



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:33 PM GMT

PDB ID : 4TQ0
Title : Crystal structure of human ATG5-ATG16N69
Authors : Kim, J.H.; Hong, S.B.; Song, H.K.
Deposited on : 2014-06-10
Resolution : 2.70 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

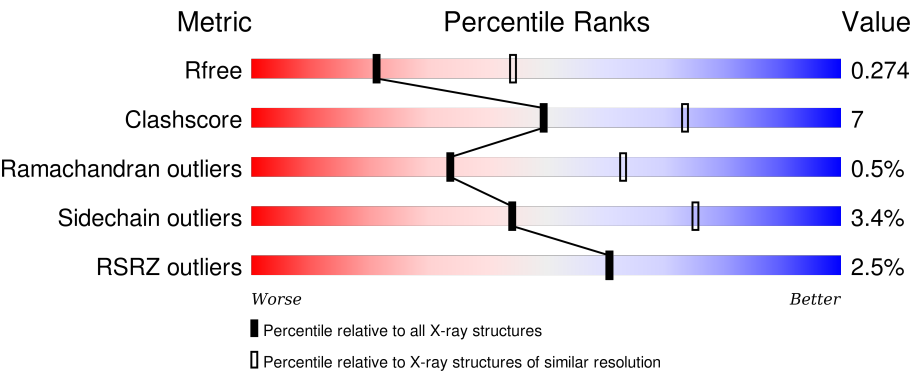
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	289	<div><div>%</div><div><div></div><div>71%</div><div>13%</div><div>•</div><div>16%</div></div></div>
1	C	289	<div><div>2%</div><div><div></div><div>67%</div><div>18%</div><div>•</div><div>14%</div></div></div>
1	E	289	<div><div>4%</div><div><div></div><div>67%</div><div>15%</div><div>•</div><div>16%</div></div></div>
2	B	69	<div><div>%</div><div><div></div><div>49%</div><div>7%</div><div>•</div><div>42%</div></div></div>
2	D	69	<div><div></div><div><div></div><div>46%</div><div>7%</div><div></div><div>46%</div></div></div>

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Mol	Chain	Length	Quality of chain
2	F	69	<div><div></div><div></div><div></div><div>43%</div><div>14%</div><div>42%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7219 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 Autophagy protein 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	244	Total	C	N	O	S	0	0	0
			1999	1306	329	353	11			
1	C	248	Total	C	N	O	S	0	0	0
			2033	1327	333	362	11			
1	E	243	Total	C	N	O	S	0	0	0
			1987	1294	330	352	11			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP Q9H1Y0
A	-12	GLY	-	expression tag	UNP Q9H1Y0
A	-11	SER	-	expression tag	UNP Q9H1Y0
A	-10	SER	-	expression tag	UNP Q9H1Y0
A	-9	HIS	-	expression tag	UNP Q9H1Y0
A	-8	HIS	-	expression tag	UNP Q9H1Y0
A	-7	HIS	-	expression tag	UNP Q9H1Y0
A	-6	HIS	-	expression tag	UNP Q9H1Y0
A	-5	HIS	-	expression tag	UNP Q9H1Y0
A	-4	HIS	-	expression tag	UNP Q9H1Y0
A	-3	SER	-	expression tag	UNP Q9H1Y0
A	-2	GLN	-	expression tag	UNP Q9H1Y0
A	-1	GLY	-	expression tag	UNP Q9H1Y0
A	0	SER	-	expression tag	UNP Q9H1Y0
C	-13	MET	-	expression tag	UNP Q9H1Y0
C	-12	GLY	-	expression tag	UNP Q9H1Y0
C	-11	SER	-	expression tag	UNP Q9H1Y0
C	-10	SER	-	expression tag	UNP Q9H1Y0
C	-9	HIS	-	expression tag	UNP Q9H1Y0
C	-8	HIS	-	expression tag	UNP Q9H1Y0
C	-7	HIS	-	expression tag	UNP Q9H1Y0
C	-6	HIS	-	expression tag	UNP Q9H1Y0
C	-5	HIS	-	expression tag	UNP Q9H1Y0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	HIS	-	expression tag	UNP Q9H1Y0
C	-3	SER	-	expression tag	UNP Q9H1Y0
C	-2	GLN	-	expression tag	UNP Q9H1Y0
C	-1	GLY	-	expression tag	UNP Q9H1Y0
C	0	SER	-	expression tag	UNP Q9H1Y0
E	-13	MET	-	expression tag	UNP Q9H1Y0
E	-12	GLY	-	expression tag	UNP Q9H1Y0
E	-11	SER	-	expression tag	UNP Q9H1Y0
E	-10	SER	-	expression tag	UNP Q9H1Y0
E	-9	HIS	-	expression tag	UNP Q9H1Y0
E	-8	HIS	-	expression tag	UNP Q9H1Y0
E	-7	HIS	-	expression tag	UNP Q9H1Y0
E	-6	HIS	-	expression tag	UNP Q9H1Y0
E	-5	HIS	-	expression tag	UNP Q9H1Y0
E	-4	HIS	-	expression tag	UNP Q9H1Y0
E	-3	SER	-	expression tag	UNP Q9H1Y0
E	-2	GLN	-	expression tag	UNP Q9H1Y0
E	-1	GLY	-	expression tag	UNP Q9H1Y0
E	0	SER	-	expression tag	UNP Q9H1Y0

