



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:38 AM UTC

PDB ID : 6AC8 / pdb_00006ac8
Title : Crystal structure of Mycobacterium smegmatis Mfd at 2.75 Å resolution
Authors : Putta, S.; Fox, G.C.; Walsh, M.A.; Rao, D.N.; Nagaraja, V.; Natesh, R.
Deposited on : 2018-07-25
Resolution : 2.75 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

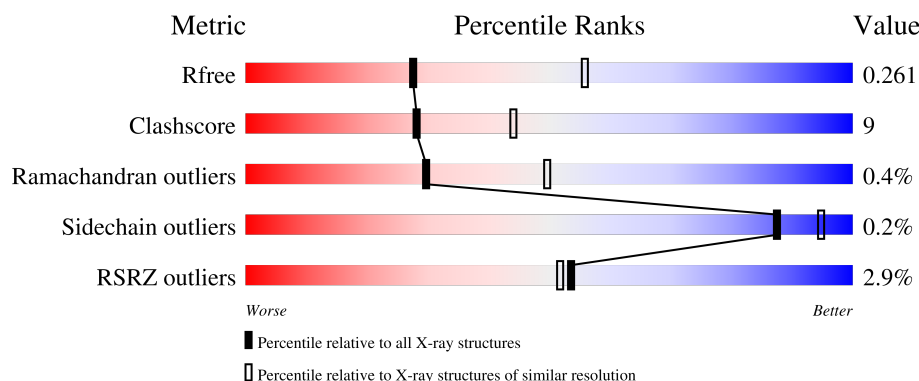
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	1235	
1	B	1235	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	1307	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17331 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 Mycobacterium smegmatis Mfd.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1171	Total	C	N	O	S	0	0	0
			8590	5395	1526	1639	30			
1	B	1174	Total	C	N	O	S	0	0	0
			8566	5391	1519	1626	30			

There are 40 discrepancies between the modelled and reference sequences:

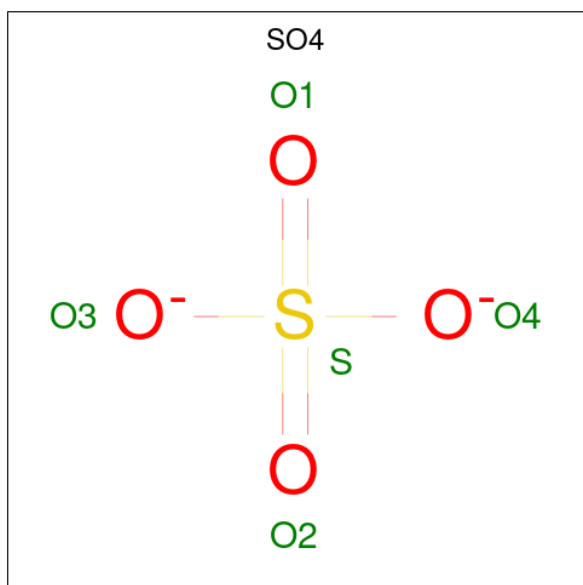
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP I7G7M2
A	-18	GLY	-	expression tag	UNP I7G7M2
A	-17	SER	-	expression tag	UNP I7G7M2
A	-16	SER	-	expression tag	UNP I7G7M2
A	-15	HIS	-	expression tag	UNP I7G7M2
A	-14	HIS	-	expression tag	UNP I7G7M2
A	-13	HIS	-	expression tag	UNP I7G7M2
A	-12	HIS	-	expression tag	UNP I7G7M2
A	-11	HIS	-	expression tag	UNP I7G7M2
A	-10	HIS	-	expression tag	UNP I7G7M2
A	-9	SER	-	expression tag	UNP I7G7M2
A	-8	SER	-	expression tag	UNP I7G7M2
A	-7	GLY	-	expression tag	UNP I7G7M2
A	-6	LEU	-	expression tag	UNP I7G7M2
A	-5	VAL	-	expression tag	UNP I7G7M2
A	-4	PRO	-	expression tag	UNP I7G7M2
A	-3	ARG	-	expression tag	UNP I7G7M2
A	-2	GLY	-	expression tag	UNP I7G7M2
A	-1	SER	-	expression tag	UNP I7G7M2
A	0	HIS	-	expression tag	UNP I7G7M2
B	-19	MET	-	initiating methionine	UNP I7G7M2
B	-18	GLY	-	expression tag	UNP I7G7M2
B	-17	SER	-	expression tag	UNP I7G7M2
B	-16	SER	-	expression tag	UNP I7G7M2
B	-15	HIS	-	expression tag	UNP I7G7M2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP I7G7M2
B	-13	HIS	-	expression tag	UNP I7G7M2
B	-12	HIS	-	expression tag	UNP I7G7M2
B	-11	HIS	-	expression tag	UNP I7G7M2
B	-10	HIS	-	expression tag	UNP I7G7M2
B	-9	SER	-	expression tag	UNP I7G7M2
B	-8	SER	-	expression tag	UNP I7G7M2
B	-7	GLY	-	expression tag	UNP I7G7M2
B	-6	LEU	-	expression tag	UNP I7G7M2
B	-5	VAL	-	expression tag	UNP I7G7M2
B	-4	PRO	-	expression tag	UNP I7G7M2
B	-3	ARG	-	expression tag	UNP I7G7M2
B	-2	GLY	-	expression tag	UNP I7G7M2
B	-1	SER	-	expression tag	UNP I7G7M2
B	0	HIS	-	expression tag	UNP I7G7M2

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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

