



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:48 PM UTC

PDB ID : 2RKO / pdb_00002rko
Title : Crystal Structure of the Vps4p-dimer
Authors : Hartmann, C.; Gruetter, M.G.
Deposited on : 2007-10-17
Resolution : 3.35 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

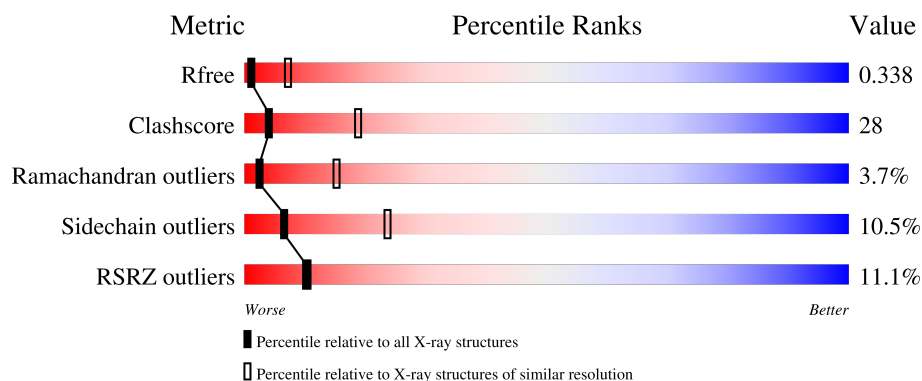
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.35 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	1099 (3.40-3.32)
Clashscore	190562	1116 (3.40-3.32)
Ramachandran outliers	187476	1101 (3.40-3.32)
Sidechain outliers	187428	1101 (3.40-3.32)
RSRZ outliers	180081	1099 (3.40-3.32)

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	331	9% 40% 35% 9% 15%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2173 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 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	280	Total	C	N	O	S	337	0	0
			2173	1380	366	419	8			

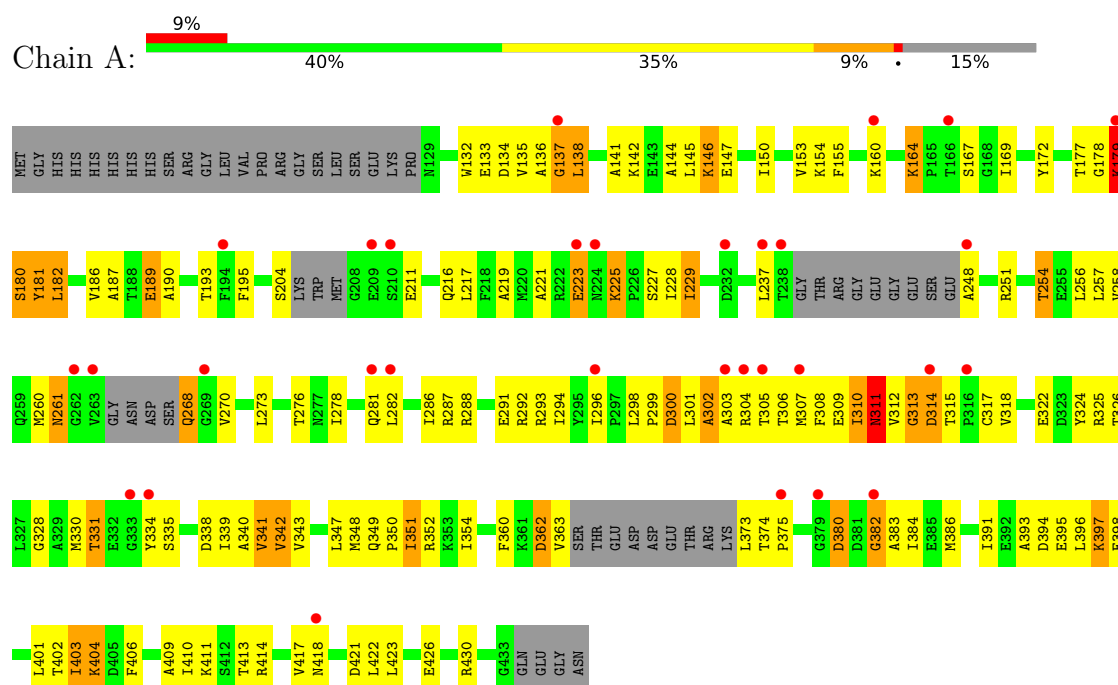
There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	MET	-	expression tag	UNP P52917
A	108	GLY	-	expression tag	UNP P52917
A	109	HIS	-	expression tag	UNP P52917
A	110	HIS	-	expression tag	UNP P52917
A	111	HIS	-	expression tag	UNP P52917
A	112	HIS	-	expression tag	UNP P52917
A	113	HIS	-	expression tag	UNP P52917
A	114	HIS	-	expression tag	UNP P52917
A	115	SER	-	expression tag	UNP P52917
A	116	ARG	-	expression tag	UNP P52917
A	117	GLY	-	expression tag	UNP P52917
A	118	LEU	-	expression tag	UNP P52917
A	119	VAL	-	expression tag	UNP P52917
A	120	PRO	-	expression tag	UNP P52917
A	121	ARG	-	expression tag	UNP P52917
A	122	GLY	-	expression tag	UNP P52917
A	123	SER	-	expression tag	UNP P52917

