



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:55 PM UTC

PDB ID : 6CBC / pdb_00006cbc
Title : Crystal structure of an N-terminal fragment of Vps13.
Authors : Kumar, N.; Horenkamp, F.A.; Reinisch, K.M.
Deposited on : 2018-02-02
Resolution : 3.00 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
Mogul : 2022.3.0, CSD as543be (2022)
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

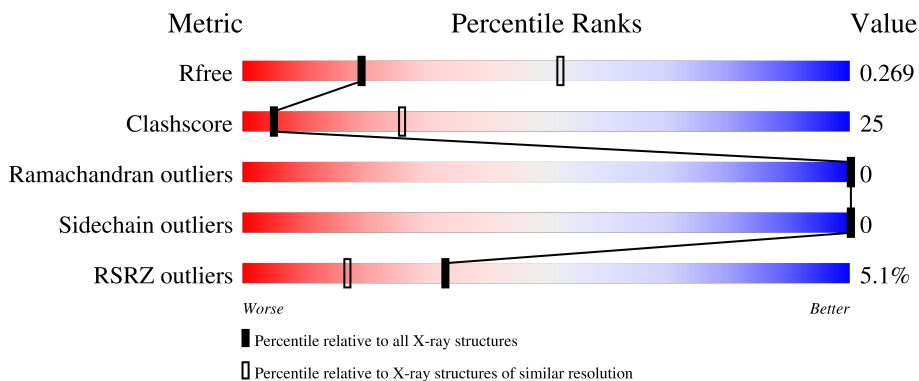
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	336	 3% 45% 27% 26%
1	B	336	 4% 51% 29% 20%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4003 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 Vacuolar protein sorting-associated protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	248	1862	1201	305	351	5	0	0	0
1	B	270	2141	1390	362	384	5	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	initiating methionine	UNP G0S3B8
A	?	-	HIS	deletion	UNP G0S3B8
A	?	-	GLU	deletion	UNP G0S3B8
A	?	-	HIS	deletion	UNP G0S3B8
A	?	-	SER	deletion	UNP G0S3B8
A	?	-	ASP	deletion	UNP G0S3B8
A	?	-	GLN	deletion	UNP G0S3B8
A	220I	HIS	ILE	engineered mutation	UNP G0S3B8
A	227	GLU	PRO	engineered mutation	UNP G0S3B8
A	327	GLY	-	expression tag	UNP G0S3B8
A	328	HIS	-	expression tag	UNP G0S3B8
A	329	HIS	-	expression tag	UNP G0S3B8
A	330	HIS	-	expression tag	UNP G0S3B8
A	331	HIS	-	expression tag	UNP G0S3B8
A	332	HIS	-	expression tag	UNP G0S3B8
A	333	HIS	-	expression tag	UNP G0S3B8
B	1	MSE	-	initiating methionine	UNP G0S3B8
B	?	-	HIS	deletion	UNP G0S3B8
B	?	-	GLU	deletion	UNP G0S3B8
B	?	-	HIS	deletion	UNP G0S3B8
B	?	-	SER	deletion	UNP G0S3B8
B	?	-	ASP	deletion	UNP G0S3B8
B	?	-	GLN	deletion	UNP G0S3B8
B	228A	HIS	ILE	engineered mutation	UNP G0S3B8
B	228B	GLU	PRO	engineered mutation	UNP G0S3B8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	327	GLY	-	expression tag	UNP G0S3B8
B	328	HIS	-	expression tag	UNP G0S3B8
B	329	HIS	-	expression tag	UNP G0S3B8
B	330	HIS	-	expression tag	UNP G0S3B8
B	331	HIS	-	expression tag	UNP G0S3B8
B	332	HIS	-	expression tag	UNP G0S3B8
B	333	HIS	-	expression tag	UNP G0S3B8

HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.93Å 85.98Å 71.03Å 90.00° 93.67° 90.00°	Depositor
Resolution (Å)	48.54 – 3.00 48.54 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (48.54-3.00) 98.4 (48.54-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.09 (at 3.01Å)	Xtrriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.230 , 0.264 0.235 , 0.269	Depositor DCC
