



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:58 PM UTC

PDB ID : 4AKF / pdb_00004akf
Title : Crystal structure of VipD from Legionella pneumophila
Authors : Lee, K.-H.; Ku, B.; Oh, B.-H.
Deposited on : 2012-02-22
Resolution : 2.90 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

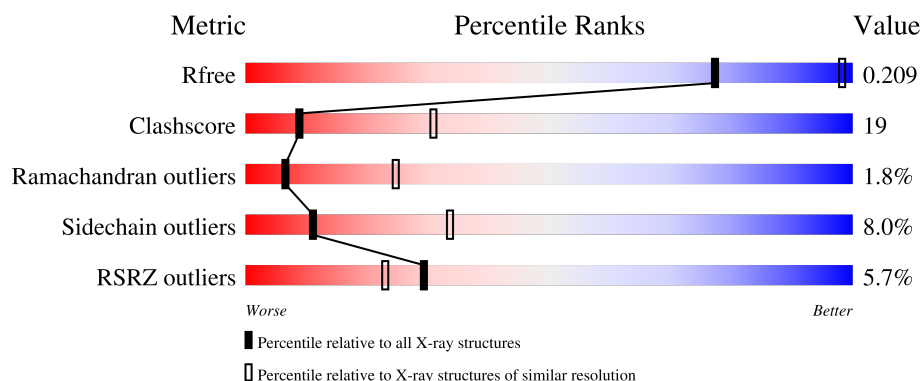
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	577	6% 61% 31% 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4466 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 VIPD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	561	Total	C	N	O	S	0	0	1
			4380	2762	754	845	19			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q5ZRP9
A	0	SER	-	expression tag	UNP Q5ZRP9

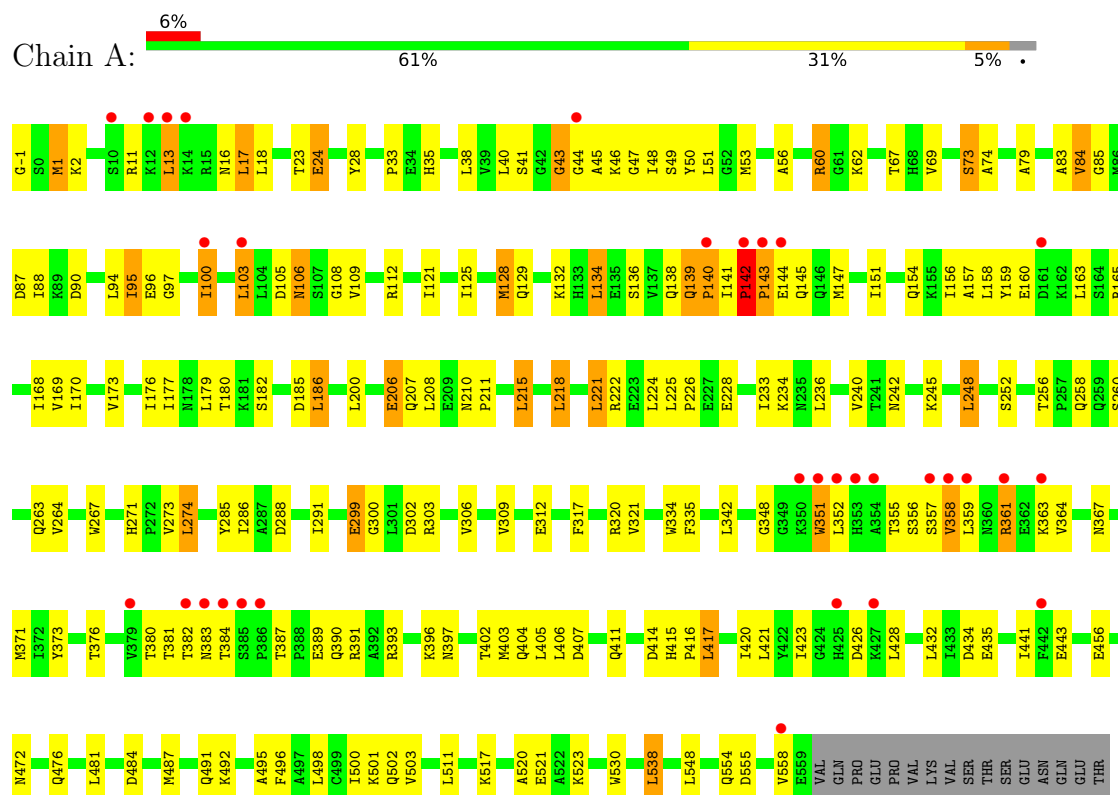
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	86	Total	O	0	0
			86	86		