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: Vacuolar protein sorting-associated protein 4



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	109.12Å 109.12Å 176.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 3.35 30.00 – 3.35	Depositor EDS
% Data completeness (in resolution range)	99.2 (30.00-3.35) 99.0 (30.00-3.35)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.11 (at 3.33Å)	Xtriage
Refinement program	CNS, REFMAC 5.2.0019	Depositor
R, R_{free}	0.268 , 0.298 0.299 , 0.338	Depositor DCC
R_{free} test set	472 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	129.1	Xtriage
Anisotropy	0.188	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 157.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2173	wwPDB-VP
Average B, all atoms (Å ²)	116.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% 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	1.53	39/2208 (1.8%)	1.24	32/2983 (1.1%)

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

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	134	ASP	CA-CB	-17.03	1.27	1.53
1	A	304	ARG	CA-CB	-15.82	1.28	1.53
1	A	418	ASN	CA-CB	-15.72	1.33	1.53
1	A	293	ARG	CA-CB	-15.59	1.30	1.53
1	A	223	GLU	CB-CG	-14.25	1.09	1.52
1	A	430	ARG	CA-CB	-12.95	1.31	1.53
1	A	288	ARG	CA-CB	-12.77	1.32	1.53
1	A	167	SER	CA-CB	-12.48	1.33	1.53
1	A	363	VAL	CB-CG1	12.47	1.93	1.52
1	A	397	LYS	CA-CB	-12.00	1.37	1.53
1	A	287	ARG	CB-CG	-11.95	1.16	1.52
1	A	394	ASP	CA-CB	-11.84	1.36	1.53
1	A	146	LYS	CB-CG	-11.57	1.17	1.52
1	A	237	LEU	CA-CB	-11.00	1.35	1.53
1	A	300	ASP	CA-CB	-10.81	1.36	1.53
1	A	423	LEU	CA-CB	-10.67	1.36	1.53
1	A	393	ALA	CA-CB	-10.51	1.36	1.53
1	A	363	VAL	CB-CG2	-10.46	1.18	1.52
1	A	404	LYS	CA-CB	-10.07	1.37	1.53
1	A	181	TYR	CA-CB	-9.92	1.37	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	380	ASP	CA-CB	-9.64	1.37	1.53
1	A	164	LYS	CA-CB	-9.38	1.34	1.53
1	A	204	SER	CA-CB	-8.36	1.36	1.53
1	A	254	THR	CB-OG1	7.75	1.56	1.43
1	A	261	ASN	CA-CB	-7.60	1.40	1.53
1	A	426	GLU	CA-CB	-7.58	1.39	1.53
1	A	291	GLU	CA-CB	-7.48	1.40	1.53
1	A	282	LEU	CB-CG	-7.39	1.38	1.53
1	A	421	ASP	CA-CB	-7.06	1.42	1.53
1	A	251	ARG	CA-CB	-6.54	1.43	1.53
1	A	211	GLU	CA-CB	-6.51	1.42	1.53
1	A	383	ALA	C-N	6.39	1.41	1.33
1	A	292	ARG	CB-CG	6.36	1.71	1.52
1	A	360	PHE	CB-CG	-6.26	1.36	1.50
1	A	133	GLU	CB-CG	-5.67	1.35	1.52
1	A	268	GLN	CA-CB	5.58	1.64	1.53
1	A	237	LEU	CG-CD2	-5.49	1.34	1.52
1	A	160	LYS	CA-CB	-5.31	1.45	1.53
1	A	382	GLY	C-N	-5.18	1.26	1.33

