



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 11:13 AM EDT

PDB ID : 2VSQ / pdb_00002vsq
Title : Structure of surfactin A synthetase C (SrfA-C), a nonribosomal peptide synthetase termination module
Authors : Tanovic, A.; Samel, S.A.; Essen, L.-O.; Marahiel, M.A.
Deposited on : 2008-04-29
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
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

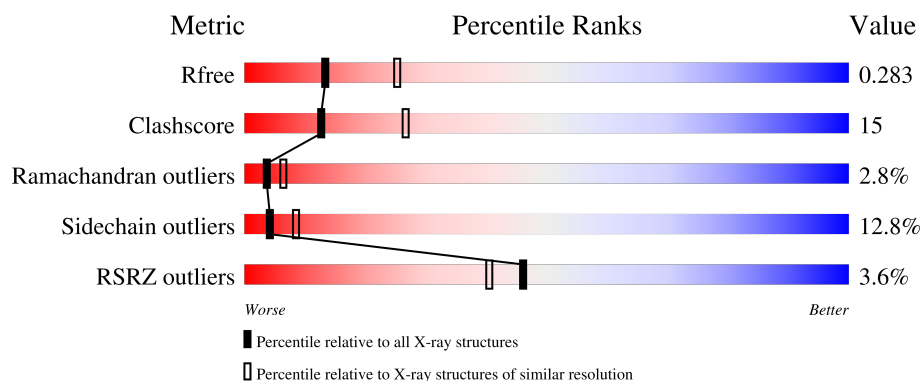
1 Overall quality at a glance

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	1304	 4% 65% 25% 6% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10109 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.

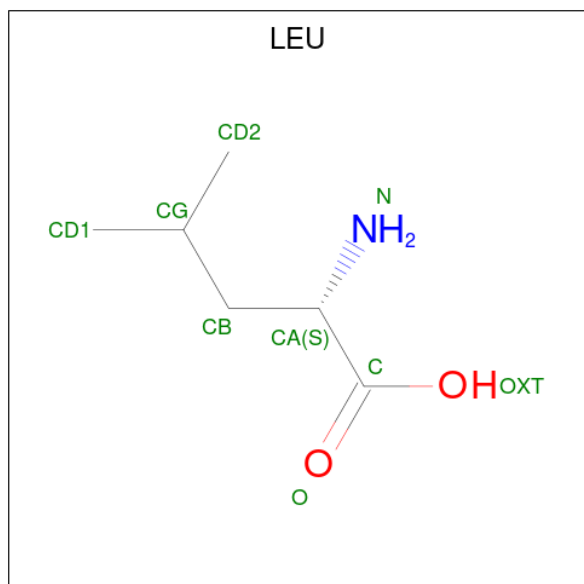
- Molecule 1 is a protein called SURFACTIN SYNTHETASE SUBUNIT 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1273	10056	6409	1685	1924	38	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	ALA	PRO	conflict	UNP Q08787
A	33	SER	THR	conflict	UNP Q08787
A	235	PRO	LEU	conflict	UNP Q08787
A	1003	ALA	SER	engineered mutation	UNP Q08787
A	1223	ILE	MET	conflict	UNP Q08787
A	1240	MET	ASN	conflict	UNP Q08787

- Molecule 2 is LEUCINE (CCD ID: LEU) (formula: $C_6H_{13}NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			9	6	1	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		

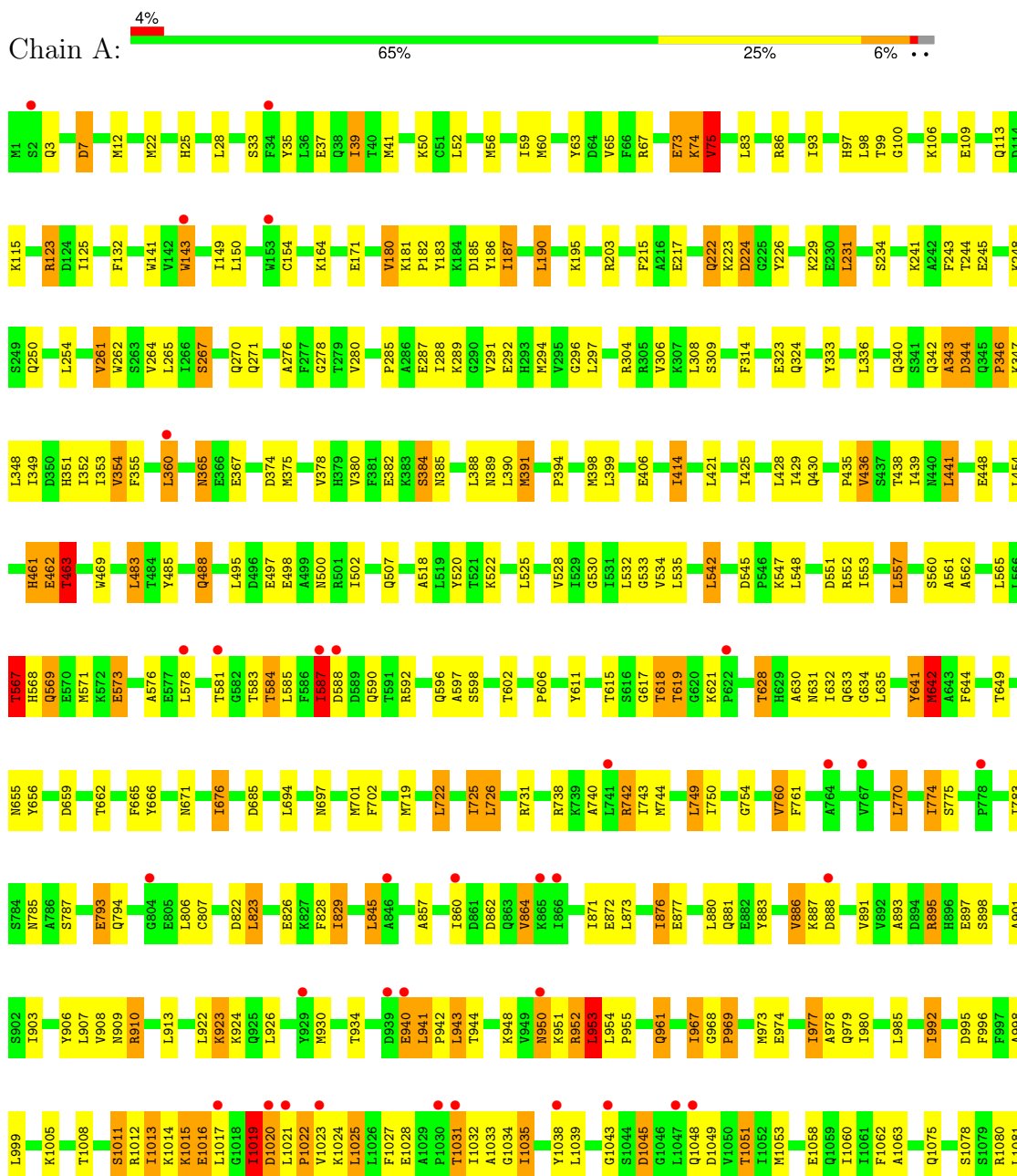
- Molecule 4 is water.

