



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 08:27 PM UTC

PDB ID : 2R02 / pdb_00002r02
Title : Crystal Structure of ALIX/AIP1 in complex with the HIV-1 YPLTSL Late Domain
Authors : Hill, C.P.; Zhai, Q.; Fisher, R.D.
Deposited on : 2007-08-17
Resolution : 2.60 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

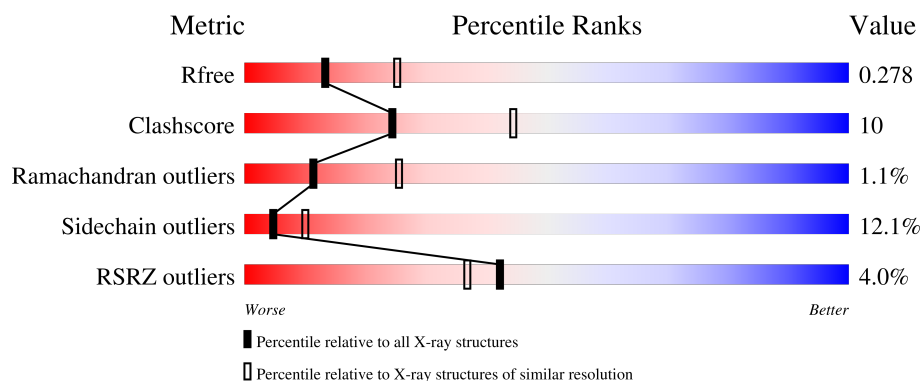
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	697	
2	B	11	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5593 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 Programmed cell death 6-interacting protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	697	Total	C	N	O	S	0	0	0
			5486	3461	938	1069	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	268	TYR	LYS	engineered mutation	UNP Q8WUM4
A	269	TYR	LYS	engineered mutation	UNP Q8WUM4

- Molecule 2 is a protein called p6-gag.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	11	Total	C	N	O	0	0	0
			90	59	14	17			

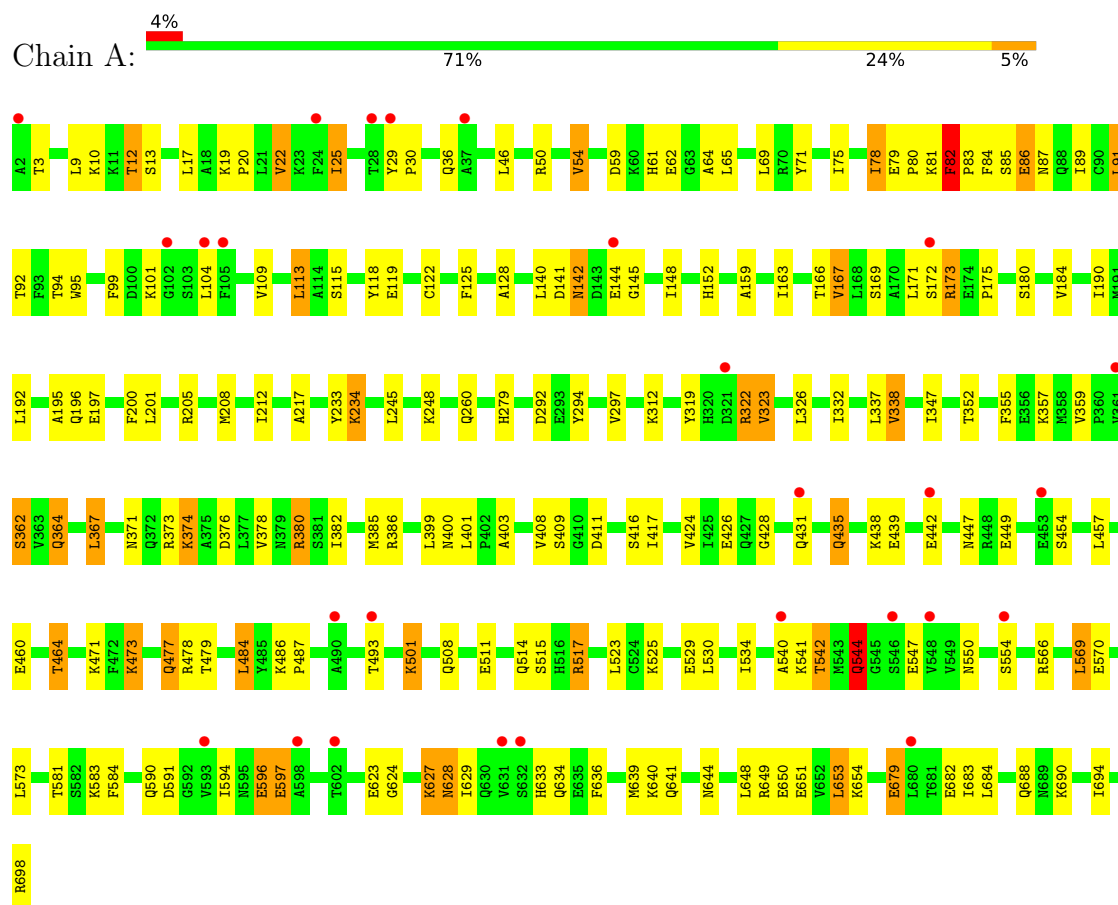
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	17	Total	O	0	0
			17	17		

