



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 11:07 PM GMT

PDB ID : 3VX8  
Title : Crystal structure of Arabidopsis thaliana Atg7NTD-Atg3 complex  
Authors : Matoba, K.; Fujioka, Y.; Noda, N.N.  
Deposited on : 2012-09-11  
Resolution : 3.11 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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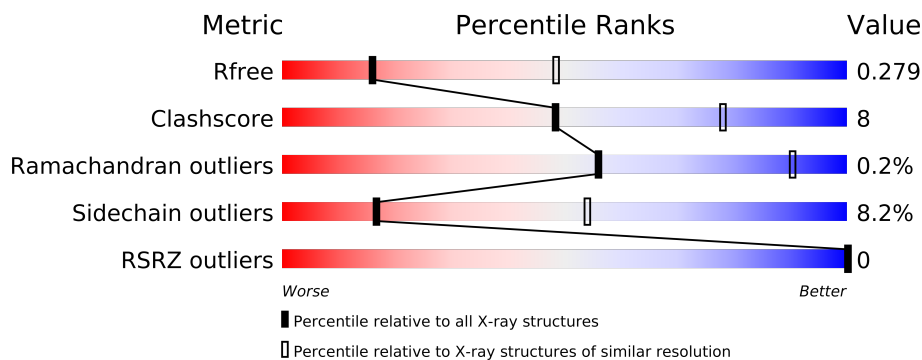
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
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 3.11 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))
$R_{free}$	66092	1466 (3.20-3.04)
Clashscore	79885	1067 (3.18-3.06)
Ramachandran outliers	78287	1034 (3.18-3.06)
Sidechain outliers	78261	1034 (3.18-3.06)
RSRZ outliers	66119	1468 (3.20-3.04)

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.

Mol	Chain	Length	Quality of chain
1	A	323	
1	D	323	
2	B	292	
2	C	292	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7608 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 Ubiquitin-like modifier-activating enzyme atg7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	300	Total	C	N	O	S	0	0	0
			2380	1533	396	444	7			
1	A	305	Total	C	N	O	S	0	0	0
			2418	1557	403	451	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	3	GLY	-	EXPRESSION TAG	UNP Q94CD5
D	4	PRO	-	EXPRESSION TAG	UNP Q94CD5
D	5	HIS	-	EXPRESSION TAG	UNP Q94CD5
D	6	MET	-	EXPRESSION TAG	UNP Q94CD5
A	3	GLY	-	EXPRESSION TAG	UNP Q94CD5
A	4	PRO	-	EXPRESSION TAG	UNP Q94CD5
A	5	HIS	-	EXPRESSION TAG	UNP Q94CD5
A	6	MET	-	EXPRESSION TAG	UNP Q94CD5

- Molecule 2 is a protein called Autophagy-related protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	175	Total	C	N	O	S	0	0	0
			1416	909	238	259	10			
2	C	172	Total	C	N	O	S	0	0	0
			1391	892	233	256	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	22	GLY	-	EXPRESSION TAG	UNP Q0WWQ1
B	23	PRO	-	EXPRESSION TAG	UNP Q0WWQ1
B	24	HIS	-	EXPRESSION TAG	UNP Q0WWQ1
B	25	MET	-	EXPRESSION TAG	UNP Q0WWQ1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	22	GLY	-	EXPRESSION TAG	UNP Q0WWQ1
C	23	PRO	-	EXPRESSION TAG	UNP Q0WWQ1
C	24	HIS	-	EXPRESSION TAG	UNP Q0WWQ1
C	25	MET	-	EXPRESSION TAG	UNP Q0WWQ1

- Molecule 3 is water.