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: VIPD



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 3 2	Depositor
Cell constants a, b, c, α , β , γ	252.76 Å 252.76 Å 252.76 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 30.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (30.00-2.90) 99.3 (30.00-2.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 2.81 Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.214 , 0.233 0.210 , 0.209	Depositor DCC
R_{free} test set	1722 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	55.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 29.5	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	0.94	EDS
Total number of atoms	4466	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.54% 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	0.52	1/4446 (0.0%)	0.99	21/5995 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	558	VAL	C-N	-5.49	1.25	1.33

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	142	PRO	N-CA-C	10.22	123.17	110.70
1	A	299	GLU	N-CA-C	9.26	122.32	111.02
1	A	358	VAL	N-CA-C	-8.50	104.30	112.29
1	A	361	ARG	N-CA-C	-7.57	104.01	112.57
1	A	443	GLU	N-CA-C	-7.11	103.61	111.36
1	A	24	GLU	N-CA-C	-7.06	103.92	112.54
1	A	356	SER	N-CA-C	-6.97	104.50	112.87
1	A	306	VAL	N-CA-C	6.76	117.62	107.75
1	A	256	THR	CA-C-N	5.95	125.63	119.56
1	A	256	THR	C-N-CA	5.95	125.63	119.56
1	A	67	THR	N-CA-C	-5.84	106.79	114.04
1	A	225	LEU	CA-C-N	5.77	125.53	119.76
1	A	225	LEU	C-N-CA	5.77	125.53	119.76
1	A	384	THR	N-CA-C	-5.58	105.91	114.16
1	A	206	GLU	N-CA-C	-5.50	103.12	110.55
1	A	342	LEU	N-CA-C	5.49	118.02	111.71
1	A	383	ASN	N-CA-C	5.47	118.01	108.76
1	A	335	PHE	N-CA-C	-5.37	105.33	111.07
1	A	236	LEU	N-CA-C	5.19	118.68	109.96
1	A	484	ASP	N-CA-C	5.09	116.51	111.07
1	A	106	ASN	N-CA-C	-5.03	106.57	113.30

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	4380	0	4461	164	0
2	A	86	0	0	6	0
All	All	4466	0	4461	164	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 19.