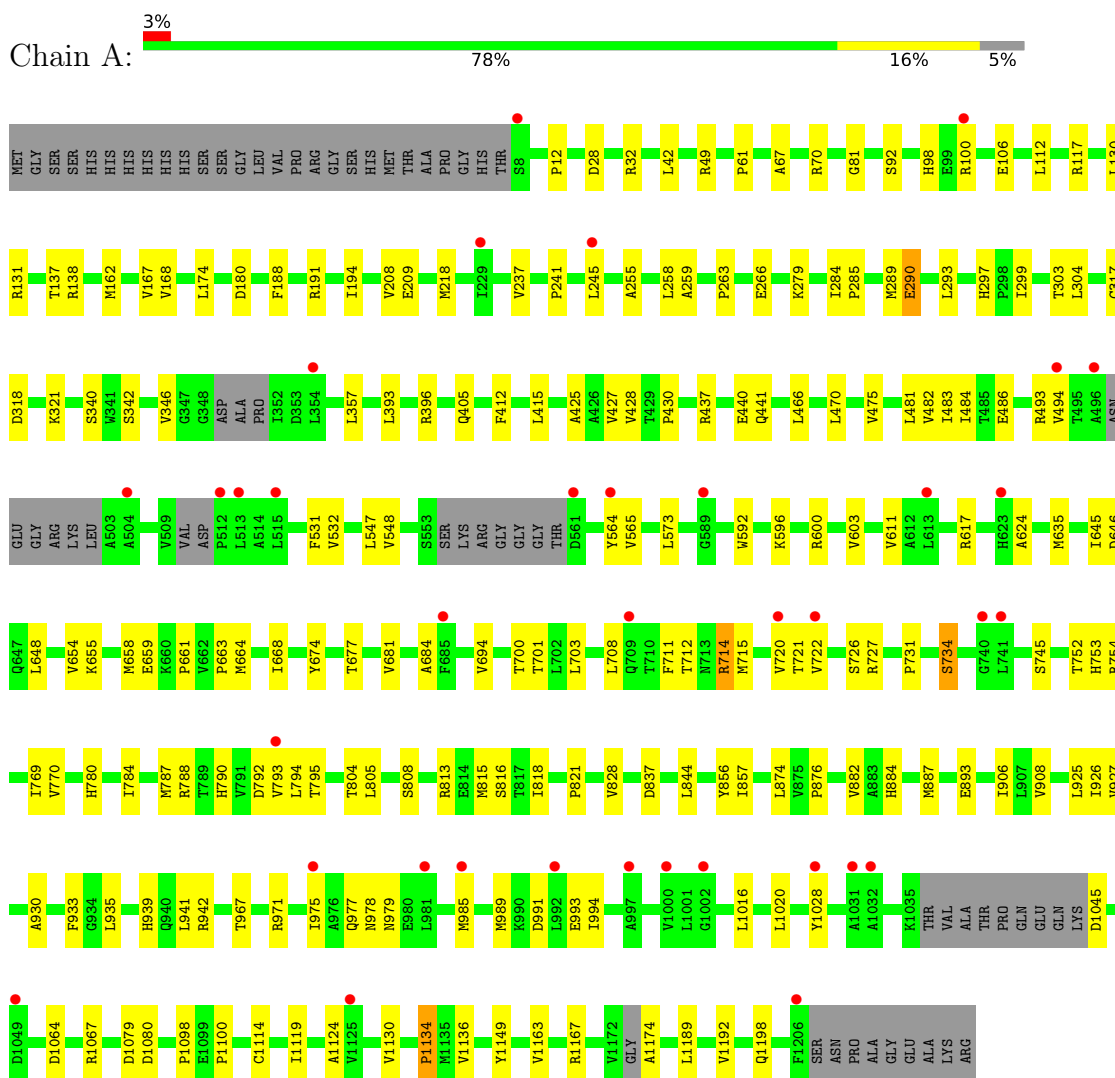
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	51	Total O 51 51	0	0
3	B	69	Total O 69 69	0	0