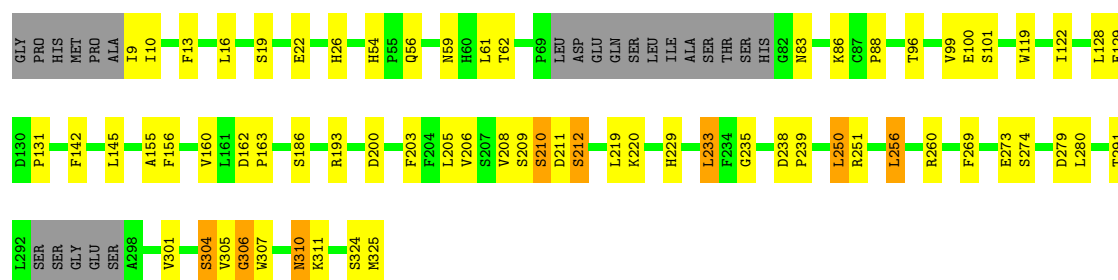
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 1 1	0	0
3	B	1	Total O 1 1	0	0
3	C	1	Total O 1 1	0	0

### 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.

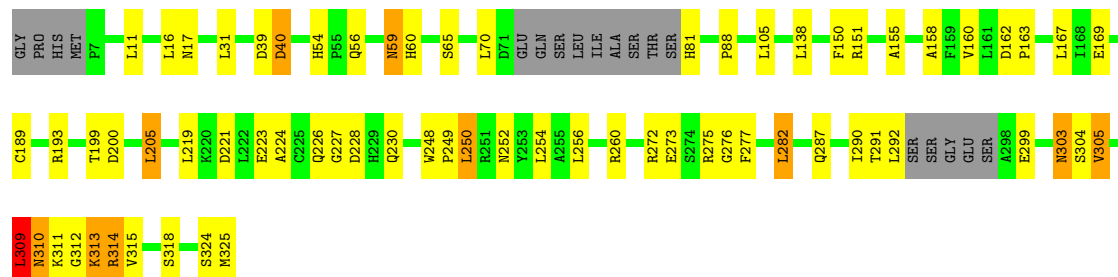
- Molecule 1: Ubiquitin-like modifier-activating enzyme atg7

Chain D:



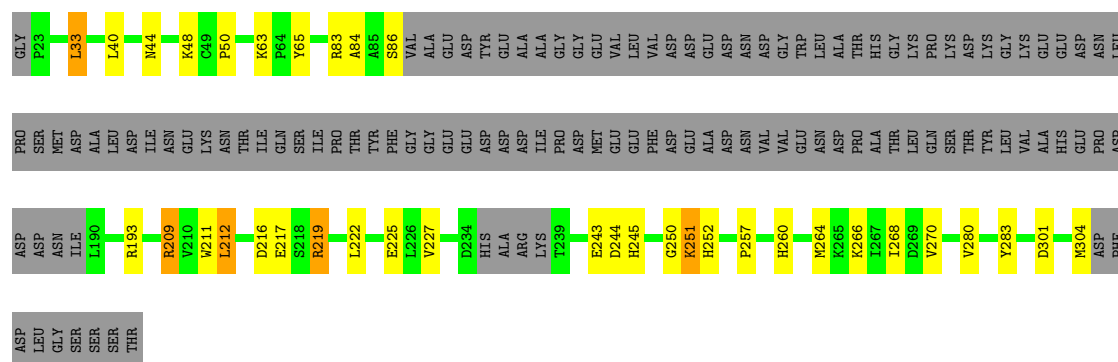
- Molecule 1: Ubiquitin-like modifier-activating enzyme atg7

Chain A:



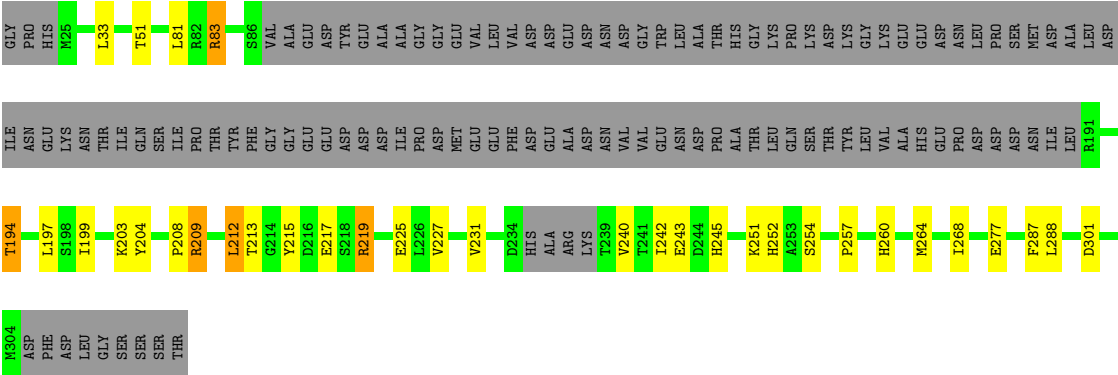
- Molecule 2: Autophagy-related protein 3

