



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:59 AM UTC

PDB ID : 6JEH / pdb_00006jeh
Title : Crystal structure of calcium free human gelsolin amyloid mutant D187Y
Authors : Zorgati, H.; Robinson, R.C.
Deposited on : 2019-02-05
Resolution : 2.95 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

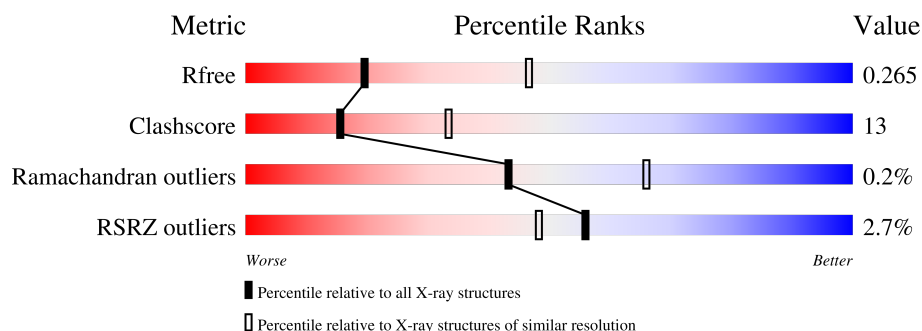
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	727	 3% 70% 27% .
1	B	727	 2% 73% 23% .

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 11005 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 Gelsolin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	708	Total	C	N	O	S	0	0	0
			5534	3497	967	1054	16			
1	B	700	Total	C	N	O	S	0	0	0
			5471	3459	956	1040	16			

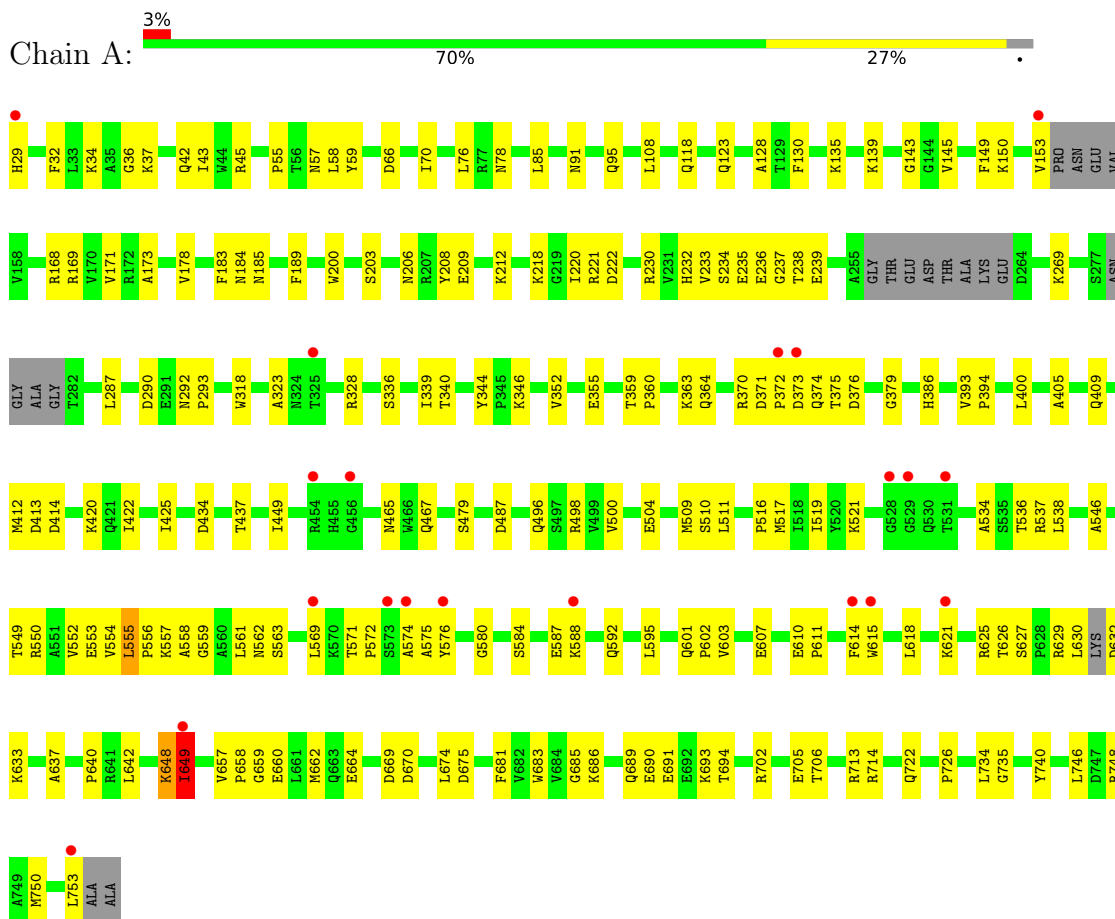
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	187	TYR	ASP	engineered mutation	UNP P06396
B	187	TYR	ASP	engineered mutation	UNP P06396