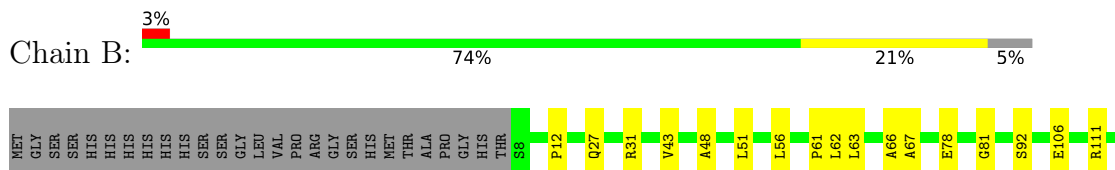
3 Residue-property plots

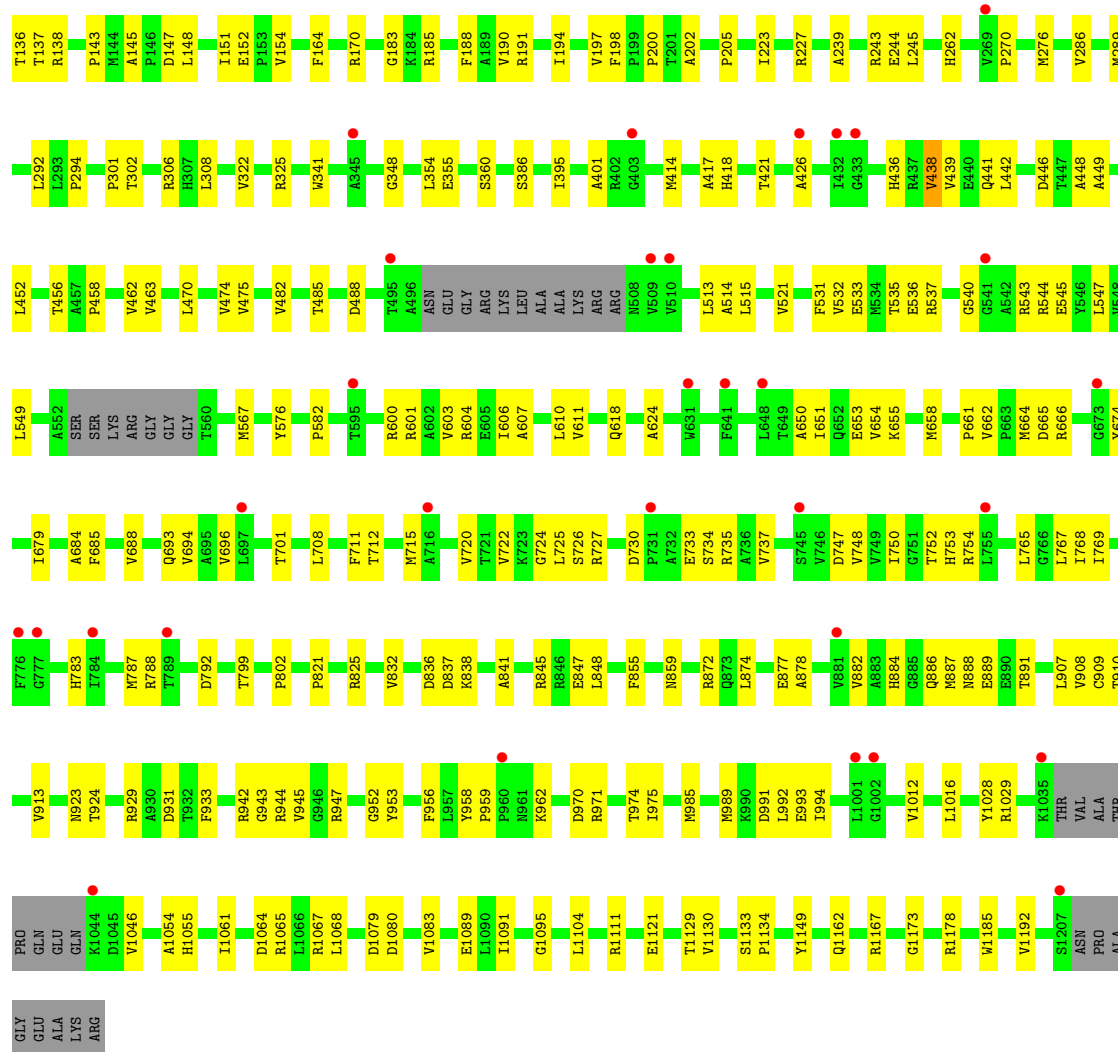
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Mycobacterium smegmatis Mfd





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.86Å 162.03Å 215.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.51 – 2.75 40.51 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.51-2.75) 99.3 (40.51-2.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.77Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575:000)	Depositor
R, R_{free}	0.224 , 0.259 0.225 , 0.261	Depositor DCC
R_{free} test set	3816 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	80.9	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17331	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.16	0/8744	0.34	0/11944
1	B	0.16	0/8721	0.36	0/11924
All	All	0.16	0/17465	0.35	0/23868

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	8590	0	8318	130	0
1	B	8566	0	8336	172	0
2	A	35	0	0	3	0
2	B	20	0	0	0	0
3	A	51	0	0	0	0
3	B	69	0	0	2	0
All	All	17331	0	16654	301	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 9.