- Molecule 2 is a protein called Autophagy-related protein 16-1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	40	Total	C	N	O	0	0	0
			354	221	73	60			
2	D	37	Total	C	N	O	0	0	0
			326	205	66	55			
2	F	40	Total	C	N	O	0	0	0
			358	224	74	60			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	43	Total	O	0	0
			43	43		
3	B	4	Total	O	0	0
			4	4		
3	C	54	Total	O	0	0
			54	54		
3	D	6	Total	O	0	0
			6	6		
3	E	49	Total	O	0	0
			49	49		

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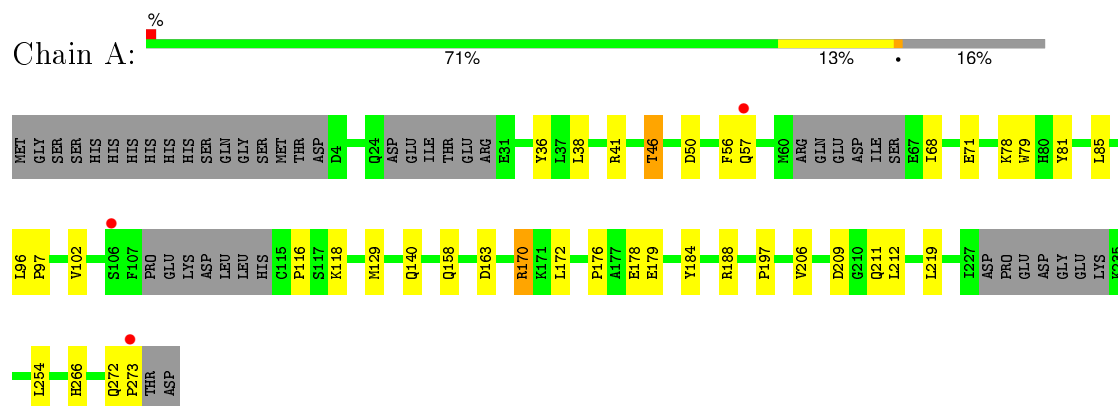
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	F	6	Total	O	0	0
			6	6		