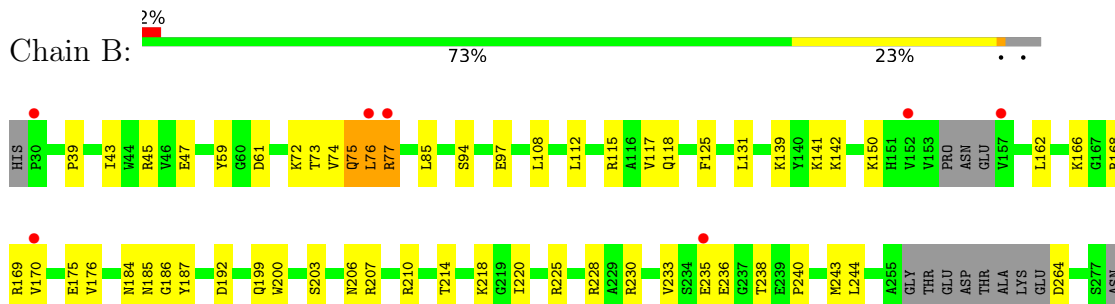
3 Residue-property plots

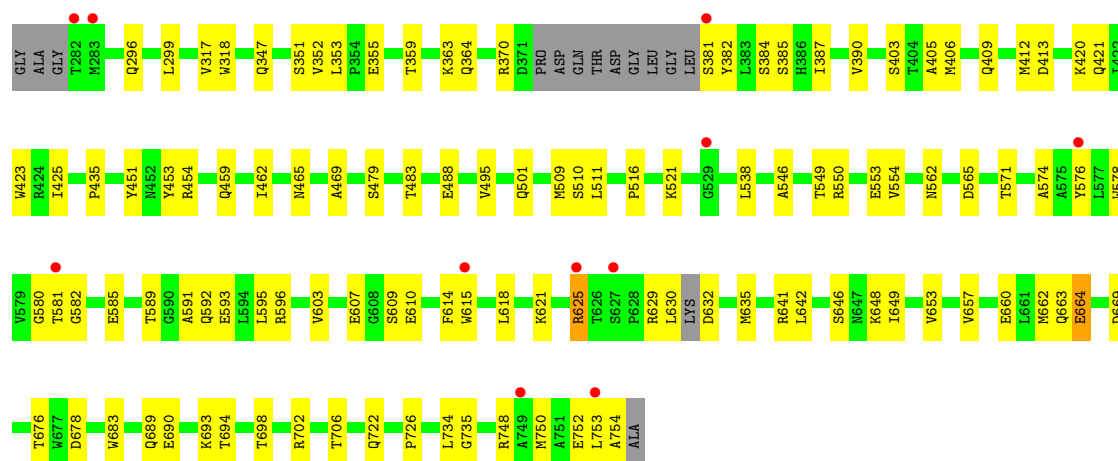
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Gelsolin



• Molecule 1: Gelsolin





4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	169.23Å 169.23Å 151.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.94 – 2.95 19.94 – 2.95	Depositor EDS
% Data completeness (in resolution range)	98.9 (19.94-2.95) 98.5 (19.94-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 2.93Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.193 , 0.249 (Not available) , 0.265	Depositor DCC
R_{free} test set	2386 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	58.5	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 53.1	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	0.93	EDS
Total number of atoms	11005	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.16% 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	0.49	1/5660 (0.0%)	0.78	6/7664 (0.1%)
1	B	0.51	2/5594 (0.0%)	0.80	9/7571 (0.1%)
All	All	0.50	3/11254 (0.0%)	0.79	15/15235 (0.1%)

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
1	B	0	2
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	585	GLU	CB-CG	6.24	1.71	1.52
1	A	649	ILE	CB-CG2	5.78	1.71	1.52
1	B	230	ARG	CB-CG	-5.43	1.36	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	664	GLU	CA-CB-CG	-11.44	91.22	114.10
1	B	230	ARG	NE-CZ-NH2	-9.35	110.79	119.20
1	B	230	ARG	NE-CZ-NH1	6.86	128.36	121.50
1	A	555	LEU	CB-CG-CD2	-6.68	90.65	110.70
1	A	621	LYS	CB-CA-C	6.30	121.10	109.70
1	B	625	ARG	CB-CG-CD	6.23	125.63	111.30
1	B	230	ARG	CG-CD-NE	-6.02	98.76	112.00
1	A	588	LYS	CB-CG-CD	-6.00	97.49	111.30
1	A	649	ILE	CG1-CB-CG2	5.94	128.53	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	664	GLU	N-CA-CB	-5.81	101.27	110.28
1	A	621	LYS	N-CA-CB	-5.68	100.97	110.80
1	A	691	GLU	CB-CA-C	-5.61	101.32	110.85
1	B	77	ARG	CA-C-N	-5.40	114.01	122.49
1	B	77	ARG	C-N-CA	-5.40	114.01	122.49
1	B	585	GLU	CA-CB-CG	5.18	124.45	114.10

