



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:16 AM UTC

PDB ID : 6FVO / pdb_00006fvo
Title : Mutant DNA polymerase sliding clamp from Mycobacterium tuberculosis with bound P7 peptide
Authors : Martiel, I.; Andre, C.; Olieric, V.; Guichard, G.; Burnouf, D.
Deposited on : 2018-03-04
Resolution : 2.69 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:

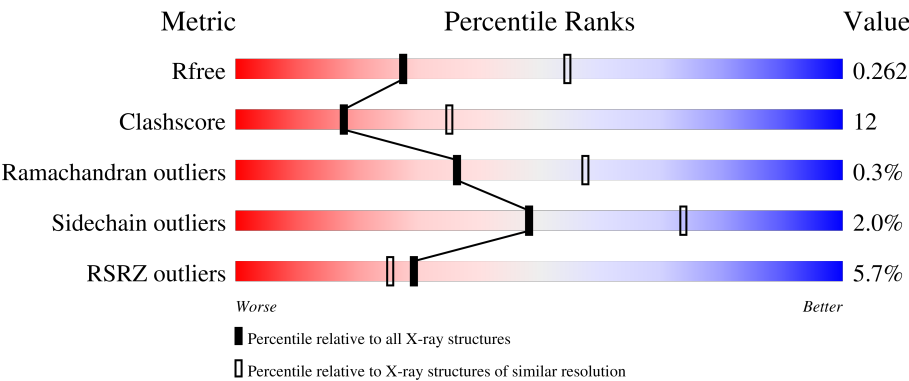
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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	402	6% 72% 24% ..
1	B	402	5% 69% 22% 6%
1	C	402	6% 72% 22% ..
1	D	402	5% 71% 22% 6%
2	H	6	50% 50%

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Mol	Chain	Length	Quality of chain
2	I	6	 83% 17%
2	J	6	 67% 17% 17%
2	K	6	 67% 17% 17%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11359 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 Beta sliding clamp.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	0	0
			2775	1761	469	539	6			
1	B	376	Total	C	N	O	S	0	0	0
			2750	1747	469	528	6			
1	C	389	Total	C	N	O	S	0	0	0
			2794	1772	473	543	6			
1	D	379	Total	C	N	O	S	0	0	0
			2740	1742	462	530	6			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	362	SER	PRO	conflict	UNP P9WNU0
B	362	SER	PRO	conflict	UNP P9WNU0
C	362	SER	PRO	conflict	UNP P9WNU0
D	362	SER	PRO	conflict	UNP P9WNU0

- Molecule 2 is a protein called P7 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	H	6	Total	C	N	O	0	0	0
			51	35	6	10			
2	I	6	Total	C	N	O	0	0	0
			51	35	6	10			
2	J	6	Total	C	N	O	0	0	0
			51	35	6	10			
2	K	6	Total	C	N	O	0	0	0
			51	35	6	10			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	2	Total Ca 2 2	0	0
3	C	1	Total Ca 1 1	0	0
3	D	1	Total Ca 1 1	0	0

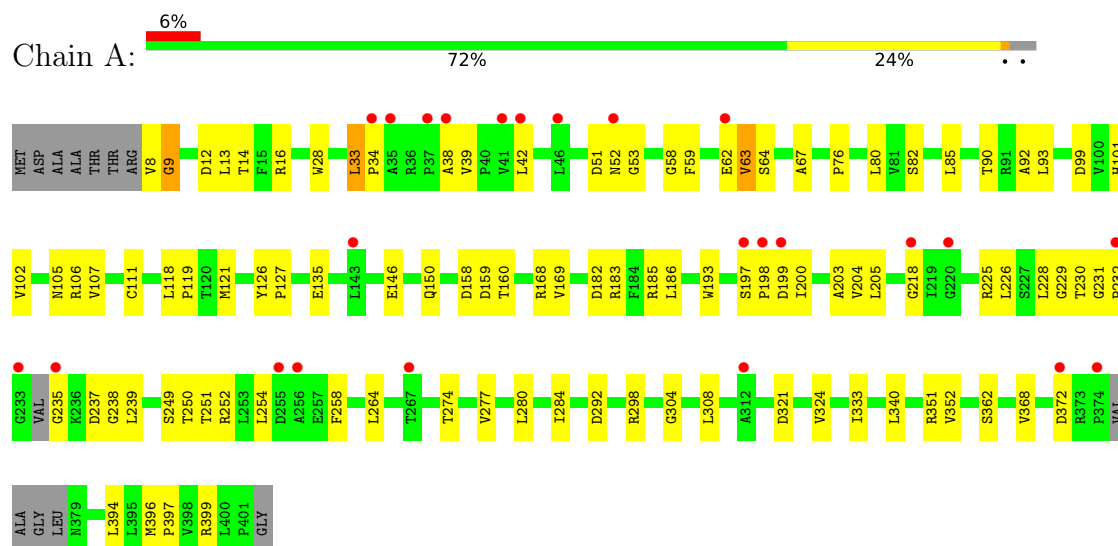
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	24	Total O 24 24	0	0
4	B	29	Total O 29 29	0	0
4	C	25	Total O 25 25	0	0
4	D	14	Total O 14 14	0	0

