	1.72	0.70
1:B:332:LEU:CD1	1:B:334:ALA:H	2.03	0.70
1:B:180:VAL:HG13	1:B:184:PHE:HB3	1.74	0.69
1:B:286:LEU:HD22	1:B:328:ILE:HD11	1.76	0.67
1:A:215:GLU:H	1:A:215:GLU:CD	2.04	0.64
1:B:317:ASP:HB2	1:B:321:LYS:HD2	1.80	0.64
1:B:169:TRP:CE3	1:B:221:LYS:HA	2.34	0.63
1:B:185:LEU:HD13	1:B:211:MET:HE3	1.80	0.62
1:B:212:ALA:HB1	1:B:213:PRO:HA	1.79	0.62
1:B:229:ALA:HB3	1:B:281:PRO:HB2	1.81	0.62
1:B:185:LEU:HD22	1:B:211:MET:HE1	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:ASP:H	1:B:321:LYS:CD	2.13	0.60
1:B:213:PRO:HD3	1:B:259:PRO:HG2	1.84	0.59
1:A:287:LYS:HA	1:A:292:ILE:HD13	1.84	0.59
1:B:195:GLN:HB2	1:B:198:GLN:HB2	1.85	0.58
1:B:328:ILE:CG2	1:B:329:GLU:N	2.66	0.57
1:B:234:TYR:HA	1:B:237:ILE:HB	1.85	0.57
1:B:181:MET:O	1:B:184:PHE:N	2.38	0.56
1:A:171:ARG:HH21	1:A:192:TYR:HE2	1.54	0.56
1:B:227:PHE:HA	1:B:230:MET:HE3	1.86	0.56
1:A:244:LEU:HD13	1:A:255:LEU:HD22	1.87	0.56
1:B:263:LYS:HE2	1:B:263:LYS:H	1.71	0.56
1:B:149:CYS:SG	1:B:158:GLN:HG2	2.46	0.55
1:B:231:PHE:HB2	1:B:237:ILE:HD11	1.87	0.55
1:A:201:GLN:O	1:A:205:GLU:HG3	2.06	0.55
1:A:272:ILE:HD13	1:A:328:ILE:HG21	1.90	0.54
1:A:241:ILE:O	1:A:245:LEU:HB2	2.08	0.54
1:B:237:ILE:CG2	1:B:238:MET:N	2.71	0.54
1:B:243:SER:HA	2:B:372:HOH:O	2.08	0.53
1:B:180:VAL:HG13	1:B:181:MET:N	2.23	0.53
1:B:169:TRP:CZ2	1:B:173:THR:HG21	2.44	0.52
1:B:226:LYS:HZ2	1:B:278:GLN:HE22	1.56	0.52
1:A:172:LEU:HD22	1:A:188:LYS:HB3	1.92	0.52
1:B:175:LYS:HE3	1:B:177:ASP:O	2.10	0.51
1:B:277:ASP:HB2	1:B:285:ILE:HD11	1.93	0.51
1:B:229:ALA:CB	1:B:281:PRO:HB2	2.41	0.51
1:A:254:TRP:CZ3	1:A:286:LEU:HD21	2.47	0.50
1:B:258:ASP:O	1:B:261:ALA:HB2	2.11	0.50
1:B:262:GLU:HB3	1:B:263:LYS:HE2	1.94	0.50
1:B:155:PRO:HA	1:B:158:GLN:NE2	2.26	0.50
1:A:182:ARG:HD2	1:A:216:GLU:O	2.12	0.50
1:A:155:PRO:HA	1:A:158:GLN:HE21	1.76	0.50
1:A:174:GLN:HG3	2:A:371:HOH:O	2.12	0.50
1:B:328:ILE:HG23	1:B:329:GLU:N	2.26	0.49
1:B:152:ASN:CG	1:B:309:MET:HE3	2.38	0.49
1:B:231:PHE:HB3	1:B:237:ILE:HD13	1.95	0.49
1:A:288:THR:HB	1:A:289:PRO:HD2	1.95	0.49
1:A:191:GLU:HG3	1:A:200:LYS:HE3	1.94	0.49
1:A:197:GLU:H	1:A:197:GLU:CD	2.20	0.48
1:B:283:CYS:HB2	1:B:296:TRP:CD1	2.49	0.48
1:B:279:ASN:O	1:B:281:PRO:HD3	2.14	0.48
1:B:236:THR:HG22	1:B:239:LEU:HD12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:THR:HG22	1:B:219:LYS:N	2.28	0.47
1:A:231:PHE:C	1:A:237:ILE:HD12	2.38	0.47
1:A:325:MET:C	1:A:326:LYS:HG2	2.38	0.47
1:B:317:ASP:N	1:B:321:LYS:HD2	2.24	0.47
1:B:226:LYS:CD	1:B:278:GLN:HE22	2.27	0.47
1:B:219:LYS:HG3	1:B:223:ASP:OD2	2.14	0.47