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	142	LYS	CB-CG-CD	11.08	136.77	111.30
1	A	134	ASP	CA-CB-CG	-10.92	101.68	112.60
1	A	147	GLU	CA-CB-CG	-10.21	93.67	114.10
1	A	146	LYS	CA-CB-CG	10.06	134.21	114.10
1	A	363	VAL	CA-CB-CG2	8.76	125.29	110.40
1	A	223	GLU	CA-CB-CG	8.70	131.50	114.10
1	A	414	ARG	CB-CA-C	-8.16	96.31	109.69
1	A	404	LYS	N-CA-CB	7.77	122.93	109.72
1	A	395	GLU	N-CA-CB	7.66	122.12	110.70
1	A	363	VAL	CA-CB-CG1	-7.11	98.31	110.40
1	A	248	ALA	N-CA-CB	-6.90	100.05	110.40
1	A	302	ALA	CB-CA-C	6.84	120.51	109.02
1	A	147	GLU	CB-CG-CD	-6.76	101.11	112.60
1	A	314	ASP	N-CA-CB	6.74	120.07	109.82
1	A	160	LYS	CB-CA-C	6.43	120.36	109.50
1	A	360	PHE	CA-CB-CG	6.38	120.18	113.80
1	A	302	ALA	N-CA-CB	-6.28	102.55	110.90
1	A	261	ASN	N-CA-CB	6.26	121.07	110.49
1	A	395	GLU	CB-CA-C	-6.17	100.76	110.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	142	LYS	CG-CD-CE	6.16	125.48	111.30
1	A	287	ARG	CA-CB-CG	6.13	126.36	114.10
1	A	237	LEU	CD1-CG-CD2	5.90	123.79	110.80
1	A	225	LYS	CB-CA-C	5.81	121.62	110.17
1	A	293	ARG	CB-CA-C	5.65	118.86	109.53
1	A	394	ASP	CA-CB-CG	-5.48	107.12	112.60
1	A	314	ASP	CA-CB-CG	5.41	118.01	112.60
1	A	423	LEU	CB-CA-C	5.24	119.49	110.79
1	A	254	THR	CA-CB-OG1	-5.17	101.84	109.60
1	A	404	LYS	CB-CA-C	-5.13	102.32	110.84
1	A	211	GLU	CB-CA-C	5.12	120.39	110.46
1	A	383	ALA	CA-C-N	5.08	128.72	122.37
1	A	383	ALA	C-N-CA	5.08	128.72	122.37

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	382	GLY	Mainchain

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	2173	0	2202	99	0
All	All	2173	0	2202	99	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 28.