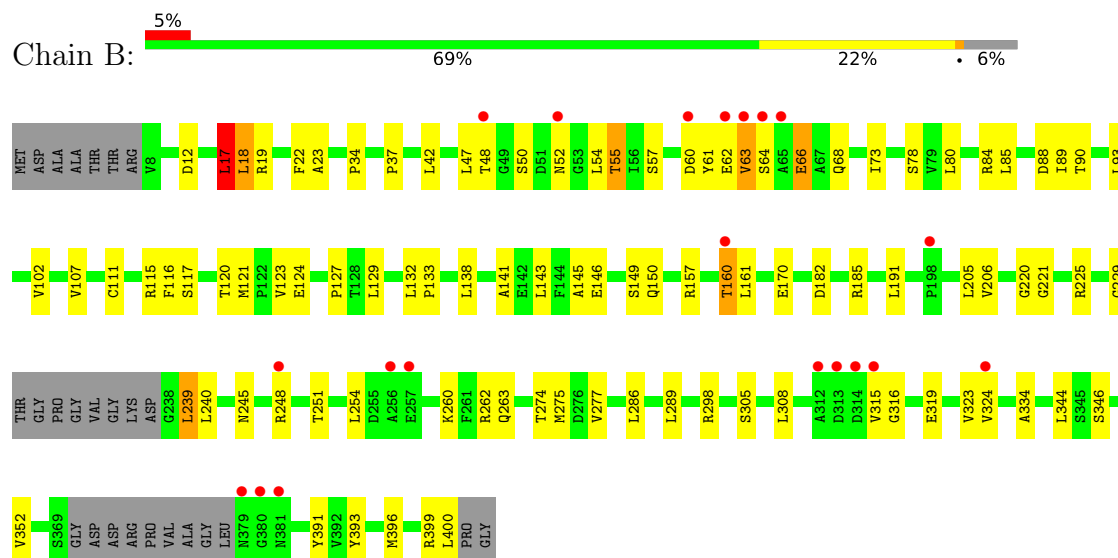
3 Residue-property plots [i](#)

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.

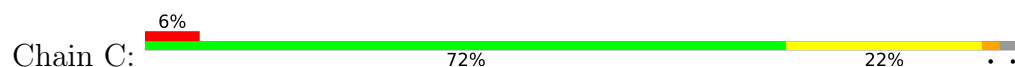
• Molecule 1: Beta sliding clamp

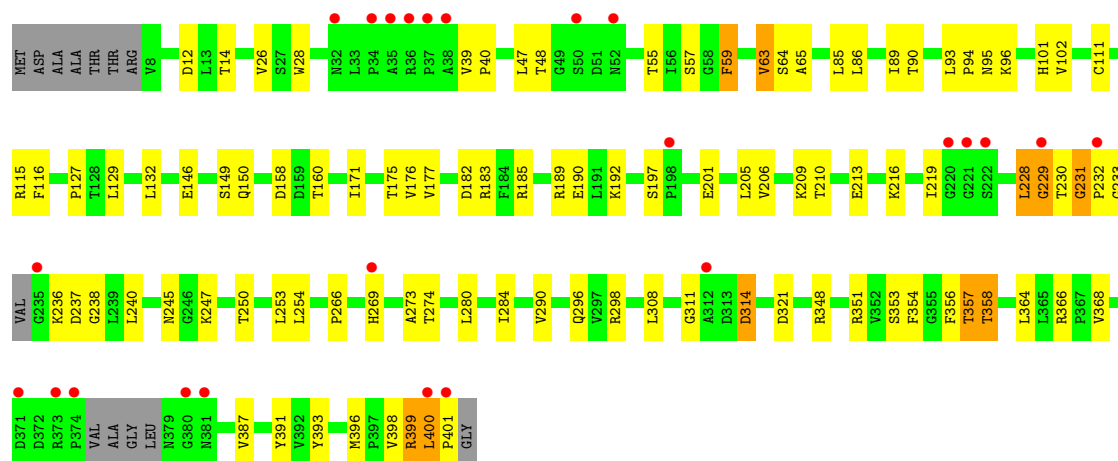


• Molecule 1: Beta sliding clamp

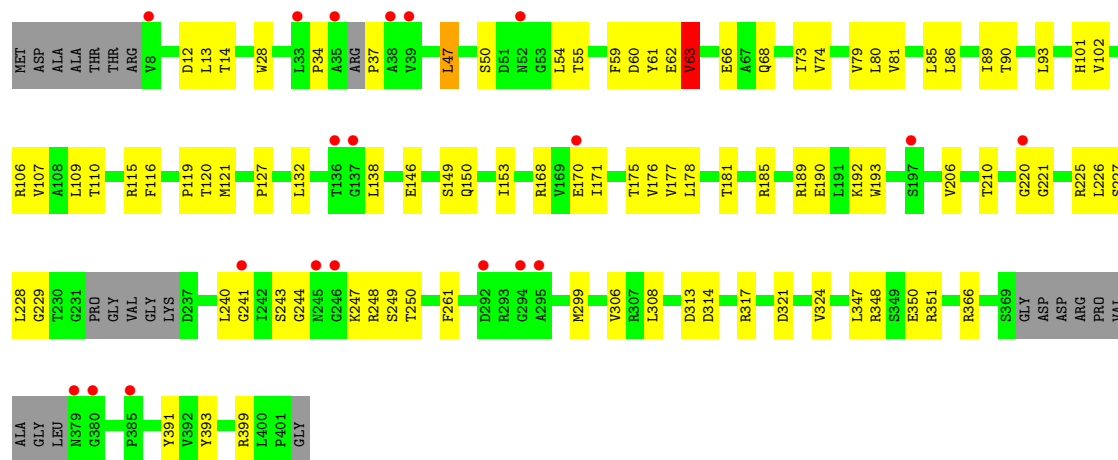


• Molecule 1: Beta sliding clamp





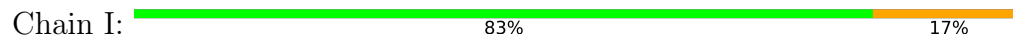
• Molecule 1: Beta sliding clamp



• Molecule 2: P7 peptide



• Molecule 2: P7 peptide

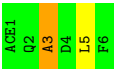


• Molecule 2: P7 peptide





● Molecule 2: P7 peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	78.49Å 81.06Å 90.02Å 68.64° 88.94° 62.56°	Depositor
Resolution (Å)	58.64 – 2.69 58.64 – 2.69	Depositor EDS
% Data completeness (in resolution range)	62.2 (58.64-2.69) 62.1 (58.64-2.69)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.69Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.210 , 0.258 0.217 , 0.262	Depositor DCC
R_{free} test set	1576 reflections (3.14%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.042	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for h,h-k,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	11359	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% 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