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	648	LYS	Peptide
1	B	488	GLU	Sidechain
1	B	75	GLN	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	5534	0	5390	164	0
1	B	5471	0	5332	134	0
All	All	11005	0	10722	292	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 13.

All (292) 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:370:ARG:NH1	1:A:627:SER:HB3	1.10	1.42
1:A:370:ARG:NH1	1:A:627:SER:CB	1.85	1.39
1:A:648:LYS:O	1:A:649:ILE:HG22	1.37	1.20
1:A:370:ARG:HH11	1:A:627:SER:CB	1.56	1.11
1:B:521:LYS:HG3	1:B:553:GLU:HG2	1.39	1.04
1:A:630:LEU:O	1:A:632:ASP:N	1.92	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:465:ASN:HB2	1:B:483:THR:HG21	1.37	1.01
1:B:75:GLN:O	1:B:76:LEU:O	1.80	0.99
1:A:34:LYS:O	1:A:37:LYS:NZ	1.98	0.96
1:B:630:LEU:O	1:B:632:ASP:N	2.01	0.94
1:A:43:ILE:HG12	1:A:70:ILE:HD12	1.53	0.90
1:A:648:LYS:O	1:A:649:ILE:CG2	2.19	0.90
1:A:185:ASN:ND2	1:A:236:GLU:OE1	2.04	0.89
1:A:713:ARG:H	1:A:713:ARG:HD3	1.38	0.89
1:A:55:PRO:HG2	1:A:58:LEU:HD12	1.56	0.88
1:A:536:THR:HG22	1:A:571:THR:HA	1.54	0.87
1:A:746:LEU:HD11	1:A:750:MET:HE3	1.59	0.85
1:A:135:LYS:HE2	1:A:662:MET:HE2	1.58	0.84
1:A:686:LYS:H	1:A:722:GLN:NE2	1.75	0.84
1:A:521:LYS:HG3	1:A:553:GLU:HG2	1.63	0.81
1:B:75:GLN:O	1:B:76:LEU:C	2.22	0.79
1:B:425:ILE:HD13	1:B:479:SER:HB3	1.65	0.79
1:A:36:GLY:H	1:A:42:GLN:HE22	1.32	0.78
1:A:420:LYS:NZ	1:A:511:LEU:O	2.16	0.77
1:A:379:GLY:H	1:A:633:LYS:HE2	1.50	0.77
1:B:45:ARG:NH1	1:B:47:GLU:OE2	2.17	0.77
1:A:519:ILE:HD12	1:A:552:VAL:HG12	1.66	0.76
1:B:576:TYR:HE2	1:B:614:PHE:HA	1.48	0.76
1:A:370:ARG:HH12	1:A:627:SER:CB	1.74	0.75
1:B:571:THR:HG23	1:B:574:ALA:H	1.53	0.74
1:A:625:ARG:NH1	1:A:627:SER:OG	2.20	0.73
1:A:269:LYS:HE2	1:A:705:GLU:O	1.88	0.72
1:A:393:VAL:HA	1:B:459:GLN:NE2	2.05	0.72
1:A:370:ARG:HH12	1:A:627:SER:HB3	0.83	0.72
1:A:70:ILE:HG13	1:A:130:PHE:HE1	1.54	0.72
1:B:185:ASN:ND2	1:B:236:GLU:OE2	2.22	0.71
1:A:371:ASP:OD1	1:A:372:PRO:HD2	1.89	0.71
1:B:296:GLN:HE22	1:B:370:ARG:H	1.38	0.70
1:B:364:GLN:HE22	1:B:641:ARG:NH2	1.89	0.70
1:A:118:GLN:HB2	1:A:352:VAL:HG22	1.74	0.70
1:A:393:VAL:HA	1:B:459:GLN:HE22	1.57	0.70
1:B:576:TYR:CE2	1:B:614:PHE:HA	2.27	0.69
1:B:752:GLU:HG2	1:B:753:LEU:HD22	1.74	0.69
1:B:615:TRP:CD2	1:B:621:LYS:HD2	2.27	0.69
1:B:39:PRO:HA	1:B:73:THR:O	1.93	0.68
1:B:479:SER:O	1:B:483:THR:HG23	1.93	0.68
1:B:214:THR:O	1:B:218:LYS:HG2	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:ASN:H	1:A:123:GLN:NE2	1.92	0.68
1:B:318:TRP:CZ2	1:B:355:GLU:HB2	2.29	0.67
1:A:169:ARG:HD2	1:A:657:VAL:HG12	1.75	0.67
1:A:318:TRP:CZ2	1:A:355:GLU:HB2	2.29	0.67
1:A:139:LYS:HE2	1:A:220:ILE:HG12	1.77	0.67