All (164) 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:46:LYS:HB3	1:A:376:THR:HG21	1.37	1.05
1:A:248:LEU:H	1:A:248:LEU:HD23	1.29	0.94
1:A:53:MET:HE2	1:A:309:VAL:HG21	1.51	0.91
1:A:142:PRO:HB2	1:A:143:PRO:HD3	1.51	0.90
1:A:90:ASP:HB3	1:A:173:VAL:HG21	1.53	0.89
1:A:487:MET:HE3	1:A:492:LYS:HA	1.57	0.87
1:A:271:HIS:HD2	1:A:273:VAL:H	1.30	0.79
1:A:260:SER:O	1:A:264:VAL:HG23	1.83	0.78
1:A:207:GLN:HE22	1:A:258:GLN:HG2	1.48	0.77
1:A:364:VAL:CG1	1:A:502:GLN:HG2	2.14	0.77
1:A:263:GLN:HG3	2:A:2041:HOH:O	1.85	0.77
1:A:210:ASN:H	1:A:258:GLN:NE2	1.84	0.75
1:A:165:ARG:HB3	1:A:165:ARG:HH11	1.53	0.73
1:A:364:VAL:HG13	1:A:502:GLN:HG2	1.72	0.72
1:A:348:GLY:HA2	1:A:523:LYS:HE3	1.73	0.71
1:A:207:GLN:NE2	1:A:258:GLN:HG2	2.05	0.70
1:A:434:ASP:OD1	1:A:435:GLU:N	2.23	0.70
1:A:143:PRO:C	1:A:145:GLN:H	1.99	0.70
1:A:142:PRO:CB	1:A:143:PRO:HD3	2.21	0.70
1:A:481:LEU:HB3	1:A:487:MET:HE2	1.74	0.69
1:A:501:LYS:HE3	1:A:538:LEU:O	1.92	0.69
1:A:108:GLY:HA3	1:A:112:ARG:O	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:LEU:HD12	1:A:13:LEU:H	1.58	0.68
1:A:548:LEU:HD12	1:A:548:LEU:N	2.09	0.67
1:A:128:MET:HE3	1:A:132:LYS:HE3	1.77	0.67
1:A:321:VAL:HG23	1:A:404:GLN:HG2	1.76	0.67
1:A:46:LYS:HZ2	1:A:312:GLU:HA	1.60	0.66
1:A:180:THR:HA	1:A:186:LEU:HD11	1.77	0.66
1:A:142:PRO:HB2	1:A:143:PRO:CD	2.25	0.66
1:A:402:THR:O	1:A:406:LEU:HB2	1.95	0.66
1:A:60:ARG:NH2	1:A:407:ASP:OD1	2.29	0.65
1:A:165:ARG:HB3	1:A:165:ARG:NH1	2.12	0.64
1:A:210:ASN:H	1:A:258:GLN:HE21	1.42	0.64
1:A:415:HIS:CE1	1:A:417:LEU:HB2	2.32	0.64
1:A:517:LYS:O	1:A:521:GLU:HG3	1.98	0.64
1:A:128:MET:HE2	1:A:129:GLN:HA	1.79	0.63
1:A:-1:GLY:HA3	1:A:300:GLY:O	1.99	0.62
1:A:156:ILE:O	1:A:160:GLU:HG3	1.98	0.62
1:A:28:TYR:CE1	1:A:416:PRO:HD3	2.35	0.62
1:A:417:LEU:HD22	1:A:421:LEU:HG	1.81	0.62
1:A:226:PRO:HB2	1:A:228:GLU:OE2	2.00	0.62
1:A:134:LEU:HD12	1:A:156:ILE:HD12	1.82	0.62
1:A:288:ASP:HB3	1:A:291:ILE:HG23	1.80	0.61
1:A:44:GLY:O	1:A:74:ALA:HB2	1.99	0.61
1:A:248:LEU:H	1:A:248:LEU:CD2	2.10	0.61
1:A:16:ASN:HD22	1:A:432:LEU:HD12	1.66	0.60
1:A:285:TYR:C	1:A:286:ILE:HD12	2.26	0.60
1:A:355:THR:HG22	1:A:357:SER:H	1.67	0.60
1:A:393:ARG:O	1:A:397:ASN:HB2	2.02	0.59
1:A:396:LYS:HB3	2:A:2067:HOH:O	2.01	0.59
1:A:165:ARG:HH11	1:A:165:ARG:CB	2.15	0.59
1:A:491:GLN:HG2	2:A:2077:HOH:O	2.02	0.59
1:A:13:LEU:HD12	1:A:13:LEU:N	2.16	0.59
1:A:121:ILE:HD12	1:A:180:THR:HG22	1.84	0.58