All (301) 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:712:THR:HG22	1:A:722:VAL:HG21	1.54	0.90
1:B:712:THR:HG22	1:B:722:VAL:HG21	1.53	0.89
1:A:635:MET:SD	1:A:714:ARG:NH2	2.48	0.87
1:B:825:ARG:HD2	1:B:947:ARG:O	1.81	0.81
1:A:942:ARG:HH22	1:A:978:ASN:HB2	1.47	0.79
1:B:545:GLU:HB2	1:B:567:MET:HE3	1.64	0.78
1:A:727:ARG:NH2	1:A:893:GLU:OE2	2.19	0.75
1:B:730:ASP:HB2	1:B:733:GLU:H	1.51	0.73
1:B:243:ARG:HD3	1:B:289:MET:HE2	1.71	0.73
1:A:92:SER:HB2	1:A:138:ARG:HD2	1.71	0.72
1:B:92:SER:HB2	1:B:138:ARG:HD2	1.72	0.71
1:B:545:GLU:O	1:B:567:MET:HE2	1.90	0.71
1:B:545:GLU:HB2	1:B:567:MET:CE	2.21	0.70
1:A:668:ILE:HG22	1:A:816:SER:HB2	1.73	0.70
1:A:971:ARG:NH1	2:A:1307:SO4:O4	2.24	0.69
1:A:318:ASP:HB3	1:A:321:LYS:HD2	1.74	0.68
1:A:770:VAL:HB	1:A:795:THR:HG22	1.76	0.68
1:B:708:LEU:HD22	1:B:724:GLY:HA3	1.74	0.68
1:A:28:ASP:O	1:A:32:ARG:HG3	1.94	0.68
1:B:933:PHE:O	1:B:971:ARG:NH2	2.27	0.67
1:B:545:GLU:C	1:B:567:MET:HE2	2.20	0.67
1:A:437:ARG:NH1	1:A:440:GLU:OE1	2.27	0.67
1:B:654:VAL:HG12	1:B:666:ARG:HD2	1.76	0.67
1:A:991:ASP:HA	1:A:994:ILE:HG12	1.76	0.66
1:B:765:LEU:HD21	1:B:768:ILE:HD11	1.77	0.66
1:B:438:VAL:O	1:B:441:GLN:N	2.29	0.66
1:B:601:ARG:HA	1:B:604:ARG:HD2	1.77	0.66
1:B:696:VAL:HG22	1:B:769:ILE:HB	1.79	0.65
1:A:100:ARG:NH2	1:A:340:SER:O	2.29	0.65
1:A:727:ARG:NH1	2:A:1304:SO4:O3	2.30	0.64
1:B:943:GLY:O	1:B:947:ARG:NE	2.29	0.64
1:B:701:THR:HG23	1:B:726:SER:HB3	1.79	0.64
1:A:194:ILE:HG12	1:A:209:GLU:HG2	1.80	0.63
1:B:439:VAL:HA	1:B:442:LEU:HD23	1.80	0.63
1:A:430:PRO:HD2	1:A:486:GLU:HG2	1.80	0.63
1:B:694:VAL:HG22	1:B:767:LEU:HB3	1.80	0.63
1:B:720:VAL:HG23	1:B:747:ASP:OD2	1.97	0.63
1:B:532:VAL:HG23	1:B:533:GLU:HG2	1.81	0.62
1:B:436:HIS:HA	1:B:439:VAL:HG22	1.82	0.62
1:A:112:LEU:HD13	1:A:241:PRO:HG2	1.81	0.62
1:A:208:VAL:HG12	1:A:218:MET:HG2	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:664:MET:HB3	1:B:792:ASP:HB3	1.82	0.62
1:B:276:MET:HG3	1:B:286:VAL:HG21	1.80	0.61
1:B:540:GLY:O	1:B:1178:ARG:NH2	2.32	0.61
1:A:117:ARG:HB3	1:A:130:LEU:HD21	1.82	0.60
1:A:1079:ASP:OD2	1:A:1080:ASP:N	2.33	0.60
1:A:708:LEU:O	1:A:712:THR:HG23	2.02	0.60
1:B:942:ARG:NH1	1:B:975:ILE:O	2.32	0.60
1:A:977:GLN:OE1	1:A:977:GLN:N	2.33	0.60
1:A:1149:TYR:HE2	1:A:1192:VAL:HG21	1.66	0.60
1:B:752:THR:OG1	1:B:753:HIS:N	2.30	0.60
1:A:645:ILE:HD12	1:A:648:LEU:HD11	1.84	0.59
1:A:437:ARG:NH1	1:A:441:GLN:HG3	2.18	0.59
1:A:857:ILE:HD13	1:A:941:LEU:HD13	1.84	0.59
1:B:708:LEU:O	1:B:712:THR:HG23	2.01	0.59
1:A:255:ALA:O	1:A:259:ALA:N	2.36	0.59
1:B:414:MET:O	1:B:417:ALA:N	2.36	0.58
1:B:611:VAL:HG11	1:B:1028:TYR:HB3	1.86	0.58
1:B:448:ALA:HB1	1:B:462:VAL:HA	1.86	0.58
1:B:148:LEU:HA	1:B:151:ILE:CD1	2.34	0.58
1:B:245:LEU:HD23	1:B:289:MET:HE3	1.85	0.57
1:B:475:VAL:HG22	1:B:482:VAL:HG13	1.85	0.57
1:B:188:PHE:HB3	1:B:197:VAL:HG23	1.87	0.57
1:A:715:MET:HB3	1:A:720:VAL:HG21	1.85	0.57
1:A:1064:ASP:OD1	1:A:1067:ARG:NH2	2.37	0.57
1:B:292:LEU:HD21	1:B:341:TRP:CH2	2.40	0.57
1:B:426:ALA:HB2	1:B:458:PRO:HG3	1.88	0.56
1:B:148:LEU:HA	1:B:151:ILE:HD13	1.88	0.56
1:B:151:ILE:HD12	1:B:151:ILE:N	2.20	0.56
1:B:725:LEU:HD21	1:B:754:ARG:HE	1.71	0.56
1:A:290:GLU:HA	1:A:293:LEU:HD13	1.87	0.56
1:B:832:VAL:HG12	1:B:956:PHE:HB2	1.87	0.56
1:B:154:VAL:HB	1:B:239:ALA:HB3	1.87	0.55
1:A:787:MET:HA	1:A:790:HIS:ND1	2.21	0.55
1:B:730:ASP:HB2	1:B:733:GLU:HB2	1.88	0.55
1:A:405:GLN:OE1	1:A:405:GLN:N	2.39	0.55