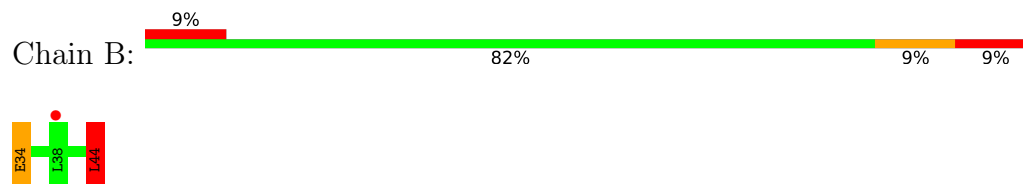
3 Residue-property plots [i](#)

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: Programmed cell death 6-interacting protein



- Molecule 2: p6-gag



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	145.92Å 99.08Å 73.23Å 90.00° 107.32° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60 50.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.1 (50.00-2.60) 98.0 (50.00-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.96 (at 2.61Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.224 , 0.285 0.221 , 0.278	Depositor DCC
R_{free} test set	1501 reflections (4.10%)	wwPDB-VP
Wilson B-factor (Å ²)	62.7	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 53.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5593	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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	0.86	10/5570 (0.2%)	1.03	11/7523 (0.1%)
2	B	2.96	1/91 (1.1%)	1.02	1/123 (0.8%)
All	All	0.93	11/5661 (0.2%)	1.03	12/7646 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	44	LEU	C-O	27.69	1.78	1.23
1	A	547	GLU	CD-OE2	14.21	1.52	1.25
1	A	627	LYS	CD-CE	11.03	1.85	1.52
1	A	627	LYS	CE-NZ	10.76	1.81	1.49
1	A	547	GLU	CD-OE1	9.80	1.44	1.25
1	A	554	SER	CB-OG	8.11	1.58	1.42
1	A	374	LYS	CE-NZ	8.06	1.73	1.49
1	A	544	GLN	CD-OE1	7.84	1.38	1.23
1	A	544	GLN	CD-NE2	6.79	1.47	1.33
1	A	362	SER	C-O	5.77	1.31	1.24
1	A	550	ASN	CG-ND2	5.49	1.44	1.33

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	217	ALA	N-CA-C	-7.61	102.33	111.69
2	B	44	LEU	CA-C-O	-7.51	108.03	120.80
1	A	627	LYS	CD-CE-NZ	-6.86	89.94	111.90
1	A	260	GLN	N-CA-C	-6.80	104.45	112.89
1	A	337	LEU	N-CA-C	-6.42	103.05	113.19
1	A	641	GLN	N-CA-C	6.28	118.66	110.43
1	A	323	VAL	N-CA-C	5.59	112.65	107.56
1	A	195	ALA	N-CA-C	-5.53	105.16	111.07
1	A	544	GLN	OE1-CD-NE2	5.47	128.07	122.60
1	A	319	TYR	N-CA-C	5.11	117.74	111.82
1	A	541	LYS	N-CA-C	5.11	118.61	112.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	332	ILE	N-CA-C	-5.02	103.10	109.58

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	5486	0	5532	109	0
2	B	90	0	95	2	0
3	A	17	0	0	0	0
All	All	5593	0	5627	111	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 10.