1:B:207:ARG:HA	1:B:210:ARG:HG3	1.76	0.67
1:B:192:ASP:OD2	1:B:225:ARG:NH2	2.27	0.67
1:A:538:LEU:HG	1:A:556:PRO:HA	1.78	0.66
1:B:185:ASN:OD1	1:B:203:SER:N	2.29	0.66
1:A:571:THR:OG1	1:A:572:PRO:HD2	1.96	0.66
1:A:45:ARG:HH22	1:A:150:LYS:HA	1.61	0.65
1:B:139:LYS:HE2	1:B:220:ILE:HG12	1.79	0.65
1:B:364:GLN:HE22	1:B:641:ARG:HH22	1.43	0.65
1:A:610:GLU:N	1:A:610:GLU:OE1	2.30	0.64
1:B:75:GLN:C	1:B:76:LEU:O	2.41	0.64
1:A:36:GLY:H	1:A:42:GLN:NE2	1.94	0.63
1:B:199:GLN:HE22	1:B:214:THR:HA	1.63	0.63
1:A:169:ARG:NH2	1:A:670:ASP:OD2	2.28	0.63
1:A:509:MET:HG2	1:A:550:ARG:HD3	1.79	0.63
1:B:387:ILE:HD11	1:B:635:MET:HE3	1.80	0.63
1:A:517:MET:HE2	1:A:519:ILE:HD11	1.81	0.62
1:B:184:ASN:OD1	1:B:185:ASN:N	2.33	0.62
1:B:420:LYS:HE2	1:B:511:LEU:HB3	1.81	0.62
1:A:143:GLY:HA3	1:A:149:PHE:HD2	1.65	0.62
1:A:323:ALA:O	1:A:328:ARG:NH1	2.32	0.61
1:A:413:ASP:OD2	1:A:726:PRO:HB2	1.99	0.61
1:A:558:ALA:HA	1:A:561:LEU:HD23	1.81	0.61
1:B:413:ASP:OD2	1:B:726:PRO:HB2	2.00	0.61
1:B:206:ASN:OD1	1:B:207:ARG:N	2.35	0.60
1:A:576:TYR:CE2	1:A:614:PHE:HB2	2.36	0.60
1:B:593:GLU:OE1	1:B:596:ARG:NH2	2.35	0.59
1:A:394:PRO:HD3	1:B:459:GLN:NE2	2.18	0.59
1:B:516:PRO:HB3	1:B:549:THR:HB	1.85	0.59
1:A:642:LEU:HD23	1:A:657:VAL:HG21	1.84	0.59
1:A:168:ARG:HH12	1:A:208:TYR:HB3	1.68	0.59
1:B:698:THR:O	1:B:702:ARG:HG2	2.02	0.58
1:A:713:ARG:H	1:A:713:ARG:CD	2.15	0.58
1:B:702:ARG:O	1:B:706:THR:HG23	2.02	0.58
1:A:557:LYS:HG3	1:A:559:GLY:H	1.67	0.58
1:B:85:LEU:HD11	1:B:108:LEU:HB3	1.85	0.58
1:B:74:VAL:HG21	1:B:385:SER:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:425:ILE:HD13	1:B:479:SER:CB	2.32	0.58
1:A:370:ARG:HH11	1:A:627:SER:HB2	1.63	0.57
1:B:43:ILE:HD12	1:B:59:TYR:CG	2.40	0.57
1:A:504:GLU:OE1	1:A:550:ARG:NH1	2.38	0.56
1:A:584:SER:OG	1:A:587:GLU:OE2	2.20	0.56
1:B:683:TRP:CZ2	1:B:722:GLN:HB2	2.40	0.56
1:A:85:LEU:HD11	1:A:108:LEU:HB3	1.88	0.56
1:A:434:ASP:HB3	1:A:437:THR:HG22	1.87	0.56
1:B:61:ASP:HB3	1:B:141:LYS:HD2	1.87	0.56
1:B:615:TRP:CD1	1:B:618:LEU:HD12	2.40	0.56
1:B:168:ARG:NH2	1:B:669:ASP:OD1	2.39	0.56
1:A:557:LYS:HE3	1:A:558:ALA:H	1.70	0.56
1:A:664:GLU:OE1	1:A:664:GLU:N	2.31	0.56
1:A:734:LEU:HD12	1:A:735:GLY:N	2.21	0.56
1:B:184:ASN:OD1	1:B:186:GLY:N	2.27	0.55
1:A:70:ILE:HG13	1:A:130:PHE:CE1	2.40	0.55
1:B:592:GLN:HA	1:B:595:LEU:HD23	1.89	0.55
1:A:43:ILE:HD12	1:A:59:TYR:CG	2.43	0.54
1:A:237:GLY:N	1:A:239:GLU:OE2	2.40	0.54
1:A:57:ASN:HD21	1:A:153:VAL:HG11	1.73	0.54
1:A:370:ARG:NH1	1:A:627:SER:OG	2.36	0.53
1:B:405:ALA:O	1:B:409:GLN:HG3	2.08	0.53
1:A:425:ILE:HG13	1:A:479:SER:HB3	1.91	0.53
1:B:591:ALA:O	1:B:595:LEU:CD2	2.56	0.53
1:A:183:PHE:HA	1:A:189:PHE:CZ	2.43	0.53
1:A:557:LYS:HG3	1:A:558:ALA:N	2.23	0.53