Chain B:



● Molecule 2: Autophagy-related protein 3

Chain C:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.52Å 132.68Å 102.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.85 – 3.11 34.85 – 3.11	Depositor EDS
% Data completeness (in resolution range)	99.1 (34.85-3.11) 99.2 (34.85-3.11)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.69 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.226 , 0.278 0.226 , 0.279	Depositor DCC
$R_{free}$ test set	1295 reflections (5.39%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.6	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , -1.8	EDS
Estimated twinning fraction	0.018 for l,-k,h	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 25339 reflections	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	7608	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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	0.50	0/2487	0.60	2/3383 (0.1%)
1	D	0.39	0/2447	0.54	0/3328
2	B	0.28	0/1452	0.51	0/1966
2	C	0.28	0/1425	0.49	0/1929
All	All	0.40	0/7811	0.55	2/10606 (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	A	227	GLY	N-CA-C	6.10	128.35	113.10
1	A	309	LEU	CA-CB-CG	5.55	128.07	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	305	VAL	Peptide

### 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	A	2418	0	2358	47	0
1	D	2380	0	2323	45	0
2	B	1416	0	1412	15	0
2	C	1391	0	1386	17	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
All	All	7608	0	7479	118	0

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 8.

All (118) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:210:SER:HB2	1:D:229:HIS:O	1.55	1.06
1:A:228:ASP:OD1	1:A:230:GLN:HB3	1.59	1.03
1:D:304:SER:O	1:D:305:VAL:HG23	1.65	0.97
2:B:193:ARG:HD2	2:B:222:LEU:HD11	1.51	0.92
1:A:230:GLN:HG3	1:A:230:GLN:O	1.72	0.89
1:D:209:SER:C	1:D:211:ASP:H	1.73	0.88
1:D:162:ASP:HB2	1:D:163:PRO:HD3	1.55	0.88
1:D:304:SER:C	1:D:305:VAL:HG23	1.96	0.85
1:D:209:SER:OG	1:D:211:ASP:HB3	1.77	0.84
1:D:209:SER:OG	1:D:211:ASP:CB	2.30	0.80
1:D:304:SER:O	1:D:305:VAL:CG2	2.30	0.79
1:A:162:ASP:HB3	1:A:163:PRO:HD3	1.64	0.78
1:A:291:THR:O	1:A:292:LEU:HB2	1.87	0.72
1:A:309:LEU:HD13	1:A:312:GLY:C	2.10	0.72
1:D:209:SER:C	1:D:211:ASP:N	2.43	0.69