1:A:355:THR:C	1:A:357:SER:N	2.62	0.57
1:A:363:LYS:HG2	2:A:2061:HOH:O	2.03	0.57
1:A:96:GLU:O	1:A:96:GLU:HG3	2.05	0.56
1:A:248:LEU:HD23	1:A:248:LEU:N	2.10	0.56
1:A:103:LEU:HA	2:A:2015:HOH:O	2.06	0.56
1:A:271:HIS:CE1	1:A:381:THR:HG21	2.41	0.56
1:A:200:LEU:O	1:A:208:LEU:HB2	2.06	0.55
1:A:222:ARG:NH1	1:A:233:ILE:HG22	2.22	0.55
1:A:103:LEU:C	1:A:103:LEU:HD23	2.31	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:LEU:HD13	1:A:233:ILE:CD1	2.37	0.55
1:A:355:THR:C	1:A:357:SER:H	2.13	0.55
1:A:143:PRO:C	1:A:145:GLN:N	2.62	0.55
1:A:415:HIS:HE1	1:A:417:LEU:HB2	1.72	0.55
1:A:84:VAL:HG13	1:A:84:VAL:O	2.08	0.54
1:A:487:MET:HE1	1:A:495:ALA:HB3	1.88	0.54
1:A:496:PHE:O	1:A:500:ILE:HG13	2.08	0.53
1:A:18:LEU:HD21	1:A:423:ILE:HG23	1.91	0.53
1:A:47:GLY:HA3	1:A:74:ALA:HB1	1.89	0.53
1:A:320:ARG:HG2	1:A:320:ARG:HH11	1.72	0.53
1:A:503:VAL:HG12	1:A:511:LEU:HD22	1.91	0.53
1:A:517:LYS:HE2	1:A:521:GLU:OE2	2.06	0.53
1:A:228:GLU:H	1:A:228:GLU:CD	2.16	0.53
1:A:359:LEU:C	1:A:361:ARG:H	2.16	0.52
1:A:23:THR:HG22	1:A:24:GLU:H	1.74	0.52
1:A:142:PRO:O	1:A:143:PRO:O	2.27	0.52
1:A:221:LEU:HD13	1:A:233:ILE:HD11	1.92	0.52
1:A:169:VAL:C	1:A:170:ILE:HD12	2.34	0.52
1:A:48:ILE:HG21	1:A:391:ARG:HG3	1.93	0.51
1:A:182:SER:HB3	1:A:185:ASP:HB2	1.92	0.51
1:A:139:GLN:N	1:A:140:PRO:HD2	2.26	0.51
1:A:221:LEU:HD22	1:A:233:ILE:CD1	2.41	0.51
1:A:271:HIS:CD2	1:A:273:VAL:H	2.19	0.51
1:A:128:MET:HE3	1:A:132:LYS:CE	2.41	0.50
1:A:43:GLY:O	1:A:46:LYS:HG3	2.11	0.49
1:A:100:ILE:HD12	1:A:100:ILE:O	2.12	0.49
1:A:23:THR:HG22	1:A:24:GLU:N	2.27	0.49
1:A:351:TRP:CD2	1:A:351:TRP:C	2.91	0.49
1:A:138:GLN:HB3	1:A:140:PRO:HD2	1.95	0.48
1:A:267:TRP:HB3	1:A:286:ILE:HG21	1.94	0.48
1:A:487:MET:CE	1:A:492:LYS:HA	2.38	0.48
1:A:2:LYS:HG2	1:A:2:LYS:O	2.12	0.48
1:A:312:GLU:HB2	1:A:317:PHE:CE2	2.49	0.48
1:A:520:ALA:HA	2:A:2057:HOH:O	2.13	0.47
1:A:142:PRO:CB	1:A:143:PRO:CD	2.90	0.47
1:A:128:MET:HE2	1:A:128:MET:C	2.39	0.47
1:A:41:SER:O	1:A:50:TYR:OH	2.33	0.47
1:A:147:MET:O	1:A:151:ILE:HG13	2.15	0.46
1:A:16:ASN:ND2	1:A:432:LEU:HD12	2.29	0.46
1:A:56:ALA:HA	1:A:403:MET:HE2	1.96	0.46
1:A:179:LEU:HD12	1:A:185:ASP:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:LEU:O	1:A:218:LEU:HB2	2.16	0.46
1:A:242:ASN:OD1	1:A:245:LYS:HG3	2.16	0.46
1:A:35:HIS:ND1	1:A:62:LYS:NZ	2.55	0.45
1:A:103:LEU:HD23	1:A:103:LEU:O	2.16	0.45
1:A:168:ILE:HG22	1:A:170:ILE:CD1	2.47	0.45
1:A:302:ASP:O	1:A:303:ARG:C	2.59	0.45