1:A:422:ILE:HG12	1:A:449:ILE:CG1	2.39	0.53
1:B:509:MET:HG2	1:B:550:ARG:HB3	1.90	0.53
1:A:690:GLU:O	1:A:694:THR:HG23	2.09	0.53
1:A:595:LEU:HD13	1:A:602:PRO:HB3	1.91	0.53
1:B:538:LEU:HB3	1:B:554:VAL:HG23	1.90	0.52
1:A:218:LYS:NZ	1:A:222:ASP:OD2	2.40	0.52
1:A:394:PRO:HD3	1:B:459:GLN:HE21	1.74	0.52
1:A:235:GLU:HG2	1:A:236:GLU:H	1.73	0.52
1:A:516:PRO:HB3	1:A:549:THR:OG1	2.07	0.52
1:A:509:MET:CE	1:A:550:ARG:HE	2.23	0.52
1:B:465:ASN:CB	1:B:483:THR:HG21	2.24	0.52
1:B:454:ARG:HG2	1:B:454:ARG:HH11	1.75	0.51
1:B:387:ILE:O	1:B:387:ILE:HG13	2.10	0.51
1:A:95:GLN:H	1:A:95:GLN:CD	2.18	0.51
1:A:171:VAL:HG21	1:A:212:LYS:HB3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:662:MET:HB3	1:B:664:GLU:OE1	2.10	0.51
1:B:317:VAL:HB	1:B:352:VAL:HG12	1.93	0.51
1:A:702:ARG:O	1:A:706:THR:HG23	2.10	0.51
1:B:118:GLN:HB2	1:B:352:VAL:HG22	1.93	0.51
1:A:400:LEU:CD2	1:A:414:ASP:HA	2.41	0.51
1:B:646:SER:OG	1:B:648:LYS:HG2	2.11	0.51
1:A:686:LYS:H	1:A:722:GLN:HE21	1.55	0.51
1:A:394:PRO:CD	1:B:459:GLN:HE21	2.23	0.51
1:A:405:ALA:O	1:A:409:GLN:HG3	2.11	0.51
1:A:576:TYR:HD1	1:A:603:VAL:HG22	1.76	0.50
1:B:142:LYS:HE3	1:B:175:GLU:HG2	1.93	0.50
1:B:469:ALA:H	1:B:501:GLN:HE21	1.59	0.50
1:B:581:THR:HG23	1:B:607:GLU:OE2	2.11	0.50
1:B:615:TRP:CE3	1:B:621:LYS:HD2	2.46	0.50
1:A:233:VAL:HG11	1:A:750:MET:HE2	1.94	0.50
1:A:336:SER:O	1:A:340:THR:HG23	2.12	0.50
1:A:519:ILE:CD1	1:A:552:VAL:HG12	2.40	0.50
1:A:412:MET:HE2	1:A:546:ALA:HB2	1.94	0.50
1:A:422:ILE:HG12	1:A:449:ILE:HG13	1.93	0.49
1:B:168:ARG:HH12	1:B:206:ASN:ND2	2.10	0.49
1:B:387:ILE:HD12	1:B:390:VAL:HG21	1.94	0.49
1:B:591:ALA:O	1:B:595:LEU:HD23	2.13	0.49
1:A:412:MET:HE1	1:A:510:SER:OG	2.11	0.49
1:B:117:VAL:HG13	1:B:353:LEU:HD23	1.93	0.49
1:A:563:SER:OG	1:A:626:THR:HG22	2.12	0.49
1:A:610:GLU:HG2	1:A:611:PRO:O	2.12	0.49
1:B:39:PRO:HA	1:B:73:THR:HG22	1.94	0.49
1:A:91:ASN:H	1:A:123:GLN:HE21	1.59	0.49
1:A:576:TYR:CD2	1:A:614:PHE:HD1	2.31	0.49
1:A:648:LYS:C	1:A:649:ILE:HG22	2.29	0.48
1:A:648:LYS:HG2	1:A:649:ILE:H	1.78	0.48
1:B:403:SER:OG	1:B:406:MET:HB3	2.14	0.48
1:B:565:ASP:OD2	1:B:629:ARG:NH2	2.43	0.48
1:A:37:LYS:H	1:A:37:LYS:HD3	1.79	0.48
1:A:465:ASN:HD21	1:A:467:GLN:HE21	1.61	0.48
1:A:562:ASN:ND2	1:A:629:ARG:HH22	2.11	0.48
1:B:200:TRP:HB2	1:B:243:MET:HE1	1.96	0.48
1:B:609:SER:N	1:B:610:GLU:OE1	2.47	0.48
1:A:43:ILE:HG23	1:A:70:ILE:CD1	2.44	0.48
1:B:689:GLN:O	1:B:693:LYS:HG3	2.14	0.48
1:A:713:ARG:HG2	1:A:714:ARG:HG3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:683:TRP:CZ2	1:A:722:GLN:HB2	2.49	0.48
1:B:94:SER:HB3	1:B:97:GLU:HB2	1.95	0.48
1:B:734:LEU:HD12	1:B:735:GLY:H	1.79	0.48