1:A:54:HIS:HD2	1:A:56:GLN:H	1.41	0.66
1:A:309:LEU:CD1	1:A:312:GLY:O	2.43	0.66
1:A:272:ARG:HD3	1:A:282:LEU:HD22	1.78	0.65
1:A:230:GLN:CG	1:A:230:GLN:O	2.44	0.65
1:A:221:ASP:O	1:A:224:ALA:HB3	1.97	0.64
1:D:83:ASN:HB3	1:D:86:LYS:HB2	1.79	0.64
1:D:209:SER:O	1:D:211:ASP:N	2.30	0.64
1:A:193:ARG:O	1:A:199:THR:HG23	1.97	0.63
1:A:54:HIS:CD2	1:A:56:GLN:H	2.16	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:13:PHE:HE1	1:D:250:LEU:HB3	1.63	0.63
1:A:228:ASP:OD1	1:A:230:GLN:CB	2.44	0.61
2:C:242:ILE:HG22	2:C:251:LYS:HG2	1.83	0.60
1:A:11:LEU:HD13	1:A:167:LEU:HD13	1.84	0.60
2:C:197:LEU:HD22	2:C:212:LEU:HB3	1.83	0.59
2:B:264:MET:O	2:B:268:ILE:HG12	2.02	0.59
1:D:22:GLU:OE1	1:D:324:SER:HB3	2.03	0.58
1:D:211:ASP:O	1:D:212:SER:HB2	2.03	0.58
1:A:310:ASN:O	1:A:311:LYS:C	2.37	0.58
2:C:51:THR:HG21	2:C:301:ASP:HB2	1.85	0.56
1:A:309:LEU:CD1	1:A:312:GLY:C	2.73	0.56
1:A:310:ASN:C	1:A:312:GLY:N	2.57	0.56
1:A:310:ASN:O	1:A:312:GLY:N	2.39	0.55
1:D:129:GLU:HA	1:D:220:LYS:HE2	1.89	0.55
2:C:194:THR:HG23	2:C:215:TYR:HB2	1.89	0.55
1:A:228:ASP:CG	1:A:230:GLN:HB3	2.26	0.54
2:B:212:LEU:HD13	2:B:227:VAL:HG11	1.89	0.54
2:B:63:LYS:HD3	2:B:65:TYR:OH	2.08	0.54
1:D:54:HIS:CD2	1:D:56:GLN:H	2.26	0.53
2:B:251:LYS:HG3	2:C:204:TYR:HB2	1.90	0.53
1:A:59:ASN:HD22	1:A:60:HIS:H	1.56	0.53
2:C:83:ARG:HA	2:C:83:ARG:HE	1.72	0.53
2:C:199:ILE:HD11	2:C:288:LEU:HD21	1.91	0.52
1:A:59:ASN:ND2	1:A:60:HIS:H	2.07	0.52
2:C:257:PRO:HA	2:C:260:HIS:CD2	2.45	0.52
1:D:205:LEU:HD23	1:D:219:LEU:HA	1.91	0.52
1:A:275:ARG:HA	2:B:301:ASP:OD2	2.10	0.52
2:B:245:HIS:HB2	2:B:252:HIS:CD2	2.45	0.52
2:B:257:PRO:HA	2:B:260:HIS:CD2	2.45	0.51
1:D:211:ASP:OD1	1:D:211:ASP:O	2.30	0.50
1:D:54:HIS:HD2	1:D:56:GLN:H	1.59	0.50
2:C:231:VAL:HG11	2:C:240:VAL:HG21	1.93	0.50
1:A:291:THR:O	1:A:292:LEU:CB	2.58	0.50
1:A:273:GLU:HB2	1:A:277:PHE:O	2.13	0.49
1:A:250:LEU:HD22	1:A:254:LEU:HG	1.93	0.49
1:D:162:ASP:HB2	1:D:163:PRO:CD	2.36	0.49
1:D:209:SER:OG	1:D:211:ASP:HB2	2.10	0.48
1:D:131:PRO:HB2	1:D:203:PHE:HB2	1.95	0.48
1:D:193:ARG:NH1	1:D:273:GLU:OE2	2.46	0.48
1:D:205:LEU:CD2	1:D:219:LEU:HA	2.44	0.47
1:D:88:PRO:HD2	1:D:200:ASP:HB2	1.97	0.47
1:A:40:ASP:N	1:A:40:ASP:OD1	2.38	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:313:LYS:O	1:A:315:VAL:N	2.48	0.46