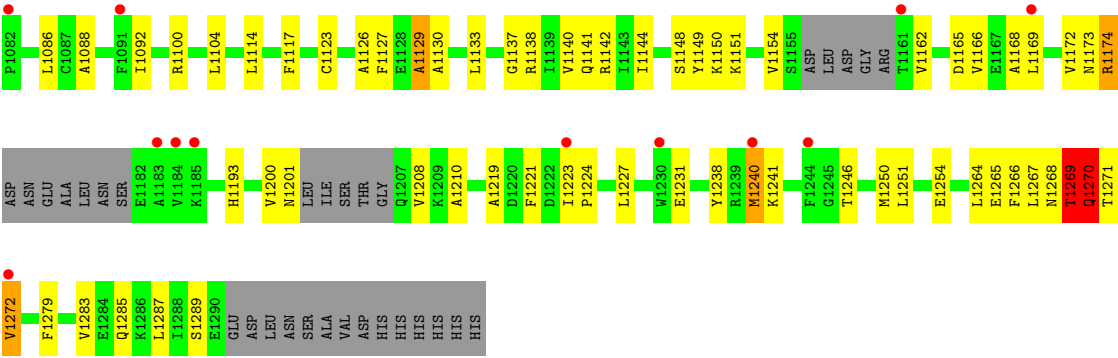
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	39	Total	O	0	0
			39	39		

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: SURFACTIN SYNTHETASE SUBUNIT 3





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.11Å 106.56Å 92.40Å 90.00° 97.35° 90.00°	Depositor
Resolution (Å)	82.50 – 2.60 82.50 – 2.60	Depositor EDS
% Data completeness (in resolution range)	93.1 (82.50-2.60) 92.5 (82.50-2.60)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.213 , 0.272 0.219 , 0.283	Depositor DCC
R_{free} test set	1040 reflections (2.17%)	wwPDB-VP
Wilson B-factor (Å ²)	69.8	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 79.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10109	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% 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: SO4

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	0.68	2/10270 (0.0%)	0.86	8/13922 (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	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	596	GLN	CD-NE2	15.26	1.65	1.33
1	A	596	GLN	CD-OE1	10.26	1.43	1.23

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	596	GLN	CG-CD-OE1	-13.08	94.65	120.80
1	A	596	GLN	CG-CD-NE2	10.89	132.73	116.40
1	A	596	GLN	OE1-CD-NE2	8.75	131.35	122.60
1	A	1240	MET	N-CA-C	6.04	118.31	108.76
1	A	462	GLU	N-CA-C	5.69	117.39	107.49
1	A	641	TYR	N-CA-C	-5.46	101.82	109.96
1	A	1027	PHE	N-CA-C	-5.41	105.30	111.14
1	A	143	TRP	CB-CA-C	-5.30	100.79	109.53

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	462	GLU	Peptide
1	A	567	THR	Peptide
1	A	587	ILE	Peptide

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	10056	0	9872	303	0
2	A	9	0	10	0	0
3	A	5	0	0	0	0
4	A	39	0	0	0	0
All	All	10109	0	9882	303	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 15.