1:B:513:LEU:HD22	1:B:1068:LEU:HD13	1.89	0.55
1:B:923:ASN:HA	1:B:945:VAL:HG22	1.89	0.55
1:A:788:ARG:HH21	1:A:795:THR:HG21	1.72	0.54
1:A:654:VAL:HG12	1:A:658:MET:HE2	1.89	0.54
1:B:841:ALA:HB2	1:B:874:LEU:HD22	1.89	0.54
1:B:607:ALA:O	1:B:611:VAL:HG12	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:848:LEU:HD11	1:B:878:ALA:HB2	1.90	0.53
1:B:51:LEU:HD11	1:B:474:VAL:HG11	1.89	0.53
1:B:712:THR:HG22	1:B:722:VAL:CG2	2.32	0.53
1:A:266:GLU:CD	1:A:266:GLU:H	2.17	0.53
1:B:137:THR:HB	1:B:322:VAL:HG13	1.91	0.53
1:A:285:PRO:HB3	1:A:289:MET:HE1	1.90	0.53
1:A:967:THR:HG23	2:A:1307:SO4:O4	2.08	0.53
1:A:218:MET:HE1	1:A:237:VAL:HG22	1.89	0.53
1:A:731:PRO:HA	1:A:734:SER:HB3	1.91	0.53
1:B:147:ASP:O	1:B:151:ILE:HD12	2.09	0.53
1:B:78:GLU:HG2	1:B:470:LEU:HA	1.90	0.52
1:B:799:THR:OG1	1:B:992:LEU:O	2.27	0.52
1:B:924:THR:HG22	1:B:953:TYR:HD2	1.75	0.52
1:A:303:THR:OG1	1:A:304:LEU:N	2.43	0.52
1:B:185:ARG:NH2	1:B:202:ALA:O	2.42	0.52
1:A:475:VAL:HG13	1:A:482:VAL:HG22	1.90	0.52
1:B:531:PHE:HE1	1:B:547:LEU:HD11	1.74	0.52
1:B:536:GLU:HG2	1:B:543:ARG:NH2	2.25	0.52
1:B:970:ASP:O	1:B:974:THR:HG23	2.09	0.52
1:B:531:PHE:CE1	1:B:547:LEU:HD11	2.44	0.52
1:B:545:GLU:CB	1:B:567:MET:CE	2.87	0.52
1:B:439:VAL:HG12	1:B:449:ALA:HB1	1.92	0.52
1:A:49:ARG:HE	1:A:317:CYS:HB3	1.74	0.52
1:B:715:MET:HE1	1:B:748:VAL:HG21	1.91	0.52
1:A:168:VAL:HG23	1:A:188:PHE:HE2	1.75	0.51
1:B:12:PRO:HD2	1:B:81:GLY:HA2	1.91	0.51
1:B:884:HIS:CE1	1:B:887:MET:HG2	2.45	0.51
1:A:415:LEU:HD22	1:A:425:ALA:HB1	1.92	0.51
1:B:56:LEU:HB2	1:B:62:LEU:HD11	1.92	0.51
1:A:12:PRO:HD2	1:A:81:GLY:HA2	1.91	0.51
1:A:884:HIS:CE1	1:A:887:MET:HG3	2.45	0.51
1:B:111:ARG:NH1	1:B:244:GLU:OE2	2.43	0.51
1:B:655:LYS:HA	1:B:658:MET:HE3	1.90	0.51
1:B:708:LEU:HA	1:B:750:ILE:HD12	1.91	0.51
1:B:931:ASP:OD1	1:B:931:ASP:N	2.43	0.51
1:A:106:GLU:CD	1:A:106:GLU:H	2.18	0.51
1:A:646:ASP:HB2	1:A:818:ILE:HG23	1.92	0.51
1:A:180:ASP:O	1:A:191:ARG:NH1	2.33	0.51
1:A:684:ALA:HB1	1:A:694:VAL:HG21	1.92	0.51
1:B:48:ALA:HB2	1:B:395:ILE:HG21	1.93	0.51
1:B:653:GLU:OE1	1:B:666:ARG:NH1	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:721:THR:OG1	1:A:745:SER:O	2.27	0.50
1:A:828:VAL:HG22	1:A:942:ARG:HD3	1.94	0.50
1:B:685:PHE:HA	1:B:688:VAL:HG22	1.94	0.50
1:B:1130:VAL:HG11	1:B:1185:TRP:HZ3	1.77	0.50
1:A:664:MET:HE2	1:A:792:ASP:HB3	1.92	0.50
1:A:681:VAL:HG12	1:A:711:PHE:CE1	2.47	0.50
1:A:787:MET:HA	1:A:790:HIS:CE1	2.47	0.50
1:A:1098:PRO:HB2	1:A:1100:PRO:HD2	1.94	0.50
1:A:481:LEU:HD21	1:A:483:ILE:HD11	1.94	0.50
1:B:847:GLU:OE1	1:B:924:THR:HG23	2.11	0.50
1:A:412:PHE:HE2	1:A:441:GLN:HB3	1.77	0.50
1:A:701:THR:HB	1:A:726:SER:HB3	1.93	0.50
1:B:27:GLN:O	1:B:31:ARG:HG2	2.11	0.50
1:B:164:PHE:CZ	1:B:190:VAL:HG23	2.47	0.50
1:B:1046:VAL:HB	1:B:1111:ARG:HG2	1.93	0.50
1:B:545:GLU:C	1:B:567:MET:CE	2.84	0.50
1:A:168:VAL:HG23	1:A:188:PHE:CE2	2.46	0.49
1:A:592:TRP:NE1	1:A:596:LYS:HD2	2.27	0.49
1:B:959:PRO:HB2	1:B:962:LYS:HG2	1.94	0.49
1:A:42:LEU:HB3	1:A:393:LEU:HD11	1.95	0.49
1:A:297:HIS:CE1	1:A:299:ILE:HD11	2.47	0.49
1:B:1054:ALA:HB1	1:B:1104:LEU:HA	1.94	0.49
1:B:43:VAL:HG11	1:B:386:SER:HA	1.94	0.49
1:B:270:PRO:O	1:B:958:TYR:OH	2.26	0.49
1:B:734:SER:HB2	1:B:754:ARG:HH21	1.77	0.49
1:B:624:ALA:HB2	1:B:661:PRO:HA	1.95	0.49
1:B:414:MET:HE2	1:B:418:HIS:HB2	1.94	0.49
1:B:1121:GLU:HG2	1:B:1133:SER:HB3	1.95	0.48
1:A:734:SER:OG	1:A:754:ARG:NH2	2.46	0.48