1:A:154:THR:OG1	1:A:157:GLU:HG3	2.15	0.47
1:A:297:ASN:HA	1:A:308:LEU:HD23	1.97	0.46
1:B:180:VAL:CG1	1:B:181:MET:N	2.77	0.46
1:B:152:ASN:HD22	1:B:152:ASN:HA	1.56	0.46
1:B:332:LEU:HD13	1:B:333:THR:N	2.30	0.46
1:A:212:ALA:CB	1:A:258:ASP:HB2	2.46	0.46
1:A:181:MET:HA	1:A:181:MET:HE2	1.99	0.45
1:B:154:THR:OG1	1:B:157:GLU:HG3	2.17	0.45
1:B:286:LEU:HD22	1:B:328:ILE:CD1	2.44	0.45
1:A:238:MET:HE2	1:A:241:ILE:HD12	1.99	0.45
1:B:317:ASP:H	1:B:321:LYS:HZ2	1.64	0.45
1:B:182:ARG:HD2	1:B:216:GLU:O	2.16	0.44
1:B:317:ASP:N	1:B:321:LYS:HZ2	2.16	0.44
1:A:245:LEU:HA	1:A:245:LEU:HD12	1.76	0.44
1:B:155:PRO:HA	1:B:158:GLN:HE21	1.82	0.44
1:A:247:ILE:HD11	1:A:320:ASP:OD1	2.18	0.43
1:B:181:MET:HE2	1:B:181:MET:HA	1.99	0.43
1:B:174:GLN:HA	1:B:174:GLN:OE1	2.19	0.43
1:B:259:PRO:C	1:B:261:ALA:H	2.26	0.43
1:B:180:VAL:CG1	1:B:184:PHE:HB3	2.45	0.43
1:B:307:TYR:CZ	1:B:318:SER:HA	2.53	0.43
1:A:215:GLU:CD	1:A:215:GLU:N	2.75	0.43
1:B:317:ASP:CB	1:B:321:LYS:HD2	2.47	0.43
1:A:196:PRO:HD2	1:A:197:GLU:OE2	2.19	0.43
1:A:149:CYS:O	1:A:280:GLU:HA	2.19	0.43
1:A:161:PHE:O	1:A:165:VAL:HG23	2.19	0.43
1:B:278:GLN:HB2	2:B:357:HOH:O	2.19	0.43
1:B:150:ILE:O	1:B:150:ILE:HG13	2.17	0.42
1:B:263:LYS:HE2	1:B:263:LYS:N	2.33	0.42
1:B:274:GLY:HA2	1:B:285:ILE:O	2.19	0.42
1:A:227:PHE:HA	1:A:230:MET:HE3	2.02	0.42
1:B:260:GLN:HA	1:B:263:LYS:HG2	2.00	0.42
1:B:309:MET:HA	1:B:314:GLU:O	2.20	0.42
1:A:288:THR:HB	1:A:289:PRO:CD	2.49	0.42
1:B:212:ALA:HA	1:B:213:PRO:C	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:GLN:OE1	1:B:238:MET:HG2	2.19	0.42
1:A:245:LEU:O	1:A:249:ASN:ND2	2.53	0.41
1:B:237:ILE:HG23	1:B:238:MET:N	2.34	0.41
1:B:328:ILE:HG23	1:B:329:GLU:HG2	2.01	0.41
1:A:244:LEU:HD13	1:A:255:LEU:CD2	2.49	0.41
1:B:156:ASN:ND2	2:B:349:HOH:O	2.53	0.41
1:B:215:GLU:H	1:B:215:GLU:CD	2.28	0.41
1:B:231:PHE:HB2	1:B:237:ILE:CD1	2.49	0.41
1:A:257:PHE:HA	2:A:368:HOH:O	2.20	0.41
1:A:186:ILE:HD13	1:A:208:SER:HA	2.03	0.41
1:B:255:LEU:HD22	1:B:256:THR:N	2.35	0.41
1:A:193:PHE:CE1	1:A:234:TYR:HB2	2.55	0.41
1:B:168:ILE:HD11	1:B:234:TYR:CD2	2.56	0.41
1:B:185:LEU:HD23	1:B:185:LEU:O	2.21	0.41
1:B:179:THR:HB	1:B:218:THR:CG2	2.51	0.41
1:B:332:LEU:HD11	1:B:334:ALA:HB3	2.02	0.41
1:A:341:MET:HA	1:A:341:MET:CE	2.44	0.40
1:B:277:ASP:HB2	1:B:285:ILE:CD1	2.50	0.40
1:B:181:MET:O	1:B:182:ARG:C	2.65	0.40
1:A:272:ILE:CD1	1:A:328:ILE:HG21	2.50	0.40
1:B:283:CYS:HB2	1:B:296:TRP:NE1	2.36	0.40
1:A:181:MET:CE	1:A:218:THR:HG22	2.52	0.40
1:A:323:PHE:HE1	1:A:328:ILE:HD11	1.86	0.40
1:B:308:LEU:HB2	1:B:316:TYR:HB2	2.03	0.40