All (111) 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:627:LYS:CD	1:A:627:LYS:CE	1.85	1.51
1:A:374:LYS:CE	1:A:374:LYS:NZ	1.73	1.48
1:A:627:LYS:CE	1:A:627:LYS:NZ	1.81	1.40
2:B:44:LEU:O	2:B:44:LEU:C	1.78	1.24
1:A:486:LYS:HG3	1:A:487:PRO:HD3	1.29	1.06
1:A:517:ARG:HH11	1:A:517:ARG:HG3	1.36	0.89
1:A:3:THR:HG21	1:A:294:TYR:O	1.72	0.87
1:A:382:ILE:HD12	1:A:570:GLU:HG3	1.56	0.87
1:A:540:ALA:HB3	1:A:640:LYS:HG3	1.57	0.85
1:A:69:LEU:HD22	1:A:347:ILE:HD12	1.59	0.83
1:A:376:ASP:O	1:A:380:ARG:HB2	1.81	0.81
1:A:540:ALA:CB	1:A:640:LYS:HG3	2.11	0.80
1:A:690:LYS:O	1:A:694:ILE:HG12	1.84	0.77
1:A:142:ASN:HD22	1:A:145:GLY:H	1.34	0.76
1:A:627:LYS:CE	1:A:627:LYS:CG	2.66	0.73
1:A:540:ALA:HB3	1:A:640:LYS:CG	2.21	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:LYS:NZ	1:A:374:LYS:CD	2.53	0.70
1:A:517:ARG:HH11	1:A:517:ARG:CG	2.05	0.68
1:A:141:ASP:O	1:A:205:ARG:CZ	2.43	0.67
1:A:208:MET:HE2	1:A:212:ILE:HG21	1.77	0.67
1:A:484:LEU:HD13	1:A:698:ARG:HD3	1.77	0.67
1:A:208:MET:HE2	1:A:212:ILE:HD13	1.78	0.66
1:A:486:LYS:HG3	1:A:487:PRO:CD	2.16	0.66
1:A:627:LYS:CD	1:A:627:LYS:NZ	2.59	0.65
1:A:142:ASN:ND2	1:A:145:GLY:H	1.97	0.63
1:A:61:HIS:CD2	1:A:62:GLU:H	2.17	0.62
1:A:196:GLN:O	1:A:197:GLU:C	2.42	0.61
1:A:431:GLN:NE2	1:A:435:GLN:HE21	1.97	0.61
1:A:118:TYR:HD1	1:A:171:LEU:HD12	1.66	0.60
1:A:22:VAL:HA	1:A:25:ILE:HD12	1.83	0.60
1:A:233:TYR:O	1:A:234:LYS:HG2	2.02	0.60
1:A:385:MET:HE1	1:A:569:LEU:HB3	1.84	0.60
1:A:113:LEU:HD11	1:A:173:ARG:HG3	1.83	0.59
1:A:197:GLU:OE2	1:A:248:LYS:NZ	2.34	0.59
1:A:431:GLN:HE21	1:A:435:GLN:HE21	1.50	0.59
1:A:69:LEU:CD2	1:A:347:ILE:HD12	2.30	0.59
1:A:517:ARG:CG	1:A:517:ARG:NH1	2.66	0.58
1:A:409:SER:OG	1:A:411:ASP:HB2	2.05	0.56
1:A:192:LEU:CD1	1:A:338:VAL:HG13	2.35	0.56
1:A:362:SER:O	1:A:594:ILE:HD12	2.07	0.55
1:A:376:ASP:O	1:A:380:ARG:HD3	2.07	0.54
1:A:190:ILE:HG12	1:A:245:LEU:HD21	1.89	0.53
1:A:192:LEU:HD11	1:A:338:VAL:HG13	1.91	0.53
1:A:208:MET:CE	1:A:212:ILE:HD13	2.38	0.53
1:A:78:ILE:HD11	1:A:355:PHE:CD1	2.43	0.53
1:A:590:GLN:HG3	1:A:591:ASP:N	2.23	0.53
1:A:373:ARG:NH2	1:A:597:GLU:HG2	2.24	0.52
1:A:367:LEU:HD21	1:A:581:THR:OG1	2.09	0.52
1:A:59:ASP:HB3	1:A:64:ALA:HB2	1.92	0.52
1:A:171:LEU:HD22	1:A:175:PRO:HG3	1.90	0.52
1:A:29:TYR:HB3	1:A:30:PRO:HD2	1.92	0.51
1:A:359:VAL:HB	1:A:364:GLN:HE22	1.76	0.51
1:A:81:LYS:O	1:A:82:PHE:C	2.53	0.51
1:A:10:LYS:HE2	1:A:122:CYS:SG	2.51	0.51
1:A:17:LEU:HD13	1:A:46:LEU:HG	1.93	0.51
1:A:530:LEU:O	1:A:534:ILE:HG13	2.10	0.51