Bond lengths and bond angles in the following residue types are not validated in this section: CA, ACE, ALC

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	0.43	1/2822 (0.0%)	0.66	4/3861 (0.1%)
1	B	0.45	3/2795 (0.1%)	0.64	0/3814
1	C	0.43	3/2840 (0.1%)	0.71	4/3883 (0.1%)
1	D	0.42	0/2785	0.64	1/3804 (0.0%)
2	H	0.18	0/37	0.28	0/46
2	I	0.35	0/37	0.57	0/46
2	J	0.13	0/37	0.25	0/46
2	K	0.10	0/37	0.33	0/46
All	All	0.43	7/11390 (0.1%)	0.66	9/15546 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	63	VAL	C-O	-6.58	1.17	1.24
1	A	63	VAL	C-O	-5.87	1.18	1.24
1	C	59	PHE	N-CA	-5.81	1.38	1.46
1	C	65	ALA	C-O	-5.54	1.17	1.23
1	B	17	LEU	C-O	-5.40	1.17	1.23
1	B	66	GLU	C-O	-5.38	1.17	1.23
1	C	64	SER	C-N	5.29	1.41	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	228	LEU	N-CA-C	8.32	124.25	113.18
1	A	9	GLY	N-CA-C	7.01	121.52	112.13
1	C	231	GLY	CA-C-N	6.72	126.98	119.32
1	C	231	GLY	C-N-CA	6.72	126.98	119.32
1	A	372	ASP	CA-C-N	6.41	131.12	120.86
1	A	372	ASP	C-N-CA	6.41	131.12	120.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	218	GLY	N-CA-C	5.63	121.19	113.99
1	D	229	GLY	N-CA-C	5.50	118.50	110.63
1	C	399	ARG	N-CA-C	5.50	121.23	110.56

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	2775	0	2729	69	0
1	B	2750	0	2769	77	0
1	C	2794	0	2774	68	0
1	D	2740	0	2720	70	0
2	H	51	0	49	4	0
2	I	51	0	49	1	0
2	J	51	0	49	2	0
2	K	51	0	49	2	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	24	0	0	0	0
4	B	29	0	0	0	0
4	C	25	0	0	1	0
4	D	14	0	0	0	0
All	All	11359	0	11188	276	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 12.