R_{free} test set	700 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	72.4	Xtrriage
Anisotropy	0.939	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 72.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4003	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.60% 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

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	0.86	3/1891 (0.2%)	1.09	15/2565 (0.6%)
1	B	0.72	0/2184	0.88	13/2953 (0.4%)
All	All	0.79	3/4075 (0.1%)	0.98	28/5518 (0.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	85	VAL	C-O	-5.14	1.18	1.24
1	A	163	GLU	C-O	-5.11	1.18	1.24
1	A	287	VAL	C-O	-5.02	1.18	1.24

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	84	ASP	N-CA-C	9.60	124.24	111.28
1	B	200	GLN	N-CA-C	8.80	120.96	111.36
1	A	200	GLN	N-CA-C	8.31	120.42	111.36
1	B	77	GLN	N-CA-C	8.00	121.78	110.50
1	A	317	GLN	CA-C-N	7.23	127.20	119.76
1	A	317	GLN	C-N-CA	7.23	127.20	119.76
1	B	39	ASN	N-CA-C	6.91	121.19	111.52
1	B	84	ASP	N-CA-C	6.46	120.00	111.28
1	B	210	SER	N-CA-C	6.34	120.39	111.52
1	A	68	ILE	CB-CA-C	-6.26	103.89	110.53
1	A	284	GLU	N-CA-C	6.15	119.73	111.24
1	A	310	HIS	N-CA-C	6.01	118.61	111.33
1	B	68	ILE	CB-CA-C	-5.86	104.45	110.13
1	A	75	SER	N-CA-C	-5.84	106.80	114.04
1	B	158	ILE	N-CA-C	5.76	116.88	108.58
1	B	158	ILE	CB-CA-C	-5.64	103.92	110.91
1	B	17	TYR	N-CA-C	5.61	120.28	113.38
1	B	224	PRO	N-CA-C	-5.60	104.19	112.26
1	A	309	ARG	N-CA-C	5.48	119.14	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	86	PHE	CA-C-N	-5.37	113.32	122.33
1	A	86	PHE	C-N-CA	-5.37	113.32	122.33
1	A	16	MSE	N-CA-C	5.31	116.75	111.07
1	B	46	ALA	N-CA-C	5.31	117.06	111.28
1	A	318	PRO	N-CA-C	5.30	119.41	111.14
1	A	55	ASN	N-CA-C	-5.28	100.63	109.24
1	B	18	VAL	N-CA-C	5.28	115.91	108.36
1	B	222	ILE	CB-CA-C	-5.23	105.27	111.97
1	A	197	THR	N-CA-C	-5.08	100.23	108.52

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	1862	0	1733	113	0
1	B	2141	0	2126	83	0
All	All	4003	0	3859	193	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 25.

All (193) 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:B:69:PRO:O	1:B:73:LEU:HB2	1.73	0.88
1:A:158:ILE:HD11	1:A:179:LEU:HB3	1.56	0.86
1:B:86:PHE:HB3	1:B:235:PHE:CD2	2.11	0.85