1:A:542:THR:C	1:A:544:GLN:H	2.19	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:GLU:O	1:A:464:THR:HG23	2.12	0.50
1:A:409:SER:OG	1:A:411:ASP:CB	2.60	0.50
1:A:501:LYS:N	1:A:501:LYS:HD3	2.26	0.50
1:A:12:THR:HG21	1:A:50:ARG:HH22	1.77	0.49
1:A:385:MET:HE1	1:A:569:LEU:CB	2.42	0.49
1:A:511:GLU:O	1:A:515:SER:HB2	2.13	0.49
1:A:84:PHE:C	1:A:86:GLU:H	2.20	0.49
1:A:400:ASN:OD1	1:A:403:ALA:HB3	2.13	0.49
1:A:424:VAL:HG21	1:A:523:LEU:CD2	2.43	0.49
1:A:426:GLU:C	1:A:428:GLY:H	2.20	0.48
1:A:525:LYS:HD3	1:A:529:GLU:OE2	2.14	0.48
1:A:624:GLY:O	1:A:628:ASN:ND2	2.47	0.48
1:A:3:THR:CG2	1:A:294:TYR:O	2.55	0.47
1:A:99:PHE:HB2	1:A:101:LYS:HG3	1.96	0.47
1:A:385:MET:HE2	1:A:566:ARG:CB	2.45	0.46
1:A:322:ARG:NH1	1:A:323:VAL:O	2.48	0.46
1:A:61:HIS:HD2	1:A:62:GLU:H	1.59	0.46
1:A:596:GLU:O	1:A:597:GLU:C	2.58	0.46
1:A:385:MET:HE2	1:A:566:ARG:HB3	1.98	0.45
1:A:636:PHE:HA	1:A:639:MET:SD	2.57	0.45
1:A:197:GLU:O	1:A:200:PHE:HB3	2.16	0.45
1:A:439:GLU:HA	1:A:442:GLU:OE2	2.17	0.45
1:A:477:GLN:HG2	1:A:477:GLN:O	2.15	0.45
1:A:478:ARG:O	1:A:479:THR:C	2.60	0.44
1:A:233:TYR:C	1:A:234:LYS:HG2	2.42	0.44
1:A:79:GLU:N	1:A:80:PRO:HD2	2.33	0.44
1:A:12:THR:HG23	1:A:95:TRP:CD2	2.53	0.44
1:A:163:ILE:O	1:A:167:VAL:HG22	2.16	0.44
1:A:378:VAL:O	1:A:382:ILE:HG12	2.17	0.44
1:A:128:ALA:HB2	1:A:159:ALA:HB3	1.99	0.44
1:A:71:TYR:OH	1:A:119:GLU:OE2	2.22	0.43
1:A:197:GLU:CD	1:A:248:LYS:HZ1	2.25	0.43
1:A:583:LYS:HE2	1:A:583:LYS:HB2	1.91	0.43
1:A:118:TYR:CD1	1:A:171:LEU:HD12	2.50	0.42
1:A:399:LEU:O	1:A:401:LEU:HG	2.19	0.42
1:A:148:ILE:HG23	1:A:152:HIS:CD2	2.54	0.42
1:A:517:ARG:HG3	1:A:517:ARG:NH1	2.15	0.42
1:A:12:THR:HG22	1:A:94:THR:O	2.19	0.42
1:A:12:THR:OG1	1:A:54:VAL:HG13	2.19	0.42
1:A:649:ARG:HG2	1:A:653:LEU:HD22	2.02	0.42
1:A:59:ASP:HB3	1:A:64:ALA:CB	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:GLU:HA	1:A:65:LEU:HB3	2.02	0.42
1:A:125:PHE:C	1:A:125:PHE:CD2	2.98	0.41
1:A:447:ASN:HD22	1:A:684:LEU:HD12	1.85	0.41
1:A:454:SER:O	1:A:457:LEU:HB2	2.20	0.41
1:A:386:ARG:NH2	1:A:570:GLU:OE2	2.52	0.41
1:A:279:HIS:HB2	1:A:326:LEU:HD21	2.00	0.41
1:A:633:HIS:O	1:A:636:PHE:N	2.54	0.41
1:A:19:LYS:N	1:A:20:PRO:HD2	2.36	0.40
1:A:172:SER:O	1:A:173:ARG:HG2	2.21	0.40
2:B:34:GLU:O	2:B:34:GLU:HG3	2.19	0.40
1:A:679:GLU:O	1:A:683:ILE:HG12	2.20	0.40
1:A:17:LEU:HD23	1:A:91:LEU:HD21	2.03	0.40
1:A:584:PHE:HB3	1:A:594:ILE:HD13	2.04	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	695/697 (100%)	638 (92%)	49 (7%)	8 (1%)	10	23
2	B	9/11 (82%)	7 (78%)	2 (22%)	0	100	100
All	All	704/708 (99%)	645 (92%)	51 (7%)	8 (1%)	11	25