1:D:22:GLU:O	1:D:26:HIS:HD2	1.99	0.46
1:D:61:LEU:HD22	1:D:156:PHE:HE1	1.80	0.46
1:A:70:LEU:HD21	1:A:81:HIS:HE1	1.81	0.46
1:A:169:GLU:HG2	1:A:287:GLN:HG2	1.97	0.46
2:C:197:LEU:CD2	2:C:212:LEU:HB3	2.46	0.46
2:C:245:HIS:HB2	2:C:252:HIS:CD2	2.51	0.45
1:D:131:PRO:HB3	1:D:219:LEU:HD12	1.98	0.45
1:D:186:SER:OG	1:D:280:LEU:HD11	2.16	0.44
1:D:206:VAL:HB	1:D:233:LEU:HB2	1.97	0.44
1:D:155:ALA:HA	1:D:306:GLY:O	2.17	0.44
1:A:60:HIS:CE1	2:B:219:ARG:HG3	2.52	0.44
1:A:309:LEU:HD13	1:A:312:GLY:O	2.11	0.44
2:C:212:LEU:HD13	2:C:227:VAL:HG11	2.00	0.44
1:A:276:GLY:HA3	2:B:50:PRO:HD2	1.99	0.44
1:D:238:ASP:HA	1:D:239:PRO:HD3	1.87	0.44
1:A:324:SER:HB3	2:B:250:GLY:HA2	1.99	0.44
1:D:163:PRO:HG2	1:D:291:THR:HB	1.99	0.43
2:C:209:ARG:HD3	2:C:254:SER:OG	2.18	0.43
2:B:84:ALA:C	2:B:86:SER:H	2.22	0.43
1:A:160:VAL:N	1:A:303:ASN:O	2.36	0.43
2:B:266:LYS:O	2:B:270:VAL:HG23	2.18	0.43
1:A:314:ARG:HB3	1:A:314:ARG:HE	1.55	0.43
1:A:303:ASN:OD1	1:A:303:ASN:N	2.51	0.42
1:A:138:LEU:HB3	1:A:155:ALA:HB3	2.01	0.42
1:D:235:GLY:HA2	1:D:269:PHE:HB3	2.01	0.42
1:A:158:ALA:HA	1:A:252:ASN:OD1	2.20	0.42
1:A:205:LEU:CD2	1:A:219:LEU:HA	2.49	0.42
1:A:310:ASN:HD22	1:A:311:LYS:N	2.16	0.42
1:D:19:SER:HB2	2:C:219:ARG:HD2	2.01	0.42
1:A:189:CYS:O	1:A:193:ARG:HG2	2.19	0.41
1:D:119:TRP:CD1	1:D:301:VAL:HG22	2.55	0.41
1:D:310:ASN:C	1:D:310:ASN:HD22	2.23	0.41
2:C:242:ILE:CG2	2:C:251:LYS:HG2	2.48	0.41
1:D:274:SER:HB3	1:D:279:ASP:HB2	2.02	0.41
1:D:26:HIS:HE1	1:D:325:MET:HA	1.86	0.41
1:A:324:SER:O	1:A:325:MET:C	2.58	0.41
2:C:33:LEU:HD12	2:C:208:PRO:HD3	2.03	0.41
1:D:160:VAL:HG22	1:D:251:ARG:HH12	1.84	0.41
1:A:88:PRO:HD2	1:A:200:ASP:HB2	2.03	0.41
1:A:150:PHE:O	1:A:318:SER:HA	2.21	0.41
1:D:96:THR:HB	1:D:101:SER:HB2	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:309:LEU:HD13	1:A:313:LYS:N	2.36	0.40
1:D:155:ALA:HB2	1:D:307:TRP:CZ3	2.57	0.40
1:A:248:TRP:N	1:A:249:PRO:CD	2.84	0.40
2:C:264:MET:HE3	2:C:287:PHE:HA	2.03	0.40
1:D:10:ILE:HG13	1:D:10:ILE:H	1.66	0.40
1:A:11:LEU:HB2	1:A:167:LEU:HB2	2.02	0.40
1:D:122:ILE:HG12	1:D:256:LEU:HG	2.03	0.40
2:B:209:ARG:HD2	2:B:211:TRP:CE2	2.57	0.40
1:D:99:VAL:HG22	1:D:142:PHE:CE2	2.57	0.40
2:B:33:LEU:HG	2:B:283:TYR:CD1	2.57	0.40