All (303) 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:876:ILE:HD11	1:A:891:VAL:CG2	1.60	1.31
1:A:876:ILE:CD1	1:A:891:VAL:HG21	1.61	1.29
1:A:587:ILE:HD13	1:A:587:ILE:O	1.53	1.09
1:A:1021:LEU:HD22	1:A:1022:PRO:HD2	1.43	1.00
1:A:60:MET:HE1	1:A:125:ILE:HG21	1.45	0.95
1:A:39:ILE:HD11	1:A:378:VAL:HG13	1.49	0.94
1:A:744:MET:HE1	1:A:749:LEU:HD11	1.52	0.92
1:A:60:MET:HE1	1:A:125:ILE:CG2	2.01	0.91
1:A:483:LEU:HD13	1:A:525:LEU:HD12	1.51	0.91
1:A:41:MET:HE2	1:A:375:MET:SD	2.11	0.90
1:A:463:THR:HG23	1:A:606:PRO:HG2	1.52	0.90
1:A:642:MET:N	1:A:642:MET:HE3	1.88	0.88
1:A:725:ILE:CG2	1:A:749:LEU:HD13	2.04	0.87
1:A:943:LEU:O	1:A:944:THR:HG22	1.76	0.85
1:A:644:PHE:HB2	1:A:671:ASN:HD22	1.40	0.84
1:A:1238:TYR:CE2	1:A:1240:MET:HE1	2.15	0.80
1:A:783:ILE:HD12	1:A:785:ASN:H	1.46	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:MET:HE2	1:A:125:ILE:HD13	1.64	0.79
1:A:611:TYR:CZ	1:A:662:THR:HG21	2.17	0.79
1:A:967:ILE:O	1:A:992:ILE:HD13	1.86	0.76
1:A:883:TYR:CE1	1:A:913:LEU:HD13	2.21	0.76
1:A:618:THR:HG23	1:A:621:LYS:O	1.87	0.75
1:A:241:LYS:O	1:A:244:THR:HG22	1.87	0.74
1:A:264:VAL:HG22	1:A:436:VAL:HG11	1.68	0.74
1:A:611:TYR:OH	1:A:662:THR:HG21	1.88	0.73
1:A:1092:ILE:HG21	1:A:1100:ARG:HD2	1.70	0.73
1:A:973:MET:O	1:A:977:ILE:HG23	1.89	0.73
1:A:725:ILE:CG2	1:A:749:LEU:CD1	2.67	0.72
1:A:522:LYS:HB2	1:A:571:MET:HE3	1.69	0.72
1:A:628:THR:HG22	1:A:631:ASN:H	1.55	0.72
1:A:740:ALA:HB1	1:A:744:MET:CE	2.20	0.72
1:A:952:ARG:HD3	1:A:953:LEU:HD23	1.72	0.71
1:A:642:MET:SD	1:A:642:MET:C	2.74	0.71
1:A:1166:VAL:HG21	1:A:1193:HIS:HA	1.72	0.71
1:A:39:ILE:CD1	1:A:378:VAL:HG13	2.20	0.71
1:A:463:THR:CG2	1:A:606:PRO:HG2	2.21	0.71
1:A:1019:ILE:HD13	1:A:1019:ILE:O	1.91	0.70
1:A:641:TYR:CD2	1:A:642:MET:HE1	2.25	0.70
1:A:365:ASN:N	1:A:365:ASN:HD22	1.89	0.70
1:A:74:LYS:O	1:A:75:VAL:HG13	1.92	0.69
1:A:583:THR:CG2	1:A:585:LEU:HD11	2.22	0.69
1:A:278:GLY:HA3	1:A:352:ILE:HG22	1.72	0.69
1:A:642:MET:HE2	1:A:750:ILE:HD11	1.74	0.69
1:A:1264:LEU:O	1:A:1268:ASN:N	2.25	0.69
1:A:888:ASP:HB2	1:A:908:VAL:HG22	1.74	0.68
1:A:876:ILE:HG22	1:A:930:MET:HE2	1.74	0.68
1:A:587:ILE:O	1:A:587:ILE:CD1	2.39	0.68
1:A:1031:THR:HG22	1:A:1034:GLY:H	1.58	0.68
1:A:754:GLY:HA3	1:A:761:PHE:HA	1.76	0.68
1:A:1268:ASN:O	1:A:1270:GLN:N	2.27	0.68
1:A:641:TYR:O	1:A:642:MET:O	2.11	0.67
1:A:876:ILE:C	1:A:876:ILE:HD12	2.20	0.67
1:A:528:VAL:HG13	1:A:665:PHE:HE1	1.60	0.67
1:A:520:TYR:O	1:A:567:THR:HG23	1.95	0.66
1:A:642:MET:HE2	1:A:750:ILE:CD1	2.26	0.66
1:A:725:ILE:HG23	1:A:749:LEU:CD1	2.25	0.66
1:A:744:MET:HE1	1:A:749:LEU:CD1	2.25	0.66
1:A:573:GLU:HG2	1:A:576:ALA:HB2	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1039:LEU:HD23	1:A:1043:GLY:HA2	1.77	0.66
1:A:1013:ILE:C	1:A:1013:ILE:HD12	2.21	0.65
1:A:977:ILE:HD11	1:A:1032:ILE:HG23	1.76	0.65
1:A:725:ILE:HG23	1:A:749:LEU:HD13	1.78	0.65
1:A:463:THR:HG23	1:A:606:PRO:CG	2.26	0.65
1:A:243:PHE:CE2	1:A:398:MET:HE1	2.33	0.64
1:A:864:VAL:HG11	1:A:893:ALA:CB	2.26	0.64
1:A:738:ARG:NH2	1:A:770:LEU:HD12	2.13	0.64
1:A:99:THR:HG22	1:A:100:GLY:N	2.11	0.63
1:A:583:THR:HG22	1:A:585:LEU:CD1	2.28	0.63
1:A:1013:ILE:O	1:A:1019:ILE:N	2.32	0.63
1:A:1023:VAL:HG23	1:A:1024:LYS:H	1.64	0.63
1:A:1149:TYR:HB2	1:A:1227:LEU:HD22	1.80	0.63
1:A:774:ILE:O	1:A:774:ILE:HG22	1.99	0.62
1:A:934:THR:HG22	1:A:961:GLN:HB3	1.81	0.62
1:A:215:PHE:H	1:A:270:GLN:HE22	1.45	0.62
1:A:864:VAL:HG23	1:A:871:ILE:HB	1.80	0.62