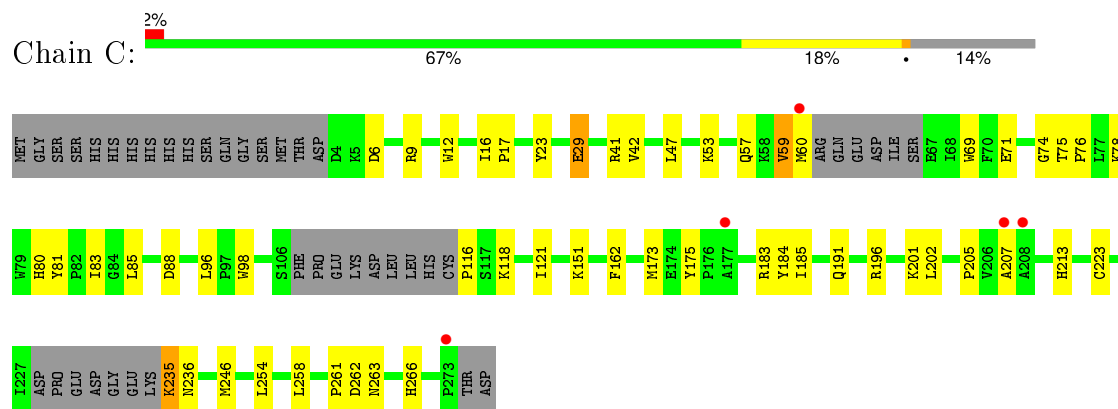
3 Residue-property plots [i](#)

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.

• Molecule 1: Autophagy protein 5



• Molecule 1: Autophagy protein 5

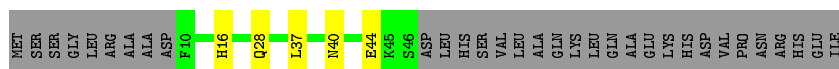




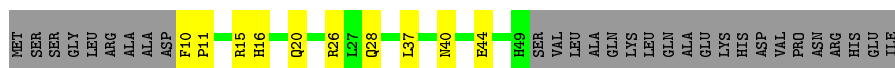
- Molecule 2: Autophagy-related protein 16-1



- Molecule 2: Autophagy-related protein 16-1



- Molecule 2: Autophagy-related protein 16-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	93.09 Å 93.09 Å 245.58 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.09 – 2.70 37.09 – 2.70	Depositor EDS
% Data completeness (in resolution range)	92.3 (37.09-2.70) 92.3 (37.09-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.17 (at 2.68 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.203 , 0.274 0.202 , 0.274	Depositor DCC
R_{free} test set	1829 reflections (6.46%)	DCC
Wilson B-factor (Å ²)	27.0	Xtriage
Anisotropy	0.066	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 37.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 28321 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7219	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.16% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.54	0/2060	0.66	1/2800 (0.0%)
1	C	0.53	0/2094	0.66	1/2845 (0.0%)
1	E	0.47	0/2047	0.67	0/2782
2	B	0.56	0/360	0.62	0/481
2	D	0.54	0/331	0.74	0/441
2	F	0.52	0/364	0.70	0/485
All	All	0.52	0/7256	0.67	2/9834 (0.0%)

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	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	C	121	ILE	CG1-CB-CG2	-6.63	96.81	111.40
1	A	188	ARG	NE-CZ-NH2	-5.24	117.68	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	57	GLN	Peptide

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	1999	0	1918	24	0
1	C	2033	0	1970	35	0
1	E	1987	0	1916	35	0
2	B	354	0	353	5	0
2	D	326	0	331	3	0
2	F	358	0	364	8	0
3	A	43	0	0	2	0
3	B	4	0	0	1	0
3	C	54	0	0	1	0
3	D	6	0	0	0	0
3	E	49	0	0	5	0
3	F	6	0	0	2	0
All	All	7219	0	6852	98	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 7.