1:B:137:THR:HG21	1:B:325:ARG:HG2	1.95	0.48
1:B:747:ASP:OD1	1:B:747:ASP:N	2.36	0.48
1:B:802:PRO:HG3	1:B:993:GLU:HB3	1.95	0.48
1:A:780:HIS:O	1:A:784:ILE:HG13	2.12	0.48
1:A:857:ILE:HD12	1:A:927:VAL:HG22	1.95	0.48
1:B:783:HIS:CE1	1:B:787:MET:HE3	2.48	0.48
1:B:882:VAL:HA	1:B:908:VAL:O	2.13	0.48
1:A:1198:GLN:OE1	1:A:1198:GLN:N	2.43	0.48
1:B:148:LEU:O	1:B:151:ILE:HD13	2.14	0.48
1:B:1061:ILE:HB	1:B:1067:ARG:HG3	1.95	0.48
1:B:1079:ASP:OD1	1:B:1080:ASP:N	2.46	0.48
1:A:655:LYS:O	1:A:659:GLU:HG2	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:752:THR:OG1	1:A:753:HIS:N	2.45	0.48
1:B:600:ARG:HA	1:B:603:VAL:HG22	1.96	0.48
1:A:162:MET:HE2	1:A:167:VAL:CG2	2.44	0.48
1:A:813:ARG:HD3	1:A:815:MET:HE3	1.96	0.48
1:A:844:LEU:HD13	1:A:906:ILE:HD13	1.96	0.48
1:B:536:GLU:HG3	1:B:545:GLU:HG2	1.95	0.48
1:B:725:LEU:C	1:B:725:LEU:HD23	2.39	0.47
1:A:844:LEU:HD21	1:A:926:ILE:HG13	1.96	0.47
1:A:1167:ARG:NH1	1:A:1174:ALA:O	2.42	0.47
1:B:684:ALA:HA	1:B:694:VAL:HG21	1.96	0.47
1:A:342:SER:O	1:A:346:VAL:HG23	2.14	0.47
1:B:67:ALA:O	1:B:137:THR:HG22	2.15	0.47
1:B:401:ALA:N	1:B:488:ASP:O	2.47	0.47
1:B:576:TYR:CZ	1:B:582:PRO:HG3	2.49	0.47
1:B:535:THR:HG21	1:B:537:ARG:CZ	2.44	0.47
1:B:1029:ARG:CZ	1:B:1029:ARG:HB2	2.44	0.47
1:B:837:ASP:HB3	1:B:874:LEU:HD21	1.96	0.47
1:A:935:LEU:H	1:A:935:LEU:HD12	1.79	0.47
1:A:882:VAL:HA	1:A:908:VAL:O	2.14	0.47
1:B:544:ARG:NH1	3:B:1402:HOH:O	2.47	0.47
1:A:788:ARG:HG2	1:A:793:VAL:HG11	1.97	0.47
1:A:1167:ARG:HD2	1:A:1174:ALA:HB3	1.97	0.47
1:B:66:ALA:O	1:B:136:THR:HA	2.15	0.47
1:A:700:THR:HG23	1:A:703:LEU:H	1.80	0.46
1:B:536:GLU:HG2	1:B:543:ARG:HH21	1.79	0.46
1:A:405:GLN:N	1:A:405:GLN:CD	2.73	0.46
1:A:655:LYS:HA	1:A:658:MET:HE3	1.96	0.46
1:B:470:LEU:HD23	1:B:485:THR:CG2	2.45	0.46
1:B:929:ARG:HG3	1:B:931:ASP:OD1	2.14	0.46
1:B:1064:ASP:OD1	1:B:1067:ARG:NH2	2.48	0.46
1:A:174:LEU:HB3	1:A:241:PRO:HA	1.97	0.46
1:A:611:VAL:HG21	1:A:1028:TYR:HB3	1.98	0.46
1:A:1045:ASP:N	1:A:1045:ASP:OD1	2.47	0.46
1:B:452:LEU:HD23	1:B:456:THR:HG23	1.97	0.46
1:B:836:ASP:OD2	1:B:838:LYS:N	2.49	0.46
1:A:266:GLU:OE1	1:A:266:GLU:N	2.40	0.46
1:A:856:TYR:O	1:A:908:VAL:HA	2.15	0.46
1:B:989:MET:HA	1:B:992:LEU:HD13	1.97	0.46
1:A:856:TYR:HB3	1:A:908:VAL:HG22	1.97	0.46
1:B:515:LEU:HD23	1:B:515:LEU:HA	1.79	0.46
1:A:677:THR:O	1:A:681:VAL:HG13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:985:MET:O	1:B:989:MET:HG2	2.15	0.46
1:B:191:ARG:HB2	1:B:194:ILE:HB	1.97	0.46
1:B:355:GLU:OE2	1:B:355:GLU:N	2.35	0.46
1:B:674:TYR:CE1	1:B:821:PRO:HB3	2.51	0.45
1:B:185:ARG:HB3	1:B:200:PRO:HA	1.97	0.45
1:A:1016:LEU:O	1:A:1020:LEU:HG	2.17	0.45
1:B:227:ARG:HD3	1:B:1089:GLU:HG3	1.97	0.45
1:B:354:LEU:HA	1:B:354:LEU:HD23	1.63	0.45
1:B:944:ARG:HA	1:B:944:ARG:HH21	1.80	0.45
1:A:67:ALA:O	1:A:137:THR:HG23	2.17	0.45
1:A:837:ASP:HB3	1:A:874:LEU:HD21	1.99	0.45
1:B:514:ALA:HB2	1:B:1065:ARG:HD3	1.98	0.45
1:B:1012:VAL:HG22	1:B:1016:LEU:HD23	1.97	0.45
1:A:531:PHE:HE1	1:A:547:LEU:HG	1.82	0.45
1:A:548:VAL:HG12	1:A:564:TYR:CD1	2.52	0.44
1:A:70:ARG:HH22	1:A:493:ARG:HD3	1.82	0.44
1:B:991:ASP:HA	1:B:994:ILE:HB	1.98	0.44
1:B:1091:ILE:HA	1:B:1095:GLY:O	2.17	0.44
1:B:1167:ARG:NH1	1:B:1173:GLY:H	2.15	0.44
1:A:162:MET:HE2	1:A:167:VAL:HG21	1.99	0.44
1:B:63:LEU:HD22	1:B:308:LEU:HD11	1.99	0.44
1:A:939:HIS:NE2	1:A:985:MET:HE2	2.33	0.44
1:B:164:PHE:HZ	1:B:190:VAL:HG23	1.82	0.44
1:B:650:ALA:O	1:B:654:VAL:HG13	2.17	0.44