1:A:292:GLN:O	1:A:295:ARG:HG2	1.76	0.84
1:A:66:LEU:HD21	1:A:80:ILE:HD12	1.62	0.82
1:A:158:ILE:HD11	1:A:179:LEU:HD23	1.59	0.81
1:B:159:HIS:CE1	1:B:238:MSE:HG3	2.18	0.79
1:A:69:PRO:HB2	1:A:72:THR:HG23	1.66	0.77
1:A:158:ILE:HG12	1:A:179:LEU:O	1.85	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:ASP:OD2	1:B:192:THR:HB	1.85	0.76
1:A:66:LEU:CD2	1:A:80:ILE:HD12	2.16	0.76
1:B:226:ARG:H	1:B:231:MSE:CE	2.00	0.75
1:B:226:ARG:H	1:B:231:MSE:HE1	1.50	0.75
1:B:226:ARG:N	1:B:231:MSE:CE	2.50	0.75
1:A:85:VAL:CG1	1:A:158:ILE:HG22	2.18	0.74
1:A:14:LEU:HD11	1:A:18:VAL:HG21	1.70	0.74
1:A:158:ILE:CD1	1:A:179:LEU:HB3	2.17	0.74
1:A:158:ILE:CG1	1:A:179:LEU:HB3	2.17	0.73
1:A:89:ALA:HB3	1:A:162:TYR:HD1	1.55	0.71
1:A:228:HIS:NE2	1:A:232:LEU:HD11	2.07	0.69
1:A:77:GLN:HG2	1:A:150:ASN:HB3	1.74	0.69
1:B:44:ARG:NH1	1:B:55:ASN:OD1	2.24	0.69
1:A:6:VAL:O	1:A:9:LEU:HG	1.93	0.69
1:A:284:GLU:O	1:A:284:GLU:HG3	1.92	0.68
1:A:164:ASP:HB3	1:A:173:PHE:CZ	2.28	0.68
1:A:66:LEU:HD23	1:A:80:ILE:HG13	1.75	0.68
1:A:85:VAL:HG12	1:A:158:ILE:HG22	1.74	0.67
1:A:168:ALA:HB2	1:A:299:MSE:HE2	1.75	0.67
1:A:215:TRP:O	1:A:293:GLN:NE2	2.28	0.67
1:A:78:VAL:CG2	1:A:151:LEU:CB	2.74	0.65
1:A:278:ALA:HB3	1:B:287:VAL:HG23	1.79	0.65
1:B:251:GLN:NE2	1:B:290:ASP:OD2	2.30	0.65
1:A:66:LEU:HD23	1:A:80:ILE:CG1	2.27	0.65
1:B:188:ASP:OD1	1:B:190:ASP:N	2.29	0.64
1:B:184:ALA:HB1	1:B:204:LYS:HE3	1.79	0.64
1:A:77:GLN:NE2	1:A:194:ALA:O	2.30	0.63
1:A:17:TYR:O	1:A:43:GLN:HG3	1.97	0.63
1:B:161:ARG:NH1	1:B:214:TYR:OH	2.32	0.62
1:A:158:ILE:HD11	1:A:179:LEU:CB	2.28	0.62
1:A:77:GLN:OE1	1:A:152:GLN:NE2	2.29	0.61
1:A:77:GLN:HG3	1:A:195:PHE:HD1	1.65	0.61
1:A:14:LEU:HD11	1:A:18:VAL:CG2	2.31	0.61
1:A:252:PHE:O	1:A:288:VAL:CG1	2.47	0.61
1:B:69:PRO:O	1:B:73:LEU:CB	2.49	0.60
1:A:209:GLU:O	1:A:210:SER:HB3	2.01	0.60
1:B:69:PRO:HG2	1:B:73:LEU:HA	1.84	0.60
1:B:61:LEU:CD1	1:B:85:VAL:HG13	2.32	0.59
1:A:280:LEU:HD12	1:B:285:ILE:HB	1.84	0.59
1:A:140:ALA:O	1:A:144:VAL:HG23	2.02	0.59
1:A:54:ILE:HG13	1:A:54:ILE:O	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:PRO:O	1:B:73:LEU:N	2.36	0.59
1:A:203:HIS:NE2	1:A:264:GLU:OE2	2.37	0.58
1:B:37:LEU:HD13	1:B:64:LEU:HD23	1.85	0.58
1:B:259:GLY:HA3	1:B:282:PHE:CD1	2.38	0.58
1:B:188:ASP:OD1	1:B:191:TRP:N	2.36	0.58
1:B:32:ASN:ND2	1:B:34:LYS:HD2	2.19	0.58
1:A:249:GLU:HG3	1:A:249:GLU:O	2.04	0.58
1:B:215:TRP:O	1:B:293:GLN:NE2	2.32	0.58