All (276) 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:C:400:LEU:HD13	1:C:401:PRO:N	1.73	1.03
1:D:63:VAL:HG21	1:D:210:THR:CG2	2.07	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:220:GLY:H	1:B:221:GLY:HA2	1.41	0.84
1:B:239:LEU:H	1:B:239:LEU:HD12	1.44	0.80
1:D:63:VAL:CG2	1:D:210:THR:HG21	2.13	0.79
1:D:63:VAL:HG21	1:D:210:THR:HG21	1.64	0.77
1:B:275:MET:HG2	1:B:324:VAL:CG1	2.15	0.76
1:C:353:SER:HB3	1:C:366:ARG:HG2	1.69	0.75
1:D:60:ASP:HB3	1:D:62:GLU:O	1.87	0.75
1:C:269:HIS:HB3	1:C:356:PHE:O	1.87	0.74
1:B:48:THR:CG2	1:B:55:THR:OG1	2.36	0.74
1:C:115:ARG:NH2	1:D:321:ASP:OD1	2.22	0.73
1:A:28:TRP:HD1	1:A:67:ALA:HB2	1.54	0.72
1:D:66:GLU:OE1	1:D:248:ARG:NH2	2.20	0.71
1:B:275:MET:HG2	1:B:324:VAL:HG13	1.72	0.70
1:A:183:ARG:NH1	2:H:4:ASP:OD2	2.25	0.70
1:B:90:THR:HA	1:B:93:LEU:HD12	1.74	0.69
1:B:132:LEU:HD21	1:B:248:ARG:HG3	1.75	0.69
1:B:239:LEU:HD12	1:B:239:LEU:N	2.07	0.69
1:B:42:LEU:HD13	1:B:60:ASP:HA	1.75	0.68
1:B:170:GLU:OE1	1:B:262:ARG:NH2	2.26	0.68
1:D:189:ARG:HD2	1:D:190:GLU:N	2.08	0.68
1:A:80:LEU:HD23	1:A:121:MET:HB2	1.75	0.66
1:C:245:ASN:ND2	4:C:603:HOH:O	2.27	0.66
1:D:149:SER:OG	1:D:348:ARG:NH2	2.26	0.66
1:C:183:ARG:NH1	2:J:6:PHE:OXT	2.28	0.66
1:B:80:LEU:HD23	1:B:121:MET:HB2	1.78	0.66
1:C:63:VAL:CG2	1:C:210:THR:HG21	2.25	0.66
1:A:396:MET:HE2	2:H:3:ALC:HB2	1.78	0.66
1:A:14:THR:HG21	1:A:101:HIS:ND1	2.11	0.65
1:D:63:VAL:HG21	1:D:210:THR:HG23	1.78	0.65
1:A:230:THR:N	1:A:231:GLY:HA2	2.10	0.65
1:A:102:VAL:HG22	1:A:107:VAL:HG22	1.77	0.65
1:B:205:LEU:HD23	1:B:254:LEU:HB2	1.79	0.65
1:D:81:VAL:HG12	1:D:120:THR:HG22	1.80	0.64
1:B:220:GLY:N	1:B:221:GLY:HA2	2.05	0.64
1:A:118:LEU:HD23	1:B:316:GLY:HA3	1.78	0.64
1:B:60:ASP:O	1:B:61:TYR:HB2	1.97	0.64
1:C:146:GLU:O	1:C:150:GLN:HG3	1.97	0.64
1:A:205:LEU:HD22	1:A:254:LEU:HB2	1.79	0.63
1:B:84:ARG:NH2	1:B:88:ASP:OD2	2.31	0.63
1:A:33:LEU:O	1:A:34:PRO:C	2.37	0.63
1:A:14:THR:CG2	1:A:101:HIS:ND1	2.61	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:63:VAL:HG21	1:C:210:THR:HG21	1.81	0.63
1:C:228:LEU:H	1:C:229:GLY:HA3	1.64	0.63
1:B:277:VAL:HG23	1:B:344:LEU:HB3	1.81	0.62
1:B:323:VAL:O	1:B:324:VAL:HG22	1.99	0.62
1:C:132:LEU:HD12	1:C:250:THR:HG22	1.81	0.62
1:C:228:LEU:O	1:C:236:LYS:CB	2.47	0.62
1:C:63:VAL:HA	1:C:250:THR:O	2.00	0.62
1:B:275:MET:HE3	1:B:324:VAL:HG21	1.80	0.62
1:A:292:ASP:OD1	1:D:399:ARG:NH1	2.33	0.61
1:C:396:MET:HE3	2:J:3:ALC:HB2	1.83	0.61
1:B:275:MET:HE3	1:B:324:VAL:CG2	2.31	0.60
1:C:189:ARG:HB2	1:C:391:TYR:HD1	1.66	0.60
1:D:60:ASP:O	1:D:61:TYR:HB2	2.00	0.60
1:B:323:VAL:O	1:B:324:VAL:CG2	2.49	0.60
1:C:205:LEU:HD23	1:C:254:LEU:HB2	1.81	0.60
1:B:85:LEU:O	1:B:89:ILE:HG12	2.02	0.60
1:D:175:THR:HG22	1:D:192:LYS:HG2	1.83	0.60
1:D:177:VAL:HG22	1:D:190:GLU:HB2	1.84	0.59
1:A:93:LEU:HD22	1:A:111:CYS:HB3	1.83	0.59
1:B:57:SER:HB3	1:B:129:LEU:HD11	1.84	0.59
1:A:8:VAL:HG21	1:A:76:PRO:HG2	1.84	0.59
1:A:280:LEU:O	1:A:284:ILE:HG13	2.02	0.59
1:B:132:LEU:HD11	1:B:248:ARG:HD2	1.85	0.59
1:B:78:SER:OG	1:B:123:VAL:HG11	2.02	0.59
1:D:306:VAL:HB	1:D:324:VAL:HG21	1.84	0.59
1:C:149:SER:OG	1:C:348:ARG:NH2	2.36	0.59
1:C:400:LEU:CD1	1:C:401:PRO:N	2.59	0.58
1:A:51:ASP:OD2	1:A:53:GLY:N	2.33	0.58
1:B:48:THR:HG22	1:B:55:THR:OG1	2.02	0.58
1:A:62:GLU:O	1:A:251:THR:HA	2.03	0.58
1:B:323:VAL:C	1:B:324:VAL:HG23	2.28	0.58
1:D:181:THR:OG1	1:D:185:ARG:O	2.20	0.58
1:C:90:THR:HA	1:C:93:LEU:HD12	1.84	0.58
1:B:132:LEU:HD11	1:B:248:ARG:CD	2.34	0.58
1:A:38:ALA:HB1	1:A:39:VAL:HA	1.85	0.58
1:A:63:VAL:HA	1:A:250:THR:O	2.04	0.57
1:B:323:VAL:C	1:B:324:VAL:CG2	2.77	0.57
1:D:189:ARG:HD2	1:D:190:GLU:H	1.69	0.57
1:C:57:SER:HB3	1:C:129:LEU:HD11	1.86	0.57
1:D:28:TRP:CD2	1:D:247:LYS:HD2	2.40	0.56
1:B:66:GLU:OE1	1:B:248:ARG:NH2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:GLU:O	1:C:201:GLU:HG2	2.06	0.55
1:D:12:ASP:HB3	1:D:102:VAL:HB	1.89	0.55
1:A:146:GLU:O	1:A:150:GLN:HG3	2.05	0.55
1:B:146:GLU:O	1:B:150:GLN:HG3	2.05	0.55
1:A:59:PHE:CE2	1:A:127:PRO:HG2	2.42	0.55
1:A:239:LEU:HD23	1:A:252:ARG:HA	1.89	0.55
1:A:351:ARG:HB2	1:A:368:VAL:HG13	1.88	0.54
1:D:85:LEU:O	1:D:89:ILE:HG13	2.07	0.54
1:D:80:LEU:HD22	1:D:121:MET:HB2	1.90	0.54
1:A:204:VAL:HG11	1:A:228:LEU:HD22	1.89	0.54