1:A:235:GLU:H	1:A:238:THR:CG2	2.27	0.47
1:A:465:ASN:ND2	1:A:467:GLN:HE21	2.12	0.47
1:B:168:ARG:HG3	1:B:168:ARG:HH11	1.79	0.47
1:B:642:LEU:O	1:B:657:VAL:HG13	2.14	0.47
1:A:36:GLY:N	1:A:42:GLN:HE22	2.08	0.47
1:A:373:ASP:O	1:A:375:THR:N	2.46	0.47
1:B:77:ARG:CG	1:B:77:ARG:HH21	2.27	0.47
1:A:76:LEU:O	1:A:78:ASN:N	2.42	0.47
1:A:534:ALA:HB3	1:A:537:ARG:HG2	1.96	0.47
1:A:139:LYS:HG2	1:A:173:ALA:HB3	1.95	0.47
1:A:498:ARG:NH1	1:A:500:VAL:HG13	2.30	0.47
1:A:558:ALA:N	1:A:618:LEU:HD12	2.30	0.47
1:B:77:ARG:NH2	1:B:77:ARG:HG3	2.29	0.47
1:B:233:VAL:HG11	1:B:750:MET:SD	2.55	0.47
1:A:576:TYR:HE2	1:A:614:PHE:HA	1.80	0.47
1:B:125:PHE:CE1	1:B:347:GLN:HG2	2.50	0.47
1:A:400:LEU:HD22	1:A:414:ASP:HA	1.97	0.46
1:B:421:GLN:HG2	1:B:423:TRP:CZ2	2.50	0.46
1:A:360:PRO:O	1:A:364:GLN:HG3	2.15	0.46
1:B:77:ARG:HG3	1:B:77:ARG:O	2.15	0.46
1:B:225:ARG:O	1:B:228:ARG:HB2	2.14	0.46
1:A:168:ARG:NH2	1:A:669:ASP:OD1	2.48	0.46
1:B:169:ARG:H	1:B:169:ARG:HG2	1.43	0.46
1:A:580:GLY:HA2	1:A:607:GLU:HB2	1.97	0.46
1:A:206:ASN:ND2	1:A:209:GLU:HG3	2.30	0.46
1:B:131:LEU:HD13	1:B:170:VAL:CG1	2.45	0.46
1:B:210:ARG:O	1:B:214:THR:HG23	2.16	0.46
1:A:221:ARG:HH12	1:A:753:LEU:HB3	1.80	0.46
1:A:558:ALA:HA	1:A:561:LEU:CD2	2.45	0.46
1:B:384:SER:O	1:B:387:ILE:HG22	2.16	0.46
1:B:359:THR:O	1:B:363:LYS:HG3	2.16	0.46
1:B:562:ASN:HA	1:B:625:ARG:O	2.16	0.46
1:A:287:LEU:HD21	1:A:290:ASP:HB3	1.98	0.46
1:A:689:GLN:O	1:A:693:LYS:HG3	2.16	0.46
1:A:576:TYR:HD2	1:A:614:PHE:HD1	1.63	0.46
1:B:748:ARG:O	1:B:752:GLU:N	2.33	0.46
1:A:230:ARG:HG2	1:A:232:HIS:CE1	2.51	0.45
1:B:462:ILE:HG23	1:B:495:VAL:CG2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:487:ASP:OD2	1:A:496:GLN:NE2	2.49	0.45
1:A:615:TRP:HD1	1:A:615:TRP:HA	1.57	0.45
1:A:740:TYR:O	1:A:748:ARG:NH2	2.45	0.45
1:B:240:PRO:O	1:B:244:LEU:HD12	2.17	0.45
1:B:412:MET:HE2	1:B:546:ALA:HB2	1.98	0.45
1:A:538:LEU:CD1	1:A:554:VAL:HG13	2.48	0.44
1:A:575:ALA:HB3	1:A:601:GLN:O	2.17	0.44
1:B:118:GLN:O	1:B:351:SER:HA	2.17	0.44
1:B:207:ARG:HG2	1:B:210:ARG:NH1	2.32	0.44
1:A:57:ASN:HD21	1:A:153:VAL:CG1	2.30	0.44
1:A:336:SER:OG	1:A:346:LYS:NZ	2.31	0.44
1:A:734:LEU:HD12	1:A:735:GLY:H	1.81	0.44
1:B:72:LYS:HE3	1:B:74:VAL:HG23	1.99	0.44
1:A:66:ASP:OD1	1:A:145:VAL:HG12	2.17	0.44
1:A:346:LYS:HD2	1:A:346:LYS:HA	1.85	0.44
1:A:37:LYS:HD3	1:A:37:LYS:N	2.31	0.44
1:A:339:ILE:HG23	1:A:344:TYR:HB2	1.98	0.44
1:A:360:PRO:HB3	1:A:637:ALA:O	2.18	0.44
1:B:43:ILE:HD12	1:B:59:TYR:CD2	2.53	0.44
1:B:454:ARG:HG2	1:B:454:ARG:NH1	2.32	0.44
1:A:648:LYS:C	1:A:649:ILE:CG2	2.89	0.44
1:A:685:GLY:HA2	1:A:722:GLN:HE21	1.83	0.44
1:B:112:LEU:O	1:B:115:ARG:HG2	2.17	0.44
1:B:690:GLU:O	1:B:694:THR:HG23	2.18	0.44