1:A:214:TYR:HB3	1:A:252:PHE:CD1	2.40	0.57
1:B:32:ASN:HD21	1:B:34:LYS:HD2	1.69	0.57
1:A:182:PHE:HE1	1:A:206:ALA:HB1	1.70	0.57
1:A:9:LEU:HD12	1:A:10:LEU:N	2.19	0.56
1:A:158:ILE:HD11	1:A:179:LEU:CD2	2.32	0.56
1:A:208:LEU:O	1:A:258:SER:HA	2.05	0.56
1:B:189:SER:HA	1:B:203:HIS:CE1	2.40	0.56
1:A:35:VAL:HG12	1:A:66:LEU:H	1.70	0.56
1:A:78:VAL:O	1:A:79:LYS:CG	2.54	0.56
1:A:54:ILE:HB	1:A:89:ALA:HB1	1.87	0.56
1:A:18:VAL:HG12	1:A:18:VAL:O	2.04	0.56
1:A:151:LEU:C	1:A:151:LEU:HD12	2.30	0.56
1:A:157:ASN:HA	1:A:180:GLU:HG2	1.88	0.56
1:B:215:TRP:CZ2	1:B:300:MSE:HE3	2.41	0.56
1:B:61:LEU:HD12	1:B:85:VAL:HG13	1.87	0.56
1:B:64:LEU:HD11	1:B:80:ILE:HD11	1.86	0.56
1:A:80:ILE:O	1:A:80:ILE:HG23	2.05	0.55
1:B:14:LEU:HD23	1:B:21:PHE:CE1	2.42	0.55
1:A:210:SER:O	1:A:210:SER:OG	2.19	0.55
1:A:88:LEU:HB2	1:A:235:PHE:CZ	2.42	0.55
1:A:78:VAL:CG2	1:A:151:LEU:HB3	2.36	0.54
1:A:164:ASP:OD2	1:A:167:SER:HB2	2.08	0.54
1:A:66:LEU:CD2	1:A:80:ILE:CD1	2.86	0.54
1:A:252:PHE:O	1:A:288:VAL:HG12	2.07	0.53
1:A:236:ARG:O	1:A:236:ARG:HG2	2.08	0.53
1:A:78:VAL:O	1:A:79:LYS:HG2	2.09	0.53
1:A:77:GLN:CG	1:A:195:PHE:HD1	2.21	0.53
1:A:158:ILE:O	1:A:158:ILE:HG13	2.09	0.53
1:B:32:ASN:OD1	1:B:34:LYS:HG3	2.09	0.53
1:A:77:GLN:HG3	1:A:195:PHE:CD1	2.44	0.52
1:B:163:GLU:HB3	1:B:221:LEU:HG	1.90	0.52
1:B:176:GLY:HA3	1:B:214:TYR:CE2	2.44	0.52
1:A:158:ILE:CD1	1:A:179:LEU:HD23	2.32	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:VAL:O	1:A:85:VAL:HG13	2.10	0.52
1:A:14:LEU:CD1	1:A:18:VAL:CG2	2.89	0.51
1:A:88:LEU:HB2	1:A:235:PHE:HZ	1.75	0.51
1:A:292:GLN:O	1:A:295:ARG:CG	2.53	0.51
1:A:296:ASP:HB3	1:A:299:MSE:HE3	1.92	0.51
1:A:77:GLN:HG2	1:A:150:ASN:O	2.11	0.50
1:A:78:VAL:HG21	1:A:151:LEU:HB3	1.94	0.50
1:B:205:LEU:HD11	1:B:260:GLN:HG2	1.93	0.50
1:B:32:ASN:OD1	1:B:34:LYS:N	2.37	0.50
1:A:78:VAL:HG22	1:A:151:LEU:HA	1.94	0.49
1:A:159:HIS:C	1:A:159:HIS:CD2	2.90	0.49
1:B:226:ARG:N	1:B:231:MSE:HE2	2.27	0.49
1:A:150:ASN:OD1	1:A:195:PHE:HA	2.12	0.49
1:B:152:GLN:HG3	1:B:185:VAL:HG12	1.94	0.49
1:B:188:ASP:OD1	1:B:190:ASP:C	2.55	0.49
1:A:292:GLN:C	1:A:295:ARG:HG2	2.37	0.49
1:A:78:VAL:CG2	1:A:151:LEU:HB2	2.43	0.49
1:A:88:LEU:HD13	1:A:235:PHE:HZ	1.78	0.49
1:A:228:HIS:CE1	1:A:232:LEU:HD11	2.48	0.49
1:B:26:LEU:HD12	1:B:26:LEU:H	1.78	0.48
1:A:35:VAL:HG12	1:A:66:LEU:HB2	1.96	0.48
1:A:80:ILE:O	1:A:80:ILE:CG2	2.61	0.48
1:B:11:ASN:ND2	1:B:23:PRO:HB3	2.29	0.48