All (98) 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)
2:F:10:PHE:O	2:F:15:ARG:NH1	2.14	0.80
1:A:176:PRO:HG2	1:A:179:GLU:HG3	1.69	0.75
2:F:28:GLN:NE2	3:F:104:HOH:O	2.13	0.73
1:E:199:ILE:HD11	1:E:222:VAL:HG12	1.71	0.71
1:C:41:ARG:NH2	1:C:88:ASP:OD1	2.24	0.70
1:E:211:GLN:O	3:E:331:HOH:O	2.08	0.69
1:E:240:ILE:HD12	1:E:245:PRO:HG2	1.74	0.69
1:C:29:GLU:H	1:C:59:VAL:HG12	1.60	0.66
1:E:188:ARG:NH1	3:E:324:HOH:O	2.31	0.63
1:E:191:GLN:NE2	1:E:223:CYS:SG	2.73	0.61
1:A:118:LYS:HB3	1:A:118:LYS:HZ2	1.66	0.61
1:A:46:THR:OG1	3:A:308:HOH:O	2.14	0.61
1:A:140:GLN:HE22	1:C:201:LYS:HE3	1.64	0.61
1:E:85:LEU:HD13	1:E:266:HIS:HB3	1.81	0.61
1:A:184:TYR:OH	1:C:183:ARG:NH2	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:209:ASP:N	1:E:210:GLY:HA2	2.16	0.59
1:E:116:PRO:O	1:E:117:SER:HB3	2.03	0.58
1:C:85:LEU:HD13	1:C:266:HIS:HB3	1.84	0.58
1:E:93:SER:HB2	1:E:273:PRO:HB3	1.86	0.58
1:E:273:PRO:O	1:E:274:THR:OG1	2.21	0.57
1:C:29:GLU:N	1:C:59:VAL:HG12	2.21	0.56
1:E:176:PRO:HG2	1:E:179:GLU:HG3	1.87	0.56
1:A:85:LEU:HD13	1:A:266:HIS:HB3	1.88	0.55
1:E:141:VAL:HG23	1:E:172:LEU:HD23	1.90	0.54
2:D:40:ASN:O	2:D:44:GLU:HG2	2.09	0.52
1:C:42:VAL:HB	1:C:261:PRO:CD	2.39	0.52
1:E:272:GLN:NE2	3:E:330:HOH:O	2.41	0.52
2:F:26:ARG:HD3	3:F:103:HOH:O	2.10	0.52
1:E:41:ARG:NH1	2:F:28:GLN:HE22	2.07	0.51
1:E:189:ILE:HB	1:E:199:ILE:HG12	1.90	0.51
1:E:209:ASP:N	1:E:209:ASP:OD1	2.43	0.51
1:E:193:THR:HB	1:E:195:GLU:HG2	1.92	0.50
2:B:37:LEU:HD13	3:B:103:HOH:O	2.11	0.50
1:C:69:TRP:CZ3	1:C:71:GLU:HB2	2.47	0.49
1:E:42:VAL:HB	1:E:261:PRO:CD	2.42	0.49
1:A:96:LEU:HD11	2:B:28:GLN:HG2	1.95	0.48
2:F:16:HIS:O	2:F:20:GLN:HG2	2.14	0.47
1:E:76:PRO:O	3:E:332:HOH:O	2.20	0.47
1:C:74:GLY:HA3	1:E:147:LYS:HE2	1.96	0.47
1:C:59:VAL:O	1:C:60:MET:HB2	2.15	0.46
1:C:6:ASP:OD1	1:C:9:ARG:NH2	2.49	0.46
1:C:53:LYS:HZ2	1:C:57:GLN:NE2	2.12	0.46
1:C:262:ASP:O	1:C:263:ASN:HB2	2.16	0.46
1:A:38:LEU:HD11	2:B:33:GLU:HG3	1.98	0.46
1:E:58:LYS:HB3	1:E:58:LYS:HE2	1.83	0.46
1:A:78:LYS:HB2	1:A:81:TYR:CD2	2.50	0.46
1:E:184:TYR:CE1	1:E:202:LEU:HB3	2.51	0.46
2:F:40:ASN:O	2:F:44:GLU:HG3	2.15	0.46
1:C:57:GLN:C	1:C:59:VAL:H	2.19	0.46
1:C:185:ILE:HD11	1:C:205:PRO:HG3	1.97	0.45
1:A:178:GLU:HG2	1:A:179:GLU:HG2	1.99	0.45
1:A:56:PHE:CE2	1:A:102:VAL:HG21	2.52	0.45
1:C:175:TYR:OH	3:C:331:HOH:O	2.17	0.45