All (99) 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:315:THR:HG21	1:A:347:LEU:HD22	1.32	1.10
1:A:298:LEU:HD21	1:A:417:VAL:HG21	1.43	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:ILE:CG2	1:A:386:MET:HE3	2.00	0.91
1:A:193:THR:HG1	1:A:195:PHE:HE1	1.19	0.90
1:A:330:MET:HE2	1:A:330:MET:HA	1.60	0.83
1:A:338:ASP:O	1:A:341:VAL:HG13	1.79	0.82
1:A:384:ILE:HG21	1:A:386:MET:HE3	1.60	0.82
1:A:298:LEU:HD12	1:A:422:LEU:HD12	1.65	0.79
1:A:298:LEU:HD11	1:A:417:VAL:HG11	1.69	0.73
1:A:403:ILE:HD12	1:A:403:ILE:H	1.55	0.72
1:A:403:ILE:HD12	1:A:403:ILE:N	2.04	0.71
1:A:257:LEU:HD21	1:A:286:ILE:HD13	1.72	0.71
1:A:298:LEU:HD12	1:A:422:LEU:CD1	2.22	0.70
1:A:132:TRP:CD1	1:A:189:GLU:HG3	2.27	0.69
1:A:311:ASN:N	1:A:311:ASN:HD22	1.89	0.68
1:A:132:TRP:CG	1:A:189:GLU:HG3	2.29	0.68
1:A:182:LEU:HD23	1:A:296:ILE:HD13	1.77	0.67
1:A:138:LEU:HB2	1:A:141:ALA:HB3	1.77	0.67
1:A:169:ILE:HB	1:A:273:LEU:HD22	1.76	0.66
1:A:298:LEU:CD1	1:A:417:VAL:HG11	2.27	0.65
1:A:318:VAL:HG12	1:A:318:VAL:O	1.96	0.64
1:A:310:ILE:O	1:A:312:VAL:N	2.30	0.63
1:A:153:VAL:HG11	1:A:190:ALA:HB1	1.81	0.62
1:A:298:LEU:CD1	1:A:422:LEU:HD12	2.28	0.62
1:A:300:ASP:O	1:A:302:ALA:N	2.33	0.62
1:A:146:LYS:HA	1:A:150:ILE:HD12	1.80	0.62
1:A:178:GLY:O	1:A:181:TYR:N	2.32	0.61
1:A:362:ASP:HA	1:A:373:LEU:HD23	1.82	0.61
1:A:309:GLU:HA	1:A:324:TYR:CE2	2.35	0.61
1:A:351:ILE:H	1:A:351:ILE:HD13	1.67	0.60
1:A:298:LEU:CD2	1:A:417:VAL:HG21	2.25	0.59
1:A:384:ILE:HG22	1:A:386:MET:HE3	1.82	0.58
1:A:310:ILE:C	1:A:312:VAL:H	2.09	0.58
1:A:187:ALA:HB2	1:A:228:ILE:HD12	1.84	0.57
1:A:216:GLN:O	1:A:219:ALA:HB3	2.04	0.57
1:A:305:THR:HG23	1:A:324:TYR:C	2.29	0.57
1:A:254:THR:O	1:A:258:VAL:HG23	2.04	0.57
1:A:347:LEU:O	1:A:350:PRO:HD2	2.08	0.54
1:A:135:VAL:HG21	1:A:182:LEU:HD13	1.90	0.54
1:A:146:LYS:O	1:A:150:ILE:HB	2.08	0.52
1:A:403:ILE:H	1:A:403:ILE:CD1	2.16	0.52
1:A:311:ASN:N	1:A:311:ASN:ND2	2.58	0.52
1:A:348:MET:SD	1:A:351:ILE:HD11	2.49	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:LEU:H	1:A:138:LEU:HD13	1.75	0.50
1:A:138:LEU:H	1:A:138:LEU:HD22	1.76	0.50
1:A:315:THR:CG2	1:A:347:LEU:HD22	2.24	0.49
1:A:227:SER:OG	1:A:270:VAL:HG22	2.12	0.49
1:A:343:VAL:HG13	1:A:406:PHE:HE2	1.78	0.49
1:A:300:ASP:C	1:A:302:ALA:H	2.21	0.49
1:A:300:ASP:C	1:A:302:ALA:N	2.69	0.49
1:A:311:ASN:OD1	1:A:340:ALA:CB	2.60	0.49
1:A:315:THR:HG21	1:A:347:LEU:CD2	2.24	0.49
1:A:374:THR:HG23	1:A:375:PRO:HD2	1.95	0.48
1:A:135:VAL:O	1:A:136:ALA:C	2.55	0.48
1:A:193:THR:OG1	1:A:195:PHE:HE1	1.89	0.48
1:A:349:GLN:HB2	1:A:350:PRO:HD3	1.96	0.48
1:A:172:TYR:HA	1:A:276:THR:O	2.14	0.47
1:A:312:VAL:O	1:A:313:GLY:C	2.59	0.46
1:A:178:GLY:O	1:A:179:LYS:C	2.58	0.46
1:A:306:THR:O	1:A:309:GLU:HB2	2.16	0.46
1:A:322:GLU:O	1:A:325:ARG:N	2.49	0.46
1:A:331:THR:HB	1:A:334:TYR:HB2	1.98	0.45
1:A:195:PHE:HB3	1:A:217:LEU:HD11	1.98	0.45
1:A:308:PHE:CE1	1:A:339:ILE:HG23	2.52	0.45
1:A:347:LEU:HD23	1:A:401:LEU:HD11	1.97	0.45
1:A:354:ILE:HD13	1:A:375:PRO:HG2	1.99	0.44
1:A:136:ALA:O	1:A:137:GLY:C	2.60	0.44
1:A:374:THR:OG1	1:A:397:LYS:O	2.33	0.44
1:A:221:ALA:CB	1:A:229:ILE:HD11	2.47	0.44
1:A:179:LYS:N	1:A:179:LYS:HD2	2.31	0.44
1:A:187:ALA:CB	1:A:228:ILE:HD12	2.47	0.44
1:A:342:VAL:HG12	1:A:413:THR:HG21	2.00	0.43
1:A:325:ARG:O	1:A:326:THR:C	2.60	0.43
1:A:384:ILE:O	1:A:386:MET:HG2	2.18	0.43
1:A:354:ILE:HD13	1:A:375:PRO:CG	2.49	0.43
1:A:409:ALA:O	1:A:410:ILE:C	2.61	0.43
1:A:299:PRO:HB3	1:A:307:MET:HE1	2.00	0.43
1:A:317:CYS:HA	1:A:401:LEU:HB2	2.00	0.42
1:A:256:LEU:HD12	1:A:260:MET:HG3	2.02	0.42
1:A:260:MET:O	1:A:261:ASN:C	2.62	0.42
1:A:343:VAL:O	1:A:347:LEU:HG	2.20	0.42
1:A:403:ILE:O	1:A:404:LYS:C	2.62	0.42
1:A:384:ILE:HG21	1:A:386:MET:CE	2.42	0.42
1:A:300:ASP:O	1:A:303:ALA:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:TRP:CE3	1:A:135:VAL:HG11	2.55	0.41
1:A:177:THR:H	1:A:179:LYS:HE3	1.85	0.41
1:A:325:ARG:O	1:A:328:GLY:N	2.53	0.41
1:A:406:PHE:O	1:A:410:ILE:HG13	2.19	0.41
1:A:178:GLY:O	1:A:180:SER:N	2.54	0.41
1:A:339:ILE:O	1:A:343:VAL:HG23	2.20	0.41
1:A:145:LEU:HB3	1:A:186:VAL:HG21	2.02	0.41
1:A:314:ASP:O	1:A:315:THR:C	2.63	0.41
1:A:331:THR:HG23	1:A:410:ILE:HD13	2.01	0.41
1:A:422:LEU:C	1:A:422:LEU:HD23	2.46	0.41
1:A:144:ALA:CB	1:A:294:ILE:HD12	2.51	0.41
1:A:318:VAL:O	1:A:318:VAL:CG1	2.67	0.41
1:A:374:THR:CG2	1:A:375:PRO:HD2	2.51	0.40
1:A:138:LEU:HD13	1:A:138:LEU:N	2.37	0.40
1:A:144:ALA:HB1	1:A:294:ILE:CD1	2.51	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	270/331 (82%)	224 (83%)	36 (13%)	10 (4%)	2	15