1:A:86:PHE:CZ	1:A:236:ARG:HB2	2.49	0.48
1:A:211:LEU:HB3	1:A:257:VAL:HG13	1.95	0.47
1:B:69:PRO:O	1:B:73:LEU:CA	2.61	0.47
1:B:226:ARG:HA	1:B:231:MSE:HE3	1.96	0.47
1:A:86:PHE:HZ	1:A:236:ARG:HB2	1.79	0.47
1:A:228:HIS:C	1:A:228:HIS:CD2	2.93	0.47
1:A:36:ARG:O	1:A:37:LEU:HD22	2.15	0.47
1:A:63:HIS:HD1	1:A:83:GLU:HB2	1.80	0.47
1:B:228:HIS:CE1	1:B:232:LEU:HD22	2.50	0.47
1:B:227:GLU:H	1:B:231:MSE:HE3	1.80	0.47
1:B:48:ASP:O	1:B:49:GLN:C	2.57	0.46
1:B:250:HIS:HB2	1:B:252:PHE:CE1	2.50	0.46
1:B:195:PHE:C	1:B:196:ILE:HG13	2.40	0.46
1:A:14:LEU:CD1	1:A:18:VAL:HG23	2.46	0.46
1:A:171:HIS:N	1:A:172:PRO:HD3	2.30	0.46
1:B:40:LEU:HB2	1:B:61:LEU:HB2	1.97	0.46
1:B:188:ASP:HA	1:B:196:ILE:CD1	2.45	0.46
1:A:66:LEU:HD21	1:A:80:ILE:CD1	2.39	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:PRO:HG2	1:B:277:LYS:HE2	1.97	0.46
1:A:157:ASN:H	1:A:180:GLU:HA	1.81	0.45
1:A:310:HIS:CD2	1:A:310:HIS:C	2.94	0.45
1:B:86:PHE:HB3	1:B:235:PHE:CE2	2.51	0.45
1:A:151:LEU:CD1	1:A:153:ILE:HG13	2.46	0.45
1:B:64:LEU:HD13	1:B:82:ILE:HG12	1.99	0.45
1:A:38:ASP:OD1	1:A:63:HIS:HB2	2.17	0.45
1:B:188:ASP:HA	1:B:196:ILE:HD13	1.99	0.45
1:B:208:LEU:HD22	1:B:211:LEU:HD22	1.99	0.45
1:B:226:ARG:CA	1:B:231:MSE:CE	2.95	0.45
1:B:259:GLY:HA3	1:B:282:PHE:CE1	2.51	0.45
1:B:37:LEU:HD13	1:B:64:LEU:CD2	2.47	0.45
1:A:310:HIS:CD2	1:A:310:HIS:O	2.70	0.45
1:B:145:THR:O	1:B:149:ASP:CG	2.60	0.45
1:B:3:GLU:OE2	1:B:3:GLU:N	2.50	0.44
1:B:33:GLY:HA3	1:B:70:TRP:HE1	1.80	0.44
1:A:90:SER:HB2	1:A:91:PRO:HD2	1.99	0.44
1:A:309:ARG:NH1	1:B:298:LEU:HD23	2.32	0.44
1:B:159:HIS:ND1	1:B:238:MSE:HG3	2.33	0.44
1:B:17:TYR:HE1	1:B:46:ALA:O	2.00	0.44
1:B:42:LEU:HD12	1:B:59:GLY:C	2.42	0.44
1:A:228:HIS:NE2	1:A:232:LEU:CD1	2.80	0.44
1:B:86:PHE:HB3	1:B:235:PHE:HD2	1.77	0.44
1:B:176:GLY:HA3	1:B:214:TYR:CZ	2.53	0.44
1:A:63:HIS:ND1	1:A:83:GLU:HB2	2.33	0.43
1:A:14:LEU:HD12	1:A:18:VAL:HG23	1.99	0.43
1:A:37:LEU:HB3	1:A:40:LEU:HD11	2.01	0.43
1:B:235:PHE:N	1:B:235:PHE:CD1	2.85	0.43
1:B:48:ASP:O	1:B:51:LYS:N	2.39	0.43
1:B:205:LEU:HD13	1:B:262:LYS:HG2	2.00	0.43
1:B:235:PHE:N	1:B:235:PHE:HD1	2.16	0.43
1:A:253:ILE:O	1:A:287:VAL:HG13	2.18	0.43
1:A:310:HIS:O	1:A:310:HIS:HD2	2.02	0.43
1:B:32:ASN:CG	1:B:34:LYS:HG3	2.44	0.43
1:B:149:ASP:C	1:B:150:ASN:HD22	2.27	0.43
1:A:191:TRP:CH2	1:A:262:LYS:HB3	2.54	0.43
1:B:14:LEU:HD21	1:B:37:LEU:HD21	2.00	0.43
1:B:18:VAL:HG13	1:B:41:GLU:O	2.19	0.42