1:A:487:MET:HE1	1:A:495:ALA:CB	2.46	0.45
1:A:17:LEU:O	1:A:33:PRO:HD3	2.15	0.45
1:A:154:GLN:O	1:A:157:ALA:HB3	2.16	0.45
1:A:334:TRP:HZ2	1:A:358:VAL:HG11	1.81	0.45
1:A:286:ILE:HD12	1:A:286:ILE:N	2.31	0.45
1:A:387:THR:O	1:A:389:GLU:N	2.50	0.45
1:A:100:ILE:HA	1:A:103:LEU:CB	2.47	0.45
1:A:334:TRP:CZ2	1:A:358:VAL:HG11	2.52	0.45
1:A:143:PRO:O	1:A:145:GLN:N	2.50	0.44
1:A:411:GLN:HE21	1:A:411:GLN:HB3	1.64	0.44
1:A:40:LEU:HD21	1:A:53:MET:HE1	2.00	0.44
1:A:351:TRP:C	1:A:351:TRP:CE3	2.96	0.44
1:A:53:MET:CE	1:A:309:VAL:HG21	2.37	0.44
1:A:435:GLU:HA	1:A:435:GLU:OE1	2.17	0.44
1:A:125:ILE:HG12	1:A:176:ILE:HD13	1.99	0.43
1:A:503:VAL:CG1	1:A:511:LEU:HD22	2.48	0.43
1:A:69:VAL:HG12	1:A:79:ALA:HB1	2.00	0.43
1:A:248:LEU:CD2	1:A:248:LEU:N	2.76	0.43
1:A:90:ASP:HB3	1:A:173:VAL:CG2	2.38	0.43
1:A:94:LEU:O	1:A:95:ILE:C	2.61	0.43
1:A:472:ASN:O	1:A:476:GLN:HG3	2.17	0.43
1:A:273:VAL:HG12	1:A:274:LEU:CD1	2.49	0.43
1:A:13:LEU:H	1:A:13:LEU:CD1	2.28	0.43
1:A:97:GLY:O	1:A:177:ILE:HG23	2.19	0.43
1:A:387:THR:C	1:A:389:GLU:N	2.76	0.42
1:A:417:LEU:CD2	1:A:421:LEU:HG	2.48	0.42
1:A:215:LEU:HD12	1:A:215:LEU:HA	1.92	0.42
1:A:498:LEU:O	1:A:502:GLN:HG3	2.19	0.42
1:A:83:ALA:C	1:A:85:GLY:H	2.27	0.42
1:A:359:LEU:C	1:A:361:ARG:N	2.76	0.42
1:A:83:ALA:HA	1:A:234:LYS:HB2	2.02	0.42
1:A:222:ARG:NH1	1:A:233:ILE:O	2.53	0.42
1:A:530:TRP:CD1	1:A:530:TRP:H	2.36	0.42
1:A:45:ALA:C	1:A:47:GLY:H	2.27	0.42
1:A:49:SER:C	1:A:51:LEU:N	2.78	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:SER:O	1:A:74:ALA:C	2.62	0.41
1:A:106:ASN:C	1:A:108:GLY:H	2.28	0.41
1:A:200:LEU:HD12	1:A:211:PRO:HG3	2.02	0.41
1:A:334:TRP:CD1	1:A:359:LEU:HG	2.55	0.41
1:A:1:MET:HE3	1:A:299:GLU:HB3	2.02	0.41
1:A:140:PRO:O	1:A:141:ILE:HB	2.20	0.41
1:A:179:LEU:CD1	1:A:185:ASP:HB3	2.50	0.41
1:A:180:THR:HA	1:A:186:LEU:CD1	2.48	0.41
1:A:121:ILE:O	1:A:125:ILE:HG13	2.21	0.41
1:A:371:MET:HG2	1:A:373:TYR:CZ	2.56	0.41
1:A:420:ILE:HD11	1:A:441:ILE:HG23	2.03	0.41
1:A:371:MET:HE2	1:A:371:MET:HB2	1.90	0.41
1:A:159:TYR:O	1:A:163:LEU:HG	2.20	0.41
1:A:51:LEU:HD23	1:A:51:LEU:O	2.21	0.40
1:A:221:LEU:HD22	1:A:233:ILE:HD11	2.04	0.40
1:A:414:ASP:HB3	1:A:491:GLN:HE21	1.85	0.40
1:A:43:GLY:O	1:A:46:LYS:CE	2.68	0.40
1:A:221:LEU:HD13	1:A:233:ILE:HD13	2.03	0.40
1:A:405:LEU:HD23	1:A:405:LEU:C	2.46	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	559/577 (97%)	507 (91%)	42 (8%)	10 (2%)	6 25