1:B:274:THR:HA	1:B:352:VAL:O	2.07	0.54
1:D:243:SER:OG	1:D:244:GLY:N	2.41	0.54
1:D:399:ARG:HB2	2:K:3:ALC:HZ3	1.89	0.54
1:A:90:THR:HA	1:A:93:LEU:HD12	1.89	0.53
1:D:110:THR:HG23	1:D:115:ARG:HG2	1.90	0.53
1:B:160:THR:O	1:B:161:LEU:HD23	2.08	0.53
1:C:12:ASP:HB3	1:C:102:VAL:HB	1.90	0.53
1:D:170:GLU:HB2	1:D:177:VAL:HB	1.91	0.53
1:D:347:LEU:HD23	1:D:391:TYR:HE2	1.74	0.53
1:C:85:LEU:O	1:C:89:ILE:HG12	2.07	0.53
1:D:170:GLU:OE1	1:D:261:PHE:HZ	1.91	0.53
1:A:33:LEU:O	1:A:33:LEU:HG	2.07	0.53
1:D:206:VAL:HG11	1:D:240:LEU:HD23	1.91	0.53
1:A:42:LEU:O	1:A:42:LEU:HG	2.08	0.53
1:B:245:ASN:O	1:B:245:ASN:ND2	2.41	0.52
1:A:321:ASP:OD1	1:B:115:ARG:NH2	2.43	0.52
1:B:12:ASP:HB3	1:B:102:VAL:HB	1.90	0.52
1:B:274:THR:O	1:B:324:VAL:HG13	2.09	0.52
1:A:168:ARG:HD2	1:A:258:PHE:HB3	1.90	0.52
1:A:182:ASP:OD1	1:A:185:ARG:HG2	2.10	0.52
1:A:333:ILE:HD11	1:A:397:PRO:HG3	1.92	0.52
1:C:228:LEU:N	1:C:229:GLY:HA3	2.24	0.52
1:B:61:TYR:HE1	1:B:127:PRO:HG3	1.75	0.52
1:D:14:THR:OG1	1:D:74:VAL:HG22	2.09	0.52
1:D:34:PRO:HG2	1:D:37:PRO:HB3	1.92	0.52
1:A:135:GLU:OE1	1:A:225:ARG:NH1	2.44	0.51
1:D:55:THR:HG22	1:D:68:GLN:HG2	1.93	0.51
1:D:59:PHE:HD2	1:D:127:PRO:HB2	1.75	0.51
1:A:158:ASP:OD1	1:A:159:ASP:N	2.44	0.51
1:C:28:TRP:CH2	1:C:247:LYS:HB3	2.45	0.51
1:D:89:ILE:HG22	1:D:93:LEU:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:ALA:HB3	1:B:286:LEU:HD13	1.92	0.51
1:D:63:VAL:HG23	1:D:210:THR:HG21	1.91	0.51
1:A:158:ASP:OD1	1:A:160:THR:N	2.44	0.51
1:D:60:ASP:C	1:D:62:GLU:H	2.19	0.51
1:C:182:ASP:O	1:C:183:ARG:HB2	2.11	0.51
1:D:63:VAL:CG2	1:D:210:THR:CG2	2.78	0.51
1:B:17:LEU:HD21	1:B:22:PHE:CG	2.46	0.50
1:B:133:PRO:HB3	1:B:229:GLY:HA3	1.93	0.50
1:D:102:VAL:HG22	1:D:107:VAL:HG22	1.93	0.50
1:B:60:ASP:OD1	1:B:62:GLU:HB2	2.11	0.50
1:D:193:TRP:CZ3	1:D:226:LEU:HD13	2.46	0.50
1:A:12:ASP:HB3	1:A:102:VAL:HB	1.93	0.50
1:A:82:SER:OG	1:A:85:LEU:HB3	2.12	0.50
1:C:314:ASP:OD2	1:C:314:ASP:N	2.44	0.50
1:D:132:LEU:CD2	1:D:250:THR:HG22	2.42	0.50
1:C:94:PRO:HB2	1:C:96:LYS:HD2	1.94	0.49
1:A:51:ASP:OD2	1:A:52:ASN:N	2.44	0.49
1:B:93:LEU:HD22	1:B:111:CYS:HB3	1.95	0.49
1:B:260:LYS:HB2	1:B:263:GLN:OE1	2.13	0.49
1:A:8:VAL:HG11	1:A:76:PRO:HG2	1.95	0.49
1:B:55:THR:HG22	1:B:68:GLN:OE1	2.13	0.49
1:A:274:THR:HA	1:A:352:VAL:O	2.12	0.49
1:D:185:ARG:HD2	1:D:393:TYR:OH	2.12	0.49
1:A:64:SER:O	1:A:249:SER:HA	2.13	0.48
1:C:266:PRO:HD2	1:C:364:LEU:HD13	1.95	0.48
1:B:52:ASN:O	1:B:68:GLN:NE2	2.46	0.48
1:C:274:THR:HG22	1:C:351:ARG:HH21	1.78	0.48
1:C:296:GLN:HE21	1:C:398:VAL:HG11	1.78	0.48
1:C:158:ASP:OD2	1:C:160:THR:HG22	2.12	0.48
1:A:158:ASP:CG	1:A:159:ASP:H	2.21	0.48
1:C:230:THR:N	1:C:231:GLY:HA2	2.29	0.48
1:A:394:LEU:HD22	2:H:5:LEU:HD21	1.95	0.48
1:B:185:ARG:HD2	1:B:393:TYR:OH	2.13	0.48
1:C:63:VAL:HG22	1:C:210:THR:HG21	1.94	0.48
1:D:79:VAL:HG21	1:D:107:VAL:CG2	2.43	0.48
1:B:61:TYR:HE1	1:B:127:PRO:CG	2.26	0.48
1:B:34:PRO:HG2	1:B:37:PRO:HG3	1.96	0.47
1:C:26:VAL:HG11	1:C:86:LEU:HG	1.96	0.47
1:C:357:THR:OG1	1:C:358:THR:HG23	2.14	0.47
1:D:171:ILE:HD11	1:D:193:TRP:CZ2	2.49	0.47
1:B:48:THR:HG23	1:B:55:THR:OG1	2.11	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:304:GLY:HA2	1:A:324:VAL:O	2.15	0.47
1:B:143:LEU:HD21	1:B:191:LEU:HD13	1.97	0.47
1:C:351:ARG:HB2	1:C:368:VAL:HG23	1.96	0.47
1:C:399:ARG:O	1:C:400:LEU:HB3	2.14	0.47
1:C:177:VAL:HG22	1:C:190:GLU:HB2	1.96	0.47
1:C:14:THR:OG1	1:C:101:HIS:ND1	2.44	0.46
1:C:89:ILE:HD12	1:C:116:PHE:CE2	2.50	0.46
1:C:95:ASN:O	1:C:96:LYS:HG3	2.16	0.46
1:D:146:GLU:O	1:D:150:GLN:HG3	2.14	0.46
1:C:209:LYS:O	1:C:213:GLU:HG2	2.16	0.46
1:D:60:ASP:C	1:D:62:GLU:N	2.72	0.46
1:A:399:ARG:HH22	1:D:314:ASP:HB2	1.80	0.46
1:D:80:LEU:CD2	1:D:121:MET:HB2	2.45	0.46
1:C:296:GLN:HE21	1:C:398:VAL:CG1	2.29	0.46
1:C:366:ARG:CZ	1:C:387:VAL:HG21	2.46	0.46
1:A:39:VAL:HG21	1:D:61:TYR:CD1	2.51	0.46
1:D:63:VAL:HA	1:D:250:THR:O	2.14	0.46
1:D:86:LEU:O	1:D:90:THR:HG23	2.15	0.46
1:A:198:PRO:HA	1:A:199:ASP:HA	1.49	0.46
1:B:286:LEU:HD12	1:B:289:LEU:HD12	1.98	0.46
1:D:79:VAL:HG21	1:D:107:VAL:HG21	1.97	0.46
1:D:241:GLY:HA2	1:D:249:SER:O	2.15	0.46
1:B:17:LEU:HG	1:B:18:LEU:O	2.16	0.46
1:A:284:ILE:HD13	1:A:340:LEU:HG	1.97	0.45
1:B:150:GLN:NE2	1:B:346:SER:O	2.48	0.45