1:E:71:GLU:OE2	1:E:103:HIS:HE1	1.99	0.45
1:E:141:VAL:CG2	1:E:172:LEU:HD23	2.46	0.45
1:C:184:TYR:CZ	1:C:202:LEU:HD13	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:78:LYS:HD2	1:C:81:TYR:CE2	2.52	0.45
1:A:79:TRP:CD2	1:A:116:PRO:HG2	2.52	0.45
2:F:10:PHE:HA	2:F:11:PRO:HD3	1.85	0.44
1:C:23:TYR:OH	1:E:119:ASP:OD2	2.30	0.44
1:A:41:ARG:NH1	2:B:28:GLN:OE1	2.51	0.44
1:A:163:ASP:OD1	1:A:163:ASP:N	2.50	0.44
1:E:175:TYR:HB2	1:E:176:PRO:HD2	1.99	0.44
1:E:133:ASP:OD1	1:E:137:HIS:ND1	2.48	0.44
1:C:207:ALA:HA	1:C:213:HIS:CE1	2.53	0.43
1:A:197:PRO:HG3	1:C:80:HIS:CD2	2.54	0.43
1:C:17:PRO:HG2	1:C:96:LEU:HD22	2.01	0.43
1:C:16:ILE:HG12	1:C:41:ARG:HG2	2.00	0.43
1:A:272:GLN:HA	1:A:273:PRO:HA	1.85	0.42
1:A:36:TYR:HE1	1:A:97:PRO:HG2	1.83	0.42
1:C:151:LYS:HA	1:C:151:LYS:HD3	1.89	0.42
1:A:170:ARG:NH2	3:A:302:HOH:O	2.43	0.42
1:C:16:ILE:HD11	1:C:41:ARG:CZ	2.49	0.42
1:C:53:LYS:NZ	1:C:57:GLN:NE2	2.67	0.42
1:A:254:LEU:HD23	1:A:254:LEU:HA	1.81	0.42
1:E:56:PHE:O	1:E:60:MET:HG2	2.20	0.41
1:C:235:LYS:HB2	1:C:236:ASN:H	1.53	0.41
1:E:207:ALA:N	3:E:331:HOH:O	2.23	0.41
1:C:191:GLN:NE2	1:C:223:CYS:SG	2.93	0.41
1:C:246:MET:HG3	2:D:16:HIS:CD2	2.55	0.41
1:C:12:TRP:CG	1:C:162:PHE:HD2	2.39	0.41
1:A:129:MET:HE1	1:A:172:LEU:HD13	2.02	0.41
1:C:254:LEU:HD23	1:C:254:LEU:HA	1.95	0.41
1:C:185:ILE:CD1	1:C:205:PRO:HG3	2.50	0.41
1:A:206:VAL:HG12	1:A:212:LEU:HD23	2.02	0.41
1:A:118:LYS:HZ1	1:A:158:GLN:NE2	2.19	0.41
1:E:22:LEU:HB2	1:E:56:PHE:CE1	2.56	0.41
1:E:16:ILE:HG12	1:E:41:ARG:HG2	2.03	0.41
1:E:272:GLN:OE1	1:E:273:PRO:HD2	2.21	0.41
1:C:75:THR:HA	1:C:76:PRO:HD3	1.97	0.41
1:E:117:SER:O	1:E:121:ILE:HG13	2.21	0.40
1:E:41:ARG:HH11	2:F:28:GLN:HE22	1.69	0.40
2:B:10:PHE:HA	2:B:11:PRO:HD2	1.83	0.40
1:A:219:LEU:HA	1:A:219:LEU:HD23	1.90	0.40
1:C:96:LEU:HD11	2:D:28:GLN:HG2	2.03	0.40
1:E:187:PHE:O	1:E:188:ARG:HG3	2.22	0.40
1:A:211:GLN:HE21	1:A:212:LEU:H	1.70	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:ILE:HG23	1:C:98:TRP:HZ3	1.86	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	234/289 (81%)	228 (97%)	6 (3%)	0	100	100
1	C	240/289 (83%)	236 (98%)	4 (2%)	0	100	100
1	E	235/289 (81%)	227 (97%)	4 (2%)	4 (2%)	11	29
2	B	38/69 (55%)	37 (97%)	1 (3%)	0	100	100
2	D	35/69 (51%)	35 (100%)	0	0	100	100
2	F	38/69 (55%)	37 (97%)	1 (3%)	0	100	100
All	All	820/1074 (76%)	800 (98%)	16 (2%)	4 (0%)	34	63