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	87	ASN
1	A	544	GLN
1	A	85	SER
1	A	82	PHE

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Mol	Chain	Res	Type
1	A	473	LYS
1	A	83	PRO
1	A	25	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	600/600 (100%)	528 (88%)	72 (12%)	5	10
2	B	11/11 (100%)	9 (82%)	2 (18%)	2	3
All	All	611/611 (100%)	537 (88%)	74 (12%)	5	10

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LEU
1	A	12	THR
1	A	13	SER
1	A	22	VAL
1	A	54	VAL
1	A	75	ILE
1	A	78	ILE
1	A	82	PHE
1	A	86	GLU
1	A	89	ILE
1	A	91	LEU
1	A	92	THR
1	A	104	LEU
1	A	109	VAL
1	A	113	LEU
1	A	115	SER
1	A	140	LEU
1	A	142	ASN
1	A	144	GLU
1	A	166	THR

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Mol	Chain	Res	Type
1	A	167	VAL
1	A	169	SER
1	A	173	ARG
1	A	180	SER
1	A	184	VAL
1	A	201	LEU
1	A	234	LYS
1	A	292	ASP
1	A	297	VAL
1	A	312	LYS
1	A	322	ARG
1	A	338	VAL
1	A	352	THR
1	A	357	LYS
1	A	364	GLN
1	A	367	LEU
1	A	371	ASN
1	A	380	ARG
1	A	408	VAL
1	A	416	SER
1	A	417	ILE
1	A	435	GLN
1	A	438	LYS
1	A	449	GLU
1	A	464	THR
1	A	471	LYS
1	A	473	LYS
1	A	477	GLN
1	A	484	LEU
1	A	493	THR
1	A	501	LYS
1	A	508	GLN
1	A	514	GLN
1	A	517	ARG
1	A	542	THR
1	A	569	LEU
1	A	573	LEU
1	A	596	GLU
1	A	597	GLU
1	A	623	GLU
1	A	628	ASN
1	A	629	ILE

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Mol	Chain	Res	Type
1	A	634	GLN
1	A	644	ASN
1	A	648	LEU
1	A	650	GLU
1	A	651	GLU
1	A	653	LEU
1	A	654	LYS
1	A	679	GLU
1	A	682	GLU
1	A	688	GLN
2	B	34	GLU
2	B	44	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	61	HIS
1	A	74	GLN
1	A	142	ASN
1	A	219	GLN
1	A	364	GLN
1	A	415	GLN
1	A	435	GLN
1	A	447	ASN
1	A	514	GLN
1	A	544	GLN
1	A	571	ASN
1	A	628	ASN
1	A	633	HIS
1	A	647	ASN
1	A	678	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [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	697/697 (100%)	0.38	27 (3%) 43 38	34, 85, 140, 181	2 (0%)
2	B	11/11 (100%)	1.00	1 (9%) 15 11	86, 109, 165, 166	0
All	All	708/708 (100%)	0.39	28 (3%) 42 37	34, 86, 141, 181	2 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	453	GLU	4.8
1	A	104	LEU	3.9
1	A	102	GLY	3.3
1	A	2	ALA	3.2
1	A	37	ALA	3.2
1	A	144	GLU	3.1
1	A	554	SER	3.0
1	A	172	SER	2.9
1	A	431	GLN	2.7
1	A	321	ASP	2.7
1	A	105	PHE	2.6
1	A	546	SER	2.5
1	A	28	THR	2.4
1	A	548	VAL	2.4
1	A	593	VAL	2.4
1	A	598	ALA	2.3
1	A	680	LEU	2.3
1	A	29	TYR	2.2
2	B	38	LEU	2.2
1	A	631	VAL	2.2
1	A	24	PHE	2.2
1	A	602	THR	2.1
1	A	632	SER	2.1
1	A	442	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	361	VAL	2.1
1	A	490	ALA	2.0
1	A	540	ALA	2.0
1	A	493	THR	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.