1:A:362:SER:HA	1:A:396:MET:HG2	1.98	0.45
1:C:400:LEU:HD13	1:C:401:PRO:C	2.41	0.45
1:A:58:GLY:HA2	1:A:126:TYR:OH	2.17	0.45
1:A:205:LEU:CD2	1:A:254:LEU:HB2	2.46	0.45
1:A:8:VAL:HG12	1:A:9:GLY:H	1.81	0.45
1:D:106:ARG:HD3	1:D:119:PRO:HA	1.97	0.45
1:C:171:ILE:HG12	1:C:176:VAL:HG22	1.98	0.45
1:B:47:LEU:HB3	1:B:54:LEU:HD11	1.99	0.45
1:B:396:MET:HE3	2:I:3:ALC:HB2	1.98	0.45
1:D:14:THR:CG2	1:D:101:HIS:ND1	2.80	0.45
1:A:232:PRO:HD3	1:A:235:GLY:O	2.17	0.45
1:C:280:LEU:O	1:C:284:ILE:HG13	2.17	0.45
1:D:299:MET:HG2	1:D:308:LEU:CD2	2.47	0.45
1:C:290:VAL:O	1:C:311:GLY:HA2	2.18	0.44
1:B:308:LEU:O	1:B:319:GLU:HA	2.18	0.44
1:D:138:LEU:HB3	1:D:225:ARG:HG3	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:SER:O	1:A:200:ILE:N	2.51	0.44
1:A:298:ARG:O	1:A:308:LEU:HA	2.18	0.44
1:B:133:PRO:HG3	1:B:239:LEU:HD13	1.99	0.44
1:D:313:ASP:OD2	1:D:317:ARG:NE	2.43	0.44
1:C:197:SER:OG	1:C:197:SER:O	2.35	0.44
1:D:12:ASP:O	1:D:14:THR:HG23	2.17	0.44
1:D:132:LEU:HD22	1:D:250:THR:HG22	2.00	0.44
1:D:220:GLY:N	1:D:221:GLY:HA2	2.32	0.44
1:D:170:GLU:OE1	1:D:261:PHE:CZ	2.70	0.44
1:B:275:MET:HG2	1:B:324:VAL:HG11	1.96	0.43
1:B:48:THR:HG23	1:B:48:THR:O	2.18	0.43
1:D:14:THR:HG22	1:D:101:HIS:ND1	2.34	0.43
1:A:119:PRO:HD2	1:B:315:VAL:O	2.19	0.43
1:B:138:LEU:HB3	1:B:225:ARG:HG2	2.01	0.43
1:B:205:LEU:CD2	1:B:254:LEU:HB2	2.47	0.43
1:C:182:ASP:OD2	1:C:185:ARG:HG2	2.18	0.43
1:C:273:ALA:HB3	1:C:354:PHE:HB2	2.01	0.43
1:D:109:LEU:HB3	1:D:116:PHE:HB2	2.01	0.43
1:A:16:ARG:HB2	1:A:99:ASP:OD1	2.19	0.43
1:A:169:VAL:O	1:A:203:ALA:HA	2.19	0.43
1:B:145:ALA:O	1:B:149:SER:OG	2.33	0.43
1:B:240:LEU:HB3	1:B:251:THR:CG2	2.49	0.43
1:C:321:ASP:OD2	1:D:115:ARG:NH2	2.51	0.43
1:D:168:ARG:O	1:D:178:LEU:HA	2.18	0.43
1:C:206:VAL:HG11	1:C:240:LEU:HD23	2.00	0.43
1:A:28:TRP:CD1	1:A:67:ALA:HB2	2.44	0.42
1:B:206:VAL:HG21	1:B:240:LEU:HD22	2.01	0.42
1:C:59:PHE:CD2	1:C:127:PRO:HG2	2.54	0.42
1:D:347:LEU:HD23	1:D:391:TYR:CE2	2.54	0.42
1:C:93:LEU:HD22	1:C:111:CYS:HB3	2.01	0.42
1:C:185:ARG:HD2	1:C:393:TYR:OH	2.20	0.42
1:C:232:PRO:N	1:C:233:GLY:HA3	2.34	0.42
1:A:14:THR:HG22	1:A:101:HIS:ND1	2.34	0.42
2:K:5:LEU:HD23	2:K:5:LEU:HA	1.76	0.42
1:C:216:LYS:O	1:C:219:ILE:HG13	2.20	0.42
1:C:400:LEU:HD22	1:C:401:PRO:N	2.35	0.42
1:B:19:ARG:O	1:B:23:ALA:N	2.40	0.42
1:C:39:VAL:HG23	1:C:40:PRO:HD2	2.02	0.42
1:C:237:ASP:HA	1:C:238:GLY:HA2	1.79	0.41
1:C:48:THR:HB	1:C:55:THR:OG1	2.20	0.41
1:C:298:ARG:O	1:C:308:LEU:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:GLN:HB3	1:B:391:TYR:CE1	2.55	0.41
1:B:50:SER:O	1:B:73:ILE:HG13	2.20	0.41
1:B:182:ASP:OD1	1:B:185:ARG:HG2	2.19	0.41
1:C:400:LEU:HD13	1:C:401:PRO:CA	2.45	0.41
1:D:47:LEU:HB3	1:D:54:LEU:HD11	2.02	0.41
1:A:8:VAL:CG1	1:A:9:GLY:H	2.33	0.41
1:A:105:ASN:OD1	1:A:106:ARG:HG3	2.21	0.41
1:A:264:LEU:HD13	2:H:5:LEU:HD23	2.03	0.41
1:B:141:ALA:HB3	1:B:221:GLY:HA3	2.03	0.41
1:D:50:SER:O	1:D:73:ILE:HG13	2.19	0.41
1:A:229:GLY:HA3	1:A:235:GLY:O	2.21	0.41
1:A:237:ASP:HA	1:A:238:GLY:HA2	1.86	0.41
1:B:132:LEU:HD11	1:B:248:ARG:NE	2.36	0.41
1:C:399:ARG:O	1:C:400:LEU:CB	2.69	0.41
1:D:13:LEU:O	1:D:102:VAL:HG23	2.21	0.41
1:D:150:GLN:O	1:D:153:ILE:HG22	2.21	0.41
1:D:350:GLU:HG2	1:D:351:ARG:HG3	2.02	0.41
1:B:18:LEU:HD12	1:B:18:LEU:HA	1.82	0.41
1:C:175:THR:HG22	1:C:192:LYS:HG2	2.02	0.41
1:A:33:LEU:O	1:A:33:LEU:CG	2.69	0.40
1:A:277:VAL:HG12	1:A:352:VAL:HB	2.01	0.40
1:B:107:VAL:HG23	1:B:120:THR:CG2	2.51	0.40
1:B:89:ILE:HD12	1:B:116:PHE:CE2	2.57	0.40
1:B:399:ARG:O	1:B:400:LEU:HB2	2.20	0.40
1:C:240:LEU:HB2	1:C:253:LEU:HD21	2.03	0.40
1:A:193:TRP:CZ3	1:A:226:LEU:HD12	2.57	0.40
1:B:298:ARG:HG3	1:B:334:ALA:HB2	2.04	0.40
1:A:13:LEU:O	1:A:102:VAL:HG23	2.21	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	383/402 (95%)	364 (95%)	19 (5%)	0	100	100
1	B	370/402 (92%)	358 (97%)	12 (3%)	0	100	100
1	C	383/402 (95%)	358 (94%)	22 (6%)	3 (1%)	16	34
1	D	371/402 (92%)	356 (96%)	14 (4%)	1 (0%)	36	57
2	H	3/6 (50%)	3 (100%)	0	0	100	100
2	I	3/6 (50%)	3 (100%)	0	0	100	100
2	J	3/6 (50%)	3 (100%)	0	0	100	100
2	K	3/6 (50%)	3 (100%)	0	0	100	100
All	All	1519/1632 (93%)	1448 (95%)	67 (4%)	4 (0%)	36	57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	63	VAL
1	C	229	GLY
1	D	63	VAL
1	C	400	LEU