1:A:1039:LEU:HD23	1:A:1043:GLY:CA	2.30	0.62
1:A:35:TYR:OH	1:A:1024:LYS:NZ	2.34	0.61
1:A:1172:VAL:C	1:A:1173:ASN:HD22	2.08	0.61
1:A:41:MET:CE	1:A:375:MET:SD	2.88	0.61
1:A:340:GLN:NE2	1:A:346:PRO:O	2.34	0.61
1:A:1014:LYS:O	1:A:1015:LYS:CB	2.49	0.61
1:A:1219:ALA:HB2	1:A:1246:THR:HG22	1.83	0.61
1:A:262:TRP:HE1	1:A:351:HIS:HE2	1.47	0.61
1:A:895:ARG:HD2	1:A:901:ALA:HB2	1.82	0.61
1:A:992:ILE:O	1:A:1031:THR:HG23	2.00	0.61
1:A:1053:MET:HE2	1:A:1075:GLN:HA	1.83	0.60
1:A:995:ASP:HB3	1:A:998:ALA:HB3	1.84	0.60
1:A:502:ILE:HD11	1:A:587:ILE:CG1	2.31	0.60
1:A:1060:ILE:HG21	1:A:1062:PHE:CZ	2.36	0.60
1:A:880:LEU:O	1:A:886:VAL:HG11	2.02	0.59
1:A:63:TYR:CE2	1:A:180:VAL:HG21	2.38	0.59
1:A:883:TYR:HE1	1:A:913:LEU:HD13	1.68	0.59
1:A:944:THR:HG23	1:A:944:THR:O	2.01	0.59
1:A:93:ILE:HD13	1:A:132:PHE:CE2	2.38	0.58
1:A:530:GLY:O	1:A:534:VAL:HG23	2.04	0.58
1:A:1092:ILE:HD13	1:A:1100:ARG:CD	2.33	0.58
1:A:1051:THR:HB	1:A:1088:ALA:HB3	1.83	0.58
1:A:738:ARG:NH2	1:A:770:LEU:CD1	2.68	0.57
1:A:719:MET:HE3	1:A:743:ILE:HG13	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:495:LEU:HD11	1:A:532:LEU:HD12	1.86	0.57
1:A:485:TYR:HB3	1:A:525:LEU:HD22	1.86	0.57
1:A:502:ILE:HG22	1:A:533:GLY:HA3	1.87	0.57
1:A:740:ALA:HB1	1:A:744:MET:HE2	1.87	0.57
1:A:583:THR:HG22	1:A:585:LEU:HD11	1.84	0.56
1:A:1032:ILE:HA	1:A:1035:ILE:HG23	1.87	0.56
1:A:719:MET:HA	1:A:722:LEU:HD22	1.88	0.56
1:A:485:TYR:HB3	1:A:525:LEU:CD2	2.34	0.56
1:A:567:THR:HG22	1:A:568:HIS:N	2.21	0.56
1:A:573:GLU:O	1:A:576:ALA:HB2	2.06	0.56
1:A:7:ASP:C	1:A:7:ASP:OD2	2.49	0.56
1:A:60:MET:CE	1:A:125:ILE:HD13	2.33	0.55
1:A:522:LYS:CB	1:A:571:MET:HE3	2.36	0.55
1:A:60:MET:CE	1:A:125:ILE:HG21	2.29	0.55
1:A:1021:LEU:HD13	1:A:1022:PRO:N	2.22	0.55
1:A:63:TYR:HE2	1:A:180:VAL:HG21	1.70	0.55
1:A:1014:LYS:HA	1:A:1020:ASP:HA	1.87	0.55
1:A:65:VAL:HB	1:A:149:ILE:HB	1.89	0.55
1:A:576:ALA:HB3	1:A:578:LEU:HD12	1.89	0.55
1:A:1231:GLU:HA	1:A:1238:TYR:CD2	2.42	0.54
1:A:528:VAL:HG13	1:A:665:PHE:CE1	2.41	0.54
1:A:254:LEU:H	1:A:254:LEU:HD12	1.73	0.54
1:A:343:ALA:O	1:A:344:ASP:HB2	2.08	0.54
1:A:502:ILE:HD11	1:A:587:ILE:HG12	1.90	0.54
1:A:1224:PRO:HG2	1:A:1227:LEU:HD12	1.91	0.53
1:A:439:ILE:HG22	1:A:441:LEU:HD13	1.91	0.53
1:A:942:PRO:O	1:A:950:ASN:HB3	2.09	0.53
1:A:941:LEU:HD12	1:A:941:LEU:N	2.24	0.53
1:A:731:ARG:HH12	1:A:872:GLU:CD	2.17	0.53
1:A:845:LEU:HD13	1:A:860:ILE:HG13	1.91	0.53
1:A:576:ALA:HB3	1:A:578:LEU:CD1	2.39	0.52
1:A:744:MET:CE	1:A:749:LEU:HD11	2.35	0.52
1:A:701:MET:O	1:A:725:ILE:HD12	2.10	0.52
1:A:262:TRP:CD1	1:A:353:ILE:HD11	2.44	0.52
1:A:676:ILE:HD12	1:A:676:ILE:N	2.25	0.52
1:A:502:ILE:HD11	1:A:587:ILE:HG13	1.91	0.52
1:A:725:ILE:HG21	1:A:749:LEU:CD1	2.38	0.51
1:A:56:MET:HE1	1:A:143:TRP:CH2	2.46	0.51
1:A:1092:ILE:HD13	1:A:1100:ARG:HD2	1.91	0.51
1:A:1172:VAL:HG23	1:A:1173:ASN:ND2	2.25	0.51
1:A:99:THR:CG2	1:A:100:GLY:N	2.74	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:LEU:HD12	1:A:414:ILE:HG22	1.93	0.50
1:A:542:LEU:HD23	1:A:542:LEU:C	2.36	0.50
1:A:1053:MET:HE2	1:A:1075:GLN:CA	2.41	0.50
1:A:871:ILE:HG23	1:A:930:MET:HE3	1.92	0.50
1:A:60:MET:HE1	1:A:125:ILE:HG23	1.91	0.50
1:A:725:ILE:HG21	1:A:749:LEU:HD13	1.91	0.50
1:A:906:TYR:C	1:A:907:LEU:HD12	2.37	0.50
1:A:73:GLU:O	1:A:74:LYS:HD3	2.12	0.50
1:A:439:ILE:HG22	1:A:441:LEU:CD1	2.41	0.50
1:A:568:HIS:O	1:A:569:GLN:CB	2.60	0.50
1:A:659:ASP:O	1:A:760:VAL:HG22	2.12	0.50