1:B:299:LEU:HD13	1:B:318:TRP:CE2	2.53	0.44
1:A:569:LEU:HD22	1:A:618:LEU:HD21	1.99	0.43
1:B:451:TYR:CE1	1:B:511:LEU:HD22	2.53	0.43
1:B:648:LYS:O	1:B:649:ILE:HG22	2.19	0.43
1:A:43:ILE:HD12	1:A:59:TYR:CD2	2.53	0.43
1:A:178:VAL:HG23	1:A:178:VAL:O	2.18	0.43
1:B:299:LEU:HD13	1:B:318:TRP:CD2	2.54	0.43
1:B:580:GLY:HA2	1:B:607:GLU:HB2	2.00	0.43
1:A:603:VAL:HG23	1:A:603:VAL:O	2.19	0.43
1:B:663:GLN:N	1:B:664:GLU:OE1	2.52	0.43
1:A:184:ASN:H	1:A:189:PHE:HZ	1.66	0.43
1:A:370:ARG:NH1	1:A:627:SER:HB2	2.13	0.43
1:A:376:ASP:HB2	1:B:435:PRO:HB3	2.00	0.43
1:A:592:GLN:HA	1:A:595:LEU:HD12	2.00	0.43
1:B:750:MET:O	1:B:754:ALA:HB3	2.18	0.42
1:A:571:THR:HG22	1:A:574:ALA:O	2.20	0.42
1:B:412:MET:HE1	1:B:510:SER:OG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:660:GLU:O	1:B:662:MET:HG3	2.19	0.42
1:B:603:VAL:HG23	1:B:603:VAL:O	2.20	0.42
1:A:128:ALA:O	1:A:658:PRO:HB3	2.19	0.42
1:A:640:PRO:HA	1:A:675:ASP:O	2.19	0.42
1:B:166:LYS:HD3	1:B:187:TYR:CE2	2.54	0.42
1:B:578:TRP:HZ2	1:B:607:GLU:HG3	1.84	0.42
1:A:269:LYS:O	1:A:269:LYS:HG3	2.20	0.42
1:A:555:LEU:HA	1:A:556:PRO:HD3	1.74	0.42
1:A:269:LYS:HB2	1:A:269:LYS:HE3	1.78	0.42
1:A:674:LEU:HB3	1:A:681:PHE:HB2	2.02	0.42
1:B:168:ARG:NH1	1:B:206:ASN:ND2	2.68	0.42
1:A:359:THR:O	1:A:363:LYS:HG3	2.19	0.41
1:A:614:PHE:CE2	1:A:615:TRP:NE1	2.88	0.41
1:B:235:GLU:H	1:B:238:THR:CG2	2.32	0.41
1:B:538:LEU:HD23	1:B:538:LEU:HA	1.85	0.41
1:B:589:THR:HA	1:B:592:GLN:HE21	1.85	0.41
1:A:386:HIS:HB2	1:A:659:GLY:C	2.45	0.41
1:A:200:TRP:HA	1:A:234:SER:O	2.21	0.41
1:B:61:ASP:OD1	1:B:139:LYS:NZ	2.39	0.41
1:A:135:LYS:NZ	1:A:658:PRO:O	2.45	0.41
1:A:292:ASN:HB2	1:A:293:PRO:HA	2.02	0.41
1:A:660:GLU:O	1:A:662:MET:HG3	2.20	0.41
1:B:676:THR:HG22	1:B:678:ASP:H	1.85	0.41
1:B:76:LEU:HD12	1:B:76:LEU:HA	1.74	0.41
1:A:95:GLN:OE1	1:A:95:GLN:N	2.54	0.40
1:B:45:ARG:HH22	1:B:150:LYS:HD3	1.86	0.40
1:B:381:SER:OG	1:B:382:TYR:N	2.53	0.40
1:B:451:TYR:HE2	1:B:453:TYR:HB3	1.86	0.40
1:A:233:VAL:HG11	1:A:750:MET:CE	2.51	0.40
1:A:509:MET:HE3	1:A:550:ARG:HE	1.85	0.40
1:B:85:LEU:HD23	1:B:85:LEU:HA	1.82	0.40
1:B:170:VAL:HG23	1:B:170:VAL:O	2.22	0.40
1:A:29:HIS:O	1:A:32:PHE:HB2	2.21	0.40
1:B:77:ARG:CG	1:B:77:ARG:NH2	2.83	0.40
1:B:264:ASP:CG	1:B:653:VAL:HG23	2.47	0.40
1:A:185:ASN:OD1	1:A:203:SER:N	2.54	0.40
1:B:97:GLU:HA	1:B:97:GLU:OE2	2.22	0.40
1:B:162:LEU:HD23	1:B:176:VAL:HG23	2.02	0.40
1:B:581:THR:OG1	1:B:582:GLY:N	2.55	0.40
1:B:625:ARG:HA	1:B:625:ARG:NE	2.37	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	698/727 (96%)	677 (97%)	19 (3%)	2 (0%)	36	59
1	B	688/727 (95%)	667 (97%)	20 (3%)	1 (0%)	48	72
All	All	1386/1454 (95%)	1344 (97%)	39 (3%)	3 (0%)	43	66