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	137	GLY
1	A	179	LYS
1	A	301	LEU
1	A	225	LYS
1	A	313	GLY
1	A	223	GLU

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Mol	Chain	Res	Type
1	A	311	ASN
1	A	154	LYS
1	A	155	PHE
1	A	310	ILE

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	239/282 (85%)	214 (90%)	25 (10%)	6 24

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	LEU
1	A	164	LYS
1	A	179	LYS
1	A	180	SER
1	A	182	LEU
1	A	189	GLU
1	A	229	ILE
1	A	268	GLN
1	A	278	ILE
1	A	281	GLN
1	A	311	ASN
1	A	331	THR
1	A	335	SER
1	A	341	VAL
1	A	342	VAL
1	A	351	ILE
1	A	352	ARG
1	A	362	ASP
1	A	380	ASP
1	A	391	ILE
1	A	396	LEU
1	A	398	GLU

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Mol	Chain	Res	Type
1	A	402	THR
1	A	403	ILE
1	A	411	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	191	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 ⓘ

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	279/331 (84%)	0.86	31 (11%)	10 10	40, 118, 150, 167	77 (27%)

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	263	VAL	4.8
1	A	232	ASP	4.4
1	A	269	GLY	3.9
1	A	179	LYS	3.9
1	A	262	GLY	3.5
1	A	305	THR	3.3
1	A	282	LEU	3.3
1	A	333	GLY	3.2
1	A	303	ALA	3.1
1	A	304	ARG	2.9
1	A	307	MET	2.7
1	A	281	GLN	2.7
1	A	166	THR	2.7
1	A	209	GLU	2.7
1	A	238	THR	2.6
1	A	210	SER	2.6
1	A	237	LEU	2.6
1	A	375	PRO	2.5
1	A	137	GLY	2.3
1	A	296	ILE	2.3
1	A	194	PHE	2.2
1	A	418	ASN	2.2
1	A	248	ALA	2.1
1	A	224	ASN	2.1
1	A	379	GLY	2.1
1	A	334	TYR	2.1
1	A	314	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	382	GLY	2.1
1	A	223	GLU	2.1
1	A	160	LYS	2.0
1	A	316	PRO	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.