1:A:618:THR:O	1:A:619:THR:C	2.55	0.50
1:A:738:ARG:CZ	1:A:770:LEU:HD12	2.42	0.50
1:A:182:PRO:O	1:A:294:MET:HE3	2.12	0.49
1:A:187:ILE:N	1:A:187:ILE:HD12	2.27	0.49
1:A:224:ASP:OD1	1:A:224:ASP:N	2.38	0.49
1:A:806:LEU:C	1:A:806:LEU:HD23	2.37	0.49
1:A:944:THR:O	1:A:944:THR:CG2	2.60	0.49
1:A:909:ASN:O	1:A:910:ARG:CB	2.60	0.49
1:A:1015:LYS:O	1:A:1017:LEU:N	2.46	0.49
1:A:12:MET:SD	1:A:12:MET:N	2.86	0.49
1:A:150:LEU:HD22	1:A:154:CYS:CB	2.42	0.49
1:A:285:PRO:HD2	1:A:291:VAL:HG21	1.95	0.49
1:A:845:LEU:HB3	1:A:857:ALA:HB3	1.94	0.49
1:A:264:VAL:HG22	1:A:436:VAL:CG1	2.39	0.49
1:A:314:PHE:CD1	1:A:428:LEU:HD22	2.47	0.49
1:A:702:PHE:C	1:A:702:PHE:CD2	2.90	0.49
1:A:1023:VAL:HG23	1:A:1024:LYS:N	2.26	0.49
1:A:1114:LEU:O	1:A:1141:GLN:N	2.45	0.49
1:A:583:THR:HG22	1:A:585:LEU:HD12	1.95	0.48
1:A:845:LEU:HD13	1:A:860:ILE:CG1	2.43	0.48
1:A:1162:VAL:HG12	1:A:1193:HIS:CE1	2.48	0.48
1:A:186:TYR:CE2	1:A:190:LEU:HD22	2.49	0.48
1:A:97:HIS:CE1	1:A:98:LEU:HD21	2.49	0.48
1:A:873:LEU:HD23	1:A:877:GLU:OE1	2.14	0.48
1:A:1005:LYS:HA	1:A:1008:THR:HB	1.94	0.48
1:A:1126:ALA:O	1:A:1129:ALA:HB3	2.13	0.48
1:A:1142:ARG:NH2	1:A:1267:LEU:HD23	2.29	0.48
1:A:542:LEU:C	1:A:542:LEU:CD2	2.86	0.48
1:A:641:TYR:CE2	1:A:642:MET:HE1	2.49	0.48
1:A:1172:VAL:HG23	1:A:1173:ASN:HD22	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:ILE:HB	1:A:291:VAL:HB	1.95	0.47
1:A:560:SER:O	1:A:561:ALA:HB3	2.15	0.47
1:A:922:LEU:HD22	1:A:926:LEU:HD12	1.96	0.47
1:A:553:ILE:HG22	1:A:557:LEU:HD22	1.95	0.47
1:A:1151:LYS:NZ	1:A:1201:ASN:O	2.48	0.47
1:A:1021:LEU:HD22	1:A:1022:PRO:CD	2.29	0.47
1:A:1063:ALA:HA	1:A:1117:PHE:HB3	1.97	0.47
1:A:1266:PHE:O	1:A:1269:THR:HG23	2.15	0.47
1:A:150:LEU:HD22	1:A:154:CYS:HB3	1.97	0.47
1:A:346:PRO:O	1:A:348:LEU:N	2.48	0.47
1:A:545:ASP:HB3	1:A:548:LEU:HD13	1.95	0.47
1:A:873:LEU:HD21	1:A:891:VAL:O	2.14	0.47
1:A:881:GLN:O	1:A:886:VAL:HG21	2.14	0.47
1:A:565:LEU:O	1:A:584:THR:HA	2.14	0.47
1:A:606:PRO:O	1:A:628:THR:HG23	2.15	0.47
1:A:1264:LEU:O	1:A:1265:GLU:C	2.57	0.47
1:A:261:VAL:HG21	1:A:425:ILE:CD1	2.45	0.46
1:A:488:GLN:N	1:A:488:GLN:OE1	2.48	0.46
1:A:1123:CYS:SG	1:A:1148:SER:HB3	2.55	0.46
1:A:995:ASP:HB3	1:A:998:ALA:CB	2.45	0.46
1:A:1078:SER:HA	1:A:1086:LEU:HD11	1.97	0.46
1:A:261:VAL:HG21	1:A:425:ILE:HD13	1.97	0.46
1:A:469:TRP:HE1	1:A:633:GLN:HE21	1.63	0.46
1:A:793:GLU:CD	1:A:793:GLU:H	2.24	0.46
1:A:360:LEU:N	1:A:360:LEU:HD22	2.30	0.45
1:A:1133:LEU:O	1:A:1137:GLY:N	2.49	0.45
1:A:1031:THR:CG2	1:A:1033:ALA:HB3	2.46	0.45
1:A:573:GLU:HG2	1:A:573:GLU:O	2.17	0.45
1:A:628:THR:O	1:A:631:ASN:HB2	2.16	0.45
1:A:1271:THR:O	1:A:1272:VAL:C	2.59	0.45
1:A:876:ILE:CG2	1:A:930:MET:HE2	2.42	0.45
1:A:744:MET:HE3	1:A:749:LEU:HD21	1.98	0.45
1:A:234:SER:CB	1:A:398:MET:O	2.65	0.45
1:A:495:LEU:HD11	1:A:532:LEU:CD1	2.46	0.45
1:A:1169:LEU:O	1:A:1172:VAL:HG22	2.17	0.45
1:A:641:TYR:O	1:A:641:TYR:CG	2.69	0.45
1:A:1130:ALA:HA	1:A:1133:LEU:HD12	1.98	0.45
1:A:649:THR:HG23	1:A:676:ILE:HD11	1.98	0.45
1:A:725:ILE:HG23	1:A:749:LEU:HD12	1.96	0.45
1:A:1221:PHE:HE2	1:A:1223:ILE:HD12	1.82	0.45
1:A:52:LEU:HD21	1:A:141:TRP:CH2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:MET:CE	1:A:125:ILE:CG2	2.86	0.44
1:A:183:TYR:HA	1:A:294:MET:CE	2.47	0.44
1:A:365:ASN:N	1:A:365:ASN:ND2	2.59	0.44
1:A:822:ASP:CG	1:A:823:LEU:HD13	2.41	0.44
1:A:942:PRO:O	1:A:950:ASN:CB	2.66	0.44
1:A:39:ILE:CD1	1:A:378:VAL:CG1	2.92	0.44