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	58	LYS
1	E	117	SER
1	E	208	ALA
1	E	273	PRO

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	214/266 (80%)	208 (97%)	6 (3%)	51	81
1	C	220/266 (83%)	211 (96%)	9 (4%)	37	69
1	E	213/266 (80%)	206 (97%)	7 (3%)	45	76
2	B	37/63 (59%)	35 (95%)	2 (5%)	27	56
2	D	34/63 (54%)	33 (97%)	1 (3%)	50	80
2	F	38/63 (60%)	37 (97%)	1 (3%)	54	83
All	All	756/987 (77%)	730 (97%)	26 (3%)	44	75

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

Mol	Chain	Res	Type
1	A	46	THR
1	A	50	ASP
1	A	68	ILE
1	A	71	GLU
1	A	170	ARG
1	A	209	ASP
2	B	37	LEU
2	B	44	GLU
1	C	29	GLU
1	C	47	LEU
1	C	59	VAL
1	C	116	PRO
1	C	118	LYS
1	C	173	MET
1	C	196	ARG
1	C	235	LYS
1	C	258	LEU
2	D	37	LEU
1	E	30	ARG
1	E	68	ILE
1	E	117	SER
1	E	193	THR
1	E	199	ILE
1	E	202	LEU
1	E	209	ASP
2	F	37	LEU

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

sidechains are listed below:

Mol	Chain	Res	Type
1	A	140	GLN
1	A	158	GLN
1	A	211	GLN
1	C	57	GLN
1	C	191	GLN
1	E	191	GLN
2	F	28	GLN

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 carbohydrates 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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	244/289 (84%)	-0.44	3 (1%) 81 81	9, 21, 53, 74	0
1	C	248/289 (85%)	-0.33	5 (2%) 68 69	10, 24, 53, 72	0
1	E	243/289 (84%)	-0.02	12 (4%) 33 32	15, 31, 70, 103	0
2	B	40/69 (57%)	-0.25	1 (2%) 61 61	18, 29, 58, 71	0
2	D	37/69 (53%)	-0.39	0 100 100	15, 29, 48, 51	0
2	F	40/69 (57%)	-0.42	0 100 100	14, 26, 52, 64	0
All	All	852/1074 (79%)	-0.28	21 (2%) 61 61	9, 26, 60, 103	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	177	ALA	6.1
1	E	195	GLU	5.7
1	E	138	LYS	4.7
1	C	208	ALA	4.2
1	E	194	THR	4.0
1	E	210	GLY	3.9
1	E	208	ALA	3.8
1	E	139	SER	3.3
1	E	193	THR	2.9
1	A	273	PRO	2.9
1	E	207	ALA	2.8
1	A	106	SER	2.6
1	C	273	PRO	2.5
1	E	211	GLN	2.4
1	E	180	ASN	2.4
1	A	57	GLN	2.4
1	C	207	ALA	2.3
1	C	60	MET	2.2
2	B	10	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	178	GLU	2.1
1	E	25	ASP	2.0

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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.