There are no symmetry-related clashes.

## 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	A	299/323 (93%)	284 (95%)	14 (5%)	1 (0%)	50	87
1	D	294/323 (91%)	275 (94%)	18 (6%)	1 (0%)	50	87
2	B	169/292 (58%)	164 (97%)	5 (3%)	0	100	100
2	C	166/292 (57%)	161 (97%)	5 (3%)	0	100	100
All	All	928/1230 (75%)	884 (95%)	42 (4%)	2 (0%)	56	90

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	314	ARG
1	D	306	GLY

### 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	A	271/286 (95%)	247 (91%)	24 (9%)	14	48
1	D	267/286 (93%)	250 (94%)	17 (6%)	25	66
2	B	160/259 (62%)	144 (90%)	16 (10%)	11	39
2	C	157/259 (61%)	144 (92%)	13 (8%)	16	54
All	All	855/1090 (78%)	785 (92%)	70 (8%)	17	54

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	9	ILE
1	D	16	LEU
1	D	59	ASN
1	D	62	THR
1	D	100	GLU
1	D	128	LEU
1	D	145	LEU
1	D	208	VAL
1	D	210	SER
1	D	212	SER
1	D	233	LEU
1	D	250	LEU
1	D	256	LEU
1	D	260	ARG
1	D	304	SER
1	D	310	ASN
1	D	311	LYS
1	A	16	LEU
1	A	17	ASN
1	A	31	LEU
1	A	39	ASP
1	A	40	ASP
1	A	59	ASN
1	A	65	SER
1	A	105	LEU
1	A	151	ARG
1	A	205	LEU
1	A	223	GLU
1	A	226	GLN
1	A	250	LEU

*Continued on next page...*

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Mol	Chain	Res	Type
1	A	256	LEU
1	A	260	ARG
1	A	282	LEU
1	A	290	ILE
1	A	299	GLU
1	A	303	ASN
1	A	304	SER
1	A	305	VAL
1	A	309	LEU
1	A	310	ASN
1	A	313	LYS
2	B	33	LEU
2	B	40	LEU
2	B	44	ASN
2	B	48	LYS
2	B	83	ARG
2	B	209	ARG
2	B	212	LEU
2	B	216	ASP
2	B	217	GLU
2	B	219	ARG
2	B	225	GLU
2	B	243	GLU
2	B	244	ASP
2	B	251	LYS
2	B	280	VAL
2	B	304	MET
2	C	81	LEU
2	C	83	ARG
2	C	194	THR
2	C	203	LYS
2	C	209	ARG
2	C	212	LEU
2	C	213	THR
2	C	217	GLU
2	C	219	ARG
2	C	225	GLU
2	C	243	GLU
2	C	268	ILE
2	C	277	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	26	HIS
1	D	54	HIS
1	D	59	ASN
1	D	60	HIS
1	D	83	ASN
1	D	287	GLN
1	D	310	ASN
1	D	323	ASN
1	A	17	ASN
1	A	54	HIS
1	A	59	ASN
1	A	60	HIS
1	A	81	HIS
1	A	85	ASN
1	A	116	ASN
1	A	226	GLN
1	A	230	GLN
1	A	287	GLN
1	A	310	ASN
2	B	44	ASN
2	C	247	HIS

### 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 ⓘ

There are no ligands in this entry.

## 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 ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	305/323 (94%)	-0.31	0 100 100	18, 32, 56, 87	0
1	D	300/323 (92%)	-0.06	0 100 100	33, 56, 85, 108	0
2	B	175/292 (59%)	-0.31	0 100 100	18, 33, 58, 89	0
2	C	172/292 (58%)	-0.18	0 100 100	33, 49, 74, 102	0
All	All	952/1230 (77%)	-0.21	0 100 100	18, 43, 79, 108	0

There are no RSRZ outliers to report.

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

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

### 6.3 Carbohydrates ⓘ

There are no carbohydrates 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.