1:A:968:GLY:O	1:A:969:PRO:C	2.60	0.44
1:A:435:PRO:O	1:A:438:THR:HB	2.17	0.44
1:A:985:LEU:HD21	1:A:996:PHE:CZ	2.52	0.44
1:A:1127:PHE:CE1	1:A:1208:VAL:HG13	2.53	0.44
1:A:226:TYR:CE2	1:A:384:SER:HB2	2.52	0.44
1:A:1133:LEU:HD13	1:A:1140:VAL:HG22	1.99	0.44
1:A:806:LEU:HD23	1:A:807:CYS:N	2.32	0.44
1:A:141:TRP:CE2	1:A:143:TRP:NE1	2.85	0.43
1:A:351:HIS:HB2	1:A:388:LEU:O	2.19	0.43
1:A:634:GLY:HA2	1:A:785:ASN:HD22	1.82	0.43
1:A:1045:ASP:HA	1:A:1049:ASP:HB3	2.00	0.43
1:A:1114:LEU:HD11	1:A:1138:ARG:CZ	2.49	0.43
1:A:267:SER:OG	1:A:308:LEU:HD12	2.18	0.43
1:A:719:MET:HE3	1:A:743:ILE:CG1	2.47	0.43
1:A:641:TYR:C	1:A:642:MET:HE3	2.41	0.43
1:A:725:ILE:HD12	1:A:726:LEU:N	2.34	0.43
1:A:922:LEU:O	1:A:923:LYS:C	2.61	0.43
1:A:829:ILE:HD13	1:A:829:ILE:HA	1.78	0.43
1:A:1238:TYR:CZ	1:A:1240:MET:HE1	2.52	0.43
1:A:573:GLU:O	1:A:573:GLU:CG	2.67	0.43
1:A:1154:VAL:HG13	1:A:1200:VAL:O	2.19	0.43
1:A:67:ARG:HD2	1:A:83:LEU:O	2.19	0.42
1:A:507:GLN:HE22	1:A:602:THR:HG22	1.84	0.42
1:A:1162:VAL:CG1	1:A:1193:HIS:CE1	3.02	0.42
1:A:980:ILE:HD11	1:A:1013:ILE:CG2	2.48	0.42
1:A:1021:LEU:HD21	1:A:1038:TYR:CD2	2.55	0.42
1:A:1080:ARG:O	1:A:1081:LEU:HD23	2.19	0.42
1:A:518:ALA:O	1:A:565:LEU:HD12	2.19	0.42
1:A:535:LEU:HD11	1:A:666:TYR:OH	2.19	0.42
1:A:954:LEU:HB3	1:A:955:PRO:HD2	2.00	0.42
1:A:655:ASN:OD1	1:A:655:ASN:C	2.63	0.42
1:A:190:LEU:HD11	1:A:333:TYR:CB	2.50	0.42
1:A:742:ARG:HH11	1:A:742:ARG:HG2	1.83	0.42
1:A:942:PRO:O	1:A:943:LEU:HD13	2.18	0.42
1:A:1140:VAL:HB	1:A:1210:ALA:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1173:ASN:O	1:A:1174:ARG:C	2.62	0.42
1:A:414:ILE:N	1:A:414:ILE:CD1	2.82	0.42
1:A:500:ASN:HB3	1:A:598:SER:O	2.20	0.42
1:A:980:ILE:HG22	1:A:1012:ARG:CZ	2.49	0.42
1:A:1165:ASP:HA	1:A:1168:ALA:HB3	2.00	0.42
1:A:83:LEU:HD12	1:A:86:ARG:HD2	2.01	0.42
1:A:33:SER:HB3	1:A:115:LYS:O	2.20	0.42
1:A:22:MET:HE3	1:A:297:LEU:HD21	2.01	0.41
1:A:181:LYS:NZ	1:A:185:ASP:OD2	2.27	0.41
1:A:336:LEU:HD21	1:A:349:ILE:HD11	2.01	0.41
1:A:497:GLU:HB3	1:A:597:ALA:HB1	2.02	0.41
1:A:628:THR:HG22	1:A:630:ALA:N	2.35	0.41
1:A:360:LEU:HD22	1:A:360:LEU:H	1.85	0.41
1:A:908:VAL:HG11	1:A:941:LEU:HD12	2.02	0.41
1:A:520:TYR:HB2	1:A:565:LEU:HD11	2.03	0.41
1:A:908:VAL:HG11	1:A:940:GLU:HA	2.01	0.41
1:A:93:ILE:CD1	1:A:132:PHE:CE2	3.03	0.41
1:A:203:ARG:HH21	1:A:342:GLN:HE22	1.69	0.41
1:A:355:PHE:CE2	1:A:394:PRO:HD3	2.55	0.41
1:A:1025:LEU:HD12	1:A:1028:GLU:HB2	2.03	0.41
1:A:3:GLN:HE22	1:A:123:ARG:CZ	2.34	0.41
1:A:974:GLU:O	1:A:978:ALA:N	2.50	0.41
1:A:985:LEU:HD21	1:A:996:PHE:CE1	2.55	0.41
1:A:1250:MET:C	1:A:1251:LEU:HD23	2.46	0.41
1:A:276:ALA:HA	1:A:304:ARG:O	2.21	0.41
1:A:149:ILE:O	1:A:296:GLY:HA2	2.20	0.41
1:A:498:GLU:O	1:A:502:ILE:HD13	2.21	0.41
1:A:1114:LEU:HB2	1:A:1140:VAL:HA	2.02	0.41
1:A:28:LEU:HD22	1:A:1011:SER:HB2	2.02	0.40
1:A:106:LYS:O	1:A:109:GLU:HB3	2.21	0.40
1:A:611:TYR:CE1	1:A:632:ILE:HD11	2.55	0.40
1:A:685:ASP:C	1:A:685:ASP:OD1	2.64	0.40
1:A:828:PHE:O	1:A:829:ILE:HD13	2.20	0.40
1:A:611:TYR:CE1	1:A:662:THR:HG21	2.56	0.40
1:A:862:ASP:O	1:A:862:ASP:CG	2.64	0.40
1:A:1279:PHE:O	1:A:1283:VAL:HG23	2.20	0.40
1:A:3:GLN:HE22	1:A:123:ARG:NH1	2.19	0.40
1:A:354:VAL:O	1:A:391:MET:HA	2.21	0.40
1:A:950:ASN:OD1	1:A:950:ASN:N	2.54	0.40
1:A:1150:LYS:O	1:A:1227:LEU:HD23	2.22	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	1265/1304 (97%)	1121 (89%)	109 (9%)	35 (3%)	4 6