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	374	GLN
1	B	76	LEU
1	A	649	ILE

5.3.2 Protein sidechains [i](#)

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

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	708/727 (97%)	-0.14	20 (2%)	55 47	30, 60, 127, 167	0
1	B	700/727 (96%)	-0.25	18 (2%)	57 49	26, 56, 107, 143	0
All	All	1408/1454 (96%)	-0.19	38 (2%)	56 48	26, 58, 117, 167	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	77	ARG	4.8
1	A	576	TYR	3.6
1	B	753	LEU	3.6
1	B	581	THR	3.2
1	A	569	LEU	3.0
1	B	76	LEU	3.0
1	A	372	PRO	3.0
1	A	649	ILE	2.9
1	B	615	TRP	2.9
1	B	157	VAL	2.8
1	A	456	GLY	2.8
1	A	614	PHE	2.8
1	A	529	GLY	2.7
1	B	627	SER	2.7
1	A	621	LYS	2.6
1	B	283	MET	2.6
1	B	381	SER	2.6
1	B	152	VAL	2.6
1	B	749	ALA	2.6
1	B	30	PRO	2.5
1	A	615	TRP	2.5
1	A	153	VAL	2.4
1	B	576	TYR	2.3
1	A	325	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	531	THR	2.3
1	A	753	LEU	2.3
1	A	588	LYS	2.2
1	A	573	SER	2.2
1	A	29	HIS	2.2
1	B	170	VAL	2.1
1	A	373	ASP	2.1
1	B	529	GLY	2.1
1	B	625	ARG	2.1
1	A	528	GLY	2.1
1	B	235	GLU	2.1
1	B	282	THR	2.0
1	A	574	ALA	2.0
1	A	454	ARG	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.