1:B:311:GLN:O	1:B:314:LYS:HE3	2.19	0.42
1:B:77:GLN:H	1:B:77:GLN:HG2	1.68	0.42
1:A:35:VAL:CG1	1:A:66:LEU:HB2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:THR:HG21	1:A:239:ILE:HA	2.01	0.42
1:A:298:LEU:HD12	1:A:298:LEU:HA	1.67	0.42
1:B:135:ARG:HG3	1:B:135:ARG:HH11	1.84	0.42
1:A:61:LEU:HD11	1:A:82:ILE:HG23	2.02	0.42
1:A:288:VAL:O	1:A:288:VAL:HG13	2.18	0.42
1:A:66:LEU:HD23	1:A:80:ILE:CD1	2.47	0.41
1:A:317:GLN:HA	1:A:318:PRO:HD3	1.71	0.41
1:B:45:GLU:CB	1:B:48:ASP:OD2	2.69	0.41
1:B:183:SER:O	1:B:206:ALA:HA	2.20	0.41
1:A:292:GLN:HA	1:A:295:ARG:HG2	2.01	0.41
1:A:54:ILE:HG22	1:A:162:TYR:CE1	2.55	0.41
1:A:253:ILE:O	1:A:287:VAL:HA	2.21	0.41
1:B:90:SER:OG	1:B:91:PRO:HD2	2.21	0.41
1:B:22:ASP:OD1	1:B:23:PRO:HD2	2.21	0.41
1:B:40:LEU:N	1:B:40:LEU:HD23	2.35	0.41
1:A:41:GLU:HG2	1:A:59:GLY:O	2.20	0.40
1:B:36:ARG:NH1	1:B:38:ASP:OD2	2.54	0.40
1:A:66:LEU:HD23	1:A:80:ILE:HD12	2.02	0.40
1:B:168:ALA:HB2	1:B:299:MSE:HE2	2.03	0.40
1:A:164:ASP:OD2	1:A:167:SER:CB	2.70	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	234/336 (70%)	228 (97%)	6 (3%)	0	100	100
1	B	262/336 (78%)	259 (99%)	3 (1%)	0	100	100
All	All	496/672 (74%)	487 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	184/289 (64%)	184 (100%)	0	100	100
1	B	225/289 (78%)	225 (100%)	0	100	100
All	All	409/578 (71%)	409 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	228	HIS
1	B	171	HIS
1	B	228	HIS
1	B	260	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 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

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	243/336 (72%)	0.51	11 (4%) 38 20	57, 118, 177, 220	0
1	B	265/336 (78%)	0.26	15 (5%) 29 15	42, 75, 151, 190	0
All	All	508/672 (75%)	0.38	26 (5%) 33 17	42, 95, 164, 220	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	270	SER	5.4
1	B	201	SER	4.6
1	B	271	HIS	4.1
1	B	317	GLN	3.9
1	B	268	THR	3.7
1	B	272	THR	3.6
1	A	29	GLU	3.4
1	A	291	ASP	3.3
1	B	200	GLN	3.3
1	B	167	SER	3.1
1	A	229	ASP	3.0
1	A	197	THR	3.0
1	B	153	ILE	2.7
1	A	75	SER	2.7
1	A	85	VAL	2.3
1	B	316	PHE	2.3
1	B	273	VAL	2.2
1	A	199	ILE	2.2
1	B	154	THR	2.1
1	A	162	TYR	2.1
1	A	6	VAL	2.1
1	B	94	GLU	2.1
1	B	224	PRO	2.1
1	A	220	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	181	GLU	2.0
1	B	235	PHE	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 oligosaccharides 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.