All (35) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	569	GLN
1	A	642	MET
1	A	910	ARG
1	A	1015	LYS
1	A	1016	GLU
1	A	1045	ASP
1	A	1269	THR
1	A	75	VAL
1	A	222	GLN
1	A	223	LYS
1	A	346	PRO
1	A	461	HIS
1	A	618	THR
1	A	1058	GLU
1	A	1129	ALA
1	A	1272	VAL
1	A	1289	SER
1	A	951	LYS
1	A	1270	GLN
1	A	343	ALA
1	A	344	ASP
1	A	347	LYS
1	A	463	THR
1	A	775	SER
1	A	898	SER
1	A	953	LEU
1	A	1019	ILE
1	A	562	ALA

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Mol	Chain	Res	Type
1	A	617	GLY
1	A	897	GLU
1	A	923	LYS
1	A	1022	PRO
1	A	1020	ASP
1	A	969	PRO
1	A	760	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	1077/1124 (96%)	940 (87%)	137 (13%)	4 9

All (137) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASP
1	A	25	HIS
1	A	37	GLU
1	A	39	ILE
1	A	50	LYS
1	A	59	ILE
1	A	73	GLU
1	A	74	LYS
1	A	75	VAL
1	A	113	GLN
1	A	123	ARG
1	A	164	LYS
1	A	171	GLU
1	A	180	VAL
1	A	187	ILE
1	A	190	LEU
1	A	195	LYS
1	A	217	GLU
1	A	222	GLN