1:B:294:PRO:HG3	1:B:360:SER:HB2	2.00	0.44
1:B:417:ALA:O	1:B:421:THR:OG1	2.22	0.44
1:B:600:ARG:O	1:B:604:ARG:NE	2.51	0.44
1:B:143:PRO:HB2	1:B:301:PRO:HB2	1.99	0.44
1:A:61:PRO:HA	1:A:131:ARG:O	2.18	0.44
1:A:396:ARG:CZ	1:A:396:ARG:HB3	2.47	0.44
1:A:674:TYR:CZ	1:A:821:PRO:HB3	2.53	0.44
1:B:543:ARG:HB3	1:B:1055:HIS:CD2	2.52	0.43
1:B:606:ILE:O	1:B:610:LEU:HD23	2.18	0.43
1:B:1080:ASP:HA	1:B:1083:VAL:HG22	2.01	0.43
1:B:665:ASP:CG	1:B:788:ARG:HD3	2.44	0.43
1:A:857:ILE:CD1	1:A:925:LEU:HD11	2.47	0.43
1:A:804:THR:OG1	1:A:805:LEU:N	2.50	0.43
1:A:547:LEU:HB3	1:A:565:VAL:HG22	2.00	0.43
1:A:989:MET:O	1:A:993:GLU:HG3	2.18	0.43
1:B:1149:TYR:CE1	1:B:1192:VAL:HG21	2.54	0.43
1:A:930:ALA:HA	1:A:933:PHE:HD2	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:975:ILE:HD12	1:A:975:ILE:HA	1.89	0.43
1:B:106:GLU:HG3	1:B:185:ARG:HD2	2.00	0.43
1:A:600:ARG:O	1:A:603:VAL:HG12	2.18	0.42
1:B:923:ASN:O	1:B:952:GLY:HA2	2.19	0.42
1:A:98:HIS:NE2	1:A:357:LEU:O	2.51	0.42
1:B:909:CYS:HB2	1:B:913:VAL:CG1	2.48	0.42
1:B:145:ALA:HB3	1:B:148:LEU:HG	2.00	0.42
1:B:884:HIS:CE1	1:B:886:GLN:HB2	2.54	0.42
1:A:279:LYS:HB3	1:A:284:ILE:HG13	2.02	0.42
1:B:183:GLY:HA2	1:B:223:ILE:HD12	2.00	0.42
1:B:693:GLN:HB3	1:B:747:ASP:HA	2.01	0.42
1:A:428:VAL:HA	1:A:466:LEU:O	2.19	0.42
1:A:770:VAL:O	1:A:795:THR:HA	2.20	0.42
1:A:828:VAL:CG2	1:A:942:ARG:HD3	2.49	0.42
1:B:664:MET:HE2	1:B:792:ASP:HB3	2.02	0.42
1:B:1149:TYR:HE1	1:B:1192:VAL:HG21	1.85	0.42
1:A:769:ILE:HA	1:A:794:LEU:O	2.20	0.42
1:B:449:ALA:HA	1:B:463:VAL:O	2.20	0.42
1:A:412:PHE:CE2	1:A:441:GLN:HB3	2.54	0.42
1:A:1114:CYS:HB3	1:A:1119:ILE:HB	2.02	0.42
1:A:624:ALA:HB2	1:A:661:PRO:HA	2.00	0.42
1:B:576:TYR:CE2	1:B:582:PRO:HG3	2.55	0.42
1:A:1163:VAL:HG21	1:A:1189:LEU:HD22	2.01	0.41
1:B:618:GLN:HE21	1:B:662:VAL:HG21	1.84	0.41
1:B:152:GLU:O	1:B:170:ARG:NH2	2.47	0.41
1:B:262:HIS:ND1	1:B:348:GLY:HA3	2.35	0.41
1:A:532:VAL:HB	1:A:548:VAL:HG23	2.02	0.41
1:A:258:LEU:HD23	1:A:258:LEU:HA	1.89	0.41
1:B:61:PRO:HD2	3:B:1416:HOH:O	2.20	0.41
1:B:302:THR:OG1	1:B:306:ARG:HD2	2.20	0.41
1:B:651:ILE:HG13	1:B:679:ILE:HG12	2.03	0.41
1:B:727:ARG:NH2	1:B:889:GLU:HG3	2.36	0.41
1:B:888:ASN:ND2	1:B:891:THR:HG23	2.35	0.41
1:A:1124:ALA:HA	1:A:1130:VAL:HA	2.03	0.41
1:B:1129:THR:HG21	1:B:1162:GLN:HG3	2.03	0.41
1:B:711:PHE:CG	1:B:750:ILE:HD11	2.56	0.41
1:A:245:LEU:CD2	1:A:293:LEU:HD11	2.51	0.41
1:A:1119:ILE:HA	1:A:1134:PRO:HG2	2.03	0.41
1:A:617:ARG:HG3	1:A:663:PRO:HG2	2.02	0.41
1:B:448:ALA:O	1:B:463:VAL:N	2.54	0.41
1:A:805:LEU:HA	1:A:808:SER:OG	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:845:ARG:NE	1:B:877:GLU:OE2	2.52	0.41
1:B:859:ASN:HA	1:B:910:THR:HG22	2.02	0.41
1:A:428:VAL:HG11	1:A:470:LEU:HB2	2.03	0.40
1:A:857:ILE:HD12	1:A:925:LEU:HD11	2.03	0.40
1:A:565:VAL:HG21	1:A:573:LEU:HD11	2.03	0.40
1:A:427:VAL:HA	1:A:484:ILE:O	2.20	0.40
1:B:604:ARG:H	1:B:604:ARG:HG2	1.62	0.40
1:B:734:SER:OG	1:B:735:ARG:N	2.54	0.40
1:B:855:PHE:HE1	1:B:907:LEU:HD23	1.87	0.40
1:A:49:ARG:NH2	1:A:318:ASP:OD2	2.55	0.40
1:B:198:PHE:CD2	1:B:205:PRO:HG3	2.56	0.40
1:B:521:VAL:HG11	1:B:549:LEU:HD22	2.03	0.40
1:B:825:ARG:CD	1:B:947:ARG:O	2.62	0.40
1:A:876:PRO:HB3	1:B:872:ARG:NH2	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	1157/1235 (94%)	1096 (95%)	56 (5%)	5 (0%)	30	50
1	B	1166/1235 (94%)	1102 (94%)	60 (5%)	4 (0%)	36	56
All	All	2323/2470 (94%)	2198 (95%)	116 (5%)	9 (0%)	30	50