There are no symmetry-related clashes.

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	A	179/221 (81%)	174 (97%)	4 (2%)	1 (1%)	21 25

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	189/221 (86%)	181 (96%)	6 (3%)	2 (1%)	11	12
All	All	368/442 (83%)	355 (96%)	10 (3%)	3 (1%)	16	19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	177	ASP
1	A	177	ASP
1	B	182	ARG

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	A	161/198 (81%)	145 (90%)	16 (10%)	7	9
1	B	168/198 (85%)	148 (88%)	20 (12%)	5	5
All	All	329/396 (83%)	293 (89%)	36 (11%)	6	7

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	ASN
1	A	187	GLU
1	A	197	GLU
1	A	201	GLN
1	A	215	GLU
1	A	221	LYS
1	A	224	LEU
1	A	237	ILE
1	A	244	LEU
1	A	246	LEU
1	A	251	LYS
1	A	253	HIS
1	A	272	ILE

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Mol	Chain	Res	Type
1	A	278	GLN
1	A	290	THR
1	A	321	LYS
1	B	152	ASN
1	B	157	GLU
1	B	158	GLN
1	B	162	ARG
1	B	180	VAL
1	B	185	LEU
1	B	197	GLU
1	B	201	GLN
1	B	207	ILE
1	B	211	MET
1	B	215	GLU
1	B	224	LEU
1	B	237	ILE
1	B	238	MET
1	B	244	LEU
1	B	246	LEU
1	B	255	LEU
1	B	256	THR
1	B	314	GLU
1	B	332	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	152	ASN
1	A	156	ASN
1	A	158	GLN
1	A	249	ASN
1	A	253	HIS
1	A	260	GLN
1	A	279	ASN
1	B	152	ASN
1	B	158	GLN
1	B	217	GLN
1	B	278	GLN
1	B	312	ASN

5.3.3 RNA [i](#)

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](#)

There are no ligands in this entry.

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	A	185/221 (83%)	0.24	1 (0%) 87 88	41, 58, 92, 115	0
1	B	193/221 (87%)	0.37	11 (5%) 29 32	41, 66, 104, 134	0
All	All	378/442 (85%)	0.31	12 (3%) 50 53	41, 62, 99, 134	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	265	ALA	4.0
1	B	330	THR	3.5
1	B	344	HIS	3.2
1	B	342	HIS	2.7
1	B	156	ASN	2.4
1	B	147	SER	2.4
1	B	343	HIS	2.3
1	B	170	PHE	2.2
1	A	345	HIS	2.2
1	B	325	MET	2.1
1	B	163	ALA	2.1
1	B	271	SER	2.1

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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.