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Mol	Chain	Res	Type
1	A	224	ASP
1	A	229	LYS
1	A	231	LEU
1	A	245	GLU
1	A	248	LYS
1	A	250	GLN
1	A	261	VAL
1	A	265	LEU
1	A	267	SER
1	A	271	GLN
1	A	280	VAL
1	A	287	GLU
1	A	289	LYS
1	A	292	GLU
1	A	306	VAL
1	A	309	SER
1	A	323	GLU
1	A	324	GLN
1	A	354	VAL
1	A	360	LEU
1	A	365	ASN
1	A	367	GLU
1	A	374	ASP
1	A	380	VAL
1	A	382	GLU
1	A	384	SER
1	A	385	ASN
1	A	389	ASN
1	A	390	LEU
1	A	391	MET
1	A	399	LEU
1	A	406	GLU
1	A	414	ILE
1	A	421	LEU
1	A	429	ILE
1	A	430	GLN
1	A	436	VAL
1	A	441	LEU
1	A	448	GLU
1	A	454	LEU
1	A	461	HIS
1	A	463	THR

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Mol	Chain	Res	Type
1	A	483	LEU
1	A	488	GLN
1	A	542	LEU
1	A	547	LYS
1	A	551	ASP
1	A	552	ARG
1	A	557	LEU
1	A	567	THR
1	A	573	GLU
1	A	581	THR
1	A	584	THR
1	A	587	ILE
1	A	588	ASP
1	A	590	GLN
1	A	592	ARG
1	A	615	THR
1	A	619	THR
1	A	628	THR
1	A	635	LEU
1	A	642	MET
1	A	656	TYR
1	A	676	ILE
1	A	694	LEU
1	A	697	ASN
1	A	722	LEU
1	A	725	ILE
1	A	726	LEU
1	A	742	ARG
1	A	749	LEU
1	A	770	LEU
1	A	774	ILE
1	A	787	SER
1	A	793	GLU
1	A	794	GLN
1	A	823	LEU
1	A	826	GLU
1	A	829	ILE
1	A	845	LEU
1	A	864	VAL
1	A	876	ILE
1	A	886	VAL
1	A	887	LYS

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Mol	Chain	Res	Type
1	A	895	ARG
1	A	903	ILE
1	A	924	LYS
1	A	940	GLU
1	A	941	LEU
1	A	943	LEU
1	A	948	LYS
1	A	950	ASN
1	A	952	ARG
1	A	953	LEU
1	A	961	GLN
1	A	967	ILE
1	A	977	ILE
1	A	979	GLN
1	A	992	ILE
1	A	999	LEU
1	A	1011	SER
1	A	1013	ILE
1	A	1016	GLU
1	A	1019	ILE
1	A	1025	LEU
1	A	1031	THR
1	A	1035	ILE
1	A	1048	GLN
1	A	1051	THR
1	A	1104	LEU
1	A	1144	ILE
1	A	1174	ARG
1	A	1241	LYS
1	A	1254	GLU
1	A	1269	THR
1	A	1270	GLN
1	A	1285	GLN
1	A	1287	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	32	GLN
1	A	222	GLN
1	A	270	GLN

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Mol	Chain	Res	Type
1	A	324	GLN
1	A	340	GLN
1	A	342	GLN
1	A	345	GLN
1	A	357	ASN
1	A	361	GLN
1	A	365	ASN
1	A	389	ASN
1	A	420	GLN
1	A	426	GLN
1	A	459	GLN
1	A	475	ASN
1	A	507	GLN
1	A	569	GLN
1	A	607	ASN
1	A	633	GLN
1	A	671	ASN
1	A	695	GLN
1	A	785	ASN
1	A	796	GLN
1	A	819	ASN
1	A	879	GLN
1	A	881	GLN
1	A	946	ASN
1	A	1048	GLN
1	A	1054	ASN
1	A	1059	GLN
1	A	1109	GLN
1	A	1136	GLN
1	A	1152	GLN
1	A	1171	ASN
1	A	1173	ASN
1	A	1201	ASN
1	A	1207	GLN

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](#)

2 ligands 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	$\# Z > 2$
3	SO4	A	2292	-	4,4,4	0.25	0	6,6,6	0.13	0
2	LEU	A	2291	-	6,8,8	0.77	0	5,10,10	1.95	1 (20%)

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	LEU	A	2291	-	-	6/8/8/8	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	2291	LEU	OXT-C-O	-4.27	114.39	124.08

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	2291	LEU	CA-CB-CG-CD2
2	A	2291	LEU	CA-CB-CG-CD1
2	A	2291	LEU	C-CA-CB-CG
2	A	2291	LEU	OXT-C-CA-CB

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Mol	Chain	Res	Type	Atoms
2	A	2291	LEU	O-C-CA-CB
2	A	2291	LEU	N-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

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	1273/1304 (97%)	0.50	46 (3%) 46 40	57, 70, 79, 110	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1223	ILE	3.5
1	A	939	ASP	3.4
1	A	1184	VAL	3.2
1	A	581	THR	3.1
1	A	1038	TYR	3.0
1	A	1043	GLY	3.0
1	A	888	ASP	2.9
1	A	360	LEU	2.9
1	A	1017	LEU	2.9
1	A	767	VAL	2.8
1	A	1272	VAL	2.8
1	A	2	SER	2.7
1	A	1021	LEU	2.7
1	A	1240	MET	2.7
1	A	940	GLU	2.7
1	A	804	GLY	2.7
1	A	860	ILE	2.6
1	A	1082	PRO	2.6
1	A	846	ALA	2.6
1	A	1169	LEU	2.5
1	A	1047	LEU	2.4
1	A	1183	ALA	2.4
1	A	1091	PHE	2.3
1	A	1230	TRP	2.3
1	A	587	ILE	2.3
1	A	865	LYS	2.3
1	A	1161	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	741	LEU	2.3
1	A	622	PRO	2.2
1	A	588	ASP	2.2
1	A	1023	VAL	2.2
1	A	929	TYR	2.2
1	A	1185	LYS	2.1
1	A	1244	PHE	2.1
1	A	950	ASN	2.1
1	A	1048	GLN	2.1
1	A	1020	ASP	2.1
1	A	143	TRP	2.1
1	A	153	TRP	2.1
1	A	1031	THR	2.1
1	A	778	PRO	2.1
1	A	1030	PRO	2.1
1	A	34	PHE	2.1
1	A	764	ALA	2.0
1	A	578	LEU	2.0
1	A	866	ILE	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](#)

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	SO4	A	2292	5/5	0.72	0.12	122,122,122,123	0
2	LEU	A	2291	9/9	0.77	0.24	77,78,79,79	0

6.5 Other polymers [i](#)

There are no such residues in this entry.