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	280/316 (89%)	278 (99%)	2 (1%)	76	89
1	B	285/316 (90%)	274 (96%)	11 (4%)	28	54
1	C	286/316 (90%)	282 (99%)	4 (1%)	59	80
1	D	281/316 (89%)	275 (98%)	6 (2%)	47	73
2	H	4/4 (100%)	4 (100%)	0	100	100
2	I	4/4 (100%)	4 (100%)	0	100	100
2	J	4/4 (100%)	4 (100%)	0	100	100
2	K	4/4 (100%)	4 (100%)	0	100	100
All	All	1148/1280 (90%)	1125 (98%)	23 (2%)	48	74

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	186	LEU
1	B	17	LEU
1	B	18	LEU
1	B	55	THR
1	B	63	VAL
1	B	64	SER
1	B	117	SER
1	B	124	GLU
1	B	157	ARG
1	B	160	THR
1	B	239	LEU
1	B	305	SER
1	C	47	LEU
1	C	314	ASP
1	C	357	THR
1	C	358	THR
1	D	47	LEU
1	D	63	VAL
1	D	176	VAL
1	D	227	SER
1	D	228	LEU
1	D	366	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	379	ASN
1	C	296	GLN
1	D	32	ASN
1	D	68	GLN
1	D	269	HIS
2	H	2	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ALC	I	3	2	9,11,12	0.21	0	11,13,15	0.79	1 (9%)
2	ALC	H	3	2	9,11,12	0.20	0	11,13,15	0.70	0
2	ALC	J	3	2	9,11,12	0.13	0	11,13,15	0.91	1 (9%)
2	ALC	K	3	2	9,11,12	0.20	0	11,13,15	0.92	1 (9%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ALC	I	3	2	-	3/5/14/16	0/1/1/1
2	ALC	H	3	2	-	4/5/14/16	0/1/1/1
2	ALC	J	3	2	-	4/5/14/16	0/1/1/1
2	ALC	K	3	2	-	3/5/14/16	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	3	ALC	CB-CA-C	-2.94	106.47	110.99
2	K	3	ALC	CB-CA-C	-2.43	107.25	110.99
2	I	3	ALC	CB-CA-C	-2.14	107.69	110.99