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	142	PRO
1	A	143	PRO

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Mol	Chain	Res	Type
1	A	43	GLY
1	A	140	PRO
1	A	105	ASP
1	A	144	GLU
1	A	390	GLN
1	A	88	ILE
1	A	95	ILE
1	A	84	VAL

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	476/493 (97%)	438 (92%)	38 (8%)	11	34

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	11	ARG
1	A	13	LEU
1	A	17	LEU
1	A	38	LEU
1	A	60	ARG
1	A	73	SER
1	A	87	ASP
1	A	100	ILE
1	A	103	LEU
1	A	109	VAL
1	A	128	MET
1	A	134	LEU
1	A	136	SER
1	A	139	GLN
1	A	158	LEU
1	A	186	LEU
1	A	206	GLU

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Mol	Chain	Res	Type
1	A	215	LEU
1	A	218	LEU
1	A	221	LEU
1	A	224	LEU
1	A	240	VAL
1	A	248	LEU
1	A	252	SER
1	A	274	LEU
1	A	351	TRP
1	A	352	LEU
1	A	367	ASN
1	A	380	THR
1	A	382	THR
1	A	417	LEU
1	A	426	ASP
1	A	428	LEU
1	A	456	GLU
1	A	538	LEU
1	A	554	GLN
1	A	555	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	146	GLN
1	A	178	ASN
1	A	194	ASN
1	A	207	GLN
1	A	229	ASN
1	A	243	GLN
1	A	258	GLN
1	A	259	GLN
1	A	271	HIS
1	A	294	ASN
1	A	360	ASN
1	A	390	GLN
1	A	397	ASN
1	A	411	GLN
1	A	449	GLN
1	A	471	GLN
1	A	491	GLN
1	A	554	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	561/577 (97%)	-0.01	32 (5%) 29 23	37, 64, 109, 133	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	352	LEU	6.7
1	A	353	HIS	5.5
1	A	354	ALA	4.6
1	A	386	PRO	3.9
1	A	442	PHE	3.9
1	A	13	LEU	3.7
1	A	144	GLU	3.5
1	A	383	ASN	3.3
1	A	142	PRO	3.3
1	A	161	ASP	3.2
1	A	384	THR	3.1
1	A	10	SER	3.1
1	A	100	ILE	3.1
1	A	359	LEU	3.1
1	A	351	TRP	2.9
1	A	14	LYS	2.9
1	A	350	LYS	2.8
1	A	379	VAL	2.8
1	A	12	LYS	2.8
1	A	44	GLY	2.7
1	A	143	PRO	2.7
1	A	558	VAL	2.6
1	A	361	ARG	2.5
1	A	140	PRO	2.5
1	A	357	SER	2.5
1	A	427	LYS	2.3
1	A	382	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	385	SER	2.2
1	A	103	LEU	2.2
1	A	425	HIS	2.1
1	A	358	VAL	2.1
1	A	363	LYS	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 [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.