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	H	3	ALC	N-CA-CB-CG
2	H	3	ALC	CA-CB-CG-CD2
2	I	3	ALC	CA-CB-CG-CD2
2	J	3	ALC	N-CA-CB-CG
2	J	3	ALC	CA-CB-CG-CD2
2	K	3	ALC	CA-CB-CG-CD2
2	K	3	ALC	CA-CB-CG-CD1
2	I	3	ALC	CA-CB-CG-CD1
2	J	3	ALC	CA-CB-CG-CD1
2	I	3	ALC	N-CA-CB-CG
2	K	3	ALC	N-CA-CB-CG
2	H	3	ALC	C-CA-CB-CG
2	J	3	ALC	C-CA-CB-CG
2	H	3	ALC	CA-CB-CG-CD1

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	3	ALC	1	0
2	H	3	ALC	1	0
2	J	3	ALC	1	0
2	K	3	ALC	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	389/402 (96%)	0.51	24 (6%) 26 22	16, 39, 73, 98	0
1	B	376/402 (93%)	0.22	20 (5%) 32 27	15, 31, 68, 96	0
1	C	389/402 (96%)	0.30	24 (6%) 26 22	14, 31, 75, 102	0
1	D	379/402 (94%)	0.45	20 (5%) 32 27	15, 38, 71, 94	0
2	H	4/6 (66%)	0.14	0 100 100	26, 36, 38, 45	0
2	I	4/6 (66%)	-0.00	0 100 100	32, 39, 41, 47	0
2	J	4/6 (66%)	-0.04	0 100 100	22, 31, 34, 34	0
2	K	4/6 (66%)	0.08	0 100 100	32, 36, 43, 45	0
All	All	1549/1632 (94%)	0.37	88 (5%) 29 25	14, 35, 72, 102	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	197	SER	5.2
1	B	64	SER	4.8
1	D	294	GLY	4.8
1	C	312	ALA	4.6
1	D	38	ALA	4.6
1	B	63	VAL	4.6
1	C	374	PRO	4.5
1	C	52	ASN	4.5
1	B	60	ASP	4.1
1	C	221	GLY	4.1
1	A	198	PRO	3.9
1	B	65	ALA	3.8
1	A	35	ALA	3.8
1	A	372	ASP	3.7
1	A	42	LEU	3.6
1	A	235	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	269	HIS	3.5
1	C	36	ARG	3.5
1	C	220	GLY	3.5
1	A	232	PRO	3.5
1	D	379	ASN	3.5
1	A	62	GLU	3.2
1	D	33	LEU	3.2
1	C	401	PRO	3.2
1	B	313	ASP	3.2
1	C	34	PRO	3.1
1	A	52	ASN	3.1
1	C	198	PRO	3.0
1	C	235	GLY	3.0
1	C	35	ALA	3.0
1	C	232	PRO	2.9
1	C	37	PRO	2.9
1	B	62	GLU	2.9
1	A	34	PRO	2.8
1	D	241	GLY	2.8
1	D	220	GLY	2.8
1	D	8	VAL	2.8
1	D	170	GLU	2.8
1	C	381	ASN	2.8
1	D	380	GLY	2.8
1	B	315	VAL	2.7
1	D	385	PRO	2.7
1	C	32	ASN	2.7
1	B	257	GLU	2.6
1	A	46	LEU	2.6
1	C	373	ARG	2.6
1	A	38	ALA	2.6
1	B	52	ASN	2.6
1	D	245	ASN	2.6
1	D	137	GLY	2.5
1	D	197	SER	2.5
1	C	222	SER	2.5
1	B	324	VAL	2.4
1	B	256	ALA	2.4
1	B	314	ASP	2.4
1	B	379	ASN	2.3
1	D	39	VAL	2.3
1	D	295	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	160	THR	2.3
1	A	233	GLY	2.3
1	C	380	GLY	2.3
1	A	41	VAL	2.3
1	B	380	GLY	2.2
1	A	256	ALA	2.2
1	B	312	ALA	2.2
1	C	38	ALA	2.2
1	A	374	PRO	2.2
1	C	371	ASP	2.2
1	B	48	THR	2.2
1	D	52	ASN	2.2
1	D	136	THR	2.1
1	B	198	PRO	2.1
1	A	143	LEU	2.1
1	A	255	ASP	2.1
1	B	381	ASN	2.1
1	D	246	GLY	2.1
1	D	292	ASP	2.1
1	A	37	PRO	2.1
1	A	218	GLY	2.1
1	C	229	GLY	2.1
1	A	199	ASP	2.0
1	B	248	ARG	2.0
1	C	50	SER	2.0
1	A	220	GLY	2.0
1	A	312	ALA	2.0
1	D	35	ALA	2.0
1	A	267	THR	2.0
1	C	400	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ALC	K	3	11/12	0.90	0.10	23,32,39,41	0
2	ALC	I	3	11/12	0.92	0.10	18,30,40,41	0
2	ALC	H	3	11/12	0.94	0.09	12,25,33,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ALC	J	3	11/12	0.95	0.09	20,31,34,38	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CA	C	501	1/1	0.87	0.08	52,52,52,52	0
3	CA	B	501	1/1	0.93	0.06	55,55,55,55	0
3	CA	D	501	1/1	0.95	0.07	26,26,26,26	0
3	CA	B	502	1/1	0.96	0.04	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.