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	494	VAL
1	B	737	VAL
1	A	263	PRO
1	A	979	ASN

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Mol	Chain	Res	Type
1	A	734	SER
1	B	446	ASP
1	A	1134	PRO
1	B	438	VAL
1	B	1134	PRO

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	848/988 (86%)	845 (100%)	3 (0%)	84	91
1	B	847/988 (86%)	847 (100%)	0	100	100
All	All	1695/1976 (86%)	1692 (100%)	3 (0%)	87	95

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	290	GLU
1	A	714	ARG
1	A	1136	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	526	HIS
1	A	839	GLN
1	A	884	HIS
1	A	1194	ASN
1	B	297	HIS
1	B	307	HIS
1	B	572	GLN
1	B	783	HIS
1	B	979	ASN

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Mol	Chain	Res	Type
1	B	1055	HIS
1	B	1198	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

11 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	$\# Z > 2$
2	SO4	A	1303	-	4,4,4	0.25	0	6,6,6	0.10	0
2	SO4	A	1307	-	4,4,4	0.26	0	6,6,6	0.05	0
2	SO4	A	1305	-	4,4,4	0.23	0	6,6,6	0.11	0
2	SO4	B	1302	-	4,4,4	0.23	0	6,6,6	0.09	0
2	SO4	B	1303	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	A	1304	-	4,4,4	0.24	0	6,6,6	0.06	0
2	SO4	A	1301	-	4,4,4	0.25	0	6,6,6	0.10	0
2	SO4	B	1301	-	4,4,4	0.25	0	6,6,6	0.09	0
2	SO4	A	1306	-	4,4,4	0.25	0	6,6,6	0.12	0
2	SO4	A	1302	-	4,4,4	0.25	0	6,6,6	0.09	0
2	SO4	B	1304	-	4,4,4	0.24	0	6,6,6	0.08	0

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.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1307	SO4	2	0
2	A	1304	SO4	1	0

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	1171/1235 (94%)	0.26	36 (3%) 51 49	47, 80, 117, 143	0
1	B	1174/1235 (95%)	0.30	31 (2%) 57 55	48, 86, 126, 141	0
All	All	2345/2470 (94%)	0.28	67 (2%) 53 52	47, 82, 123, 143	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	496	ALA	4.2
1	B	269	VAL	3.4
1	B	509	VAL	3.4
1	A	504	ALA	3.3
1	B	1035	LYS	3.3
1	A	354	LEU	3.2
1	B	510	VAL	3.2
1	B	426	ALA	3.2
1	A	512	PRO	3.1
1	B	641	PHE	3.1
1	B	595	THR	3.0
1	A	975	ILE	2.8
1	B	745	SER	2.8
1	A	709	GLN	2.7
1	A	515	LEU	2.6
1	A	985	MET	2.6
1	A	997	ALA	2.6
1	B	1207	SER	2.5
1	A	740	GLY	2.5
1	A	8	SER	2.5
1	B	648	LEU	2.5
1	B	755	LEU	2.5
1	A	722	VAL	2.5
1	B	1002	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	776	PHE	2.4
1	B	495	THR	2.4
1	B	784	ILE	2.4
1	A	513	LEU	2.4
1	A	1125	VAL	2.4
1	A	1206	PHE	2.4
1	A	1028	TYR	2.3
1	A	1002	GLY	2.3
1	A	741	LEU	2.3
1	B	1001	LEU	2.3
1	A	793	VAL	2.3
1	B	432	ILE	2.3
1	A	564	TYR	2.3
1	B	403	GLY	2.3
1	B	631	TRP	2.3
1	A	1000	VAL	2.3
1	A	229	ILE	2.2
1	B	789	THR	2.2
1	B	673	GLY	2.2
1	A	494	VAL	2.2
1	A	1031	ALA	2.2
1	A	245	LEU	2.2
1	B	1044	LYS	2.2
1	A	981	LEU	2.2
1	A	1049	ASP	2.2
1	B	697	LEU	2.1
1	A	589	GLY	2.1
1	B	731	PRO	2.1
1	B	960	PRO	2.1
1	A	685	PHE	2.1
1	B	716	ALA	2.1
1	B	433	GLY	2.1
1	A	720	VAL	2.1
1	B	881	VAL	2.1
1	A	613	LEU	2.0
1	B	541	GLY	2.0
1	A	100	ARG	2.0
1	A	561	ASP	2.0
1	A	1032	ALA	2.0
1	B	345	ALA	2.0
1	A	992	LEU	2.0
1	A	623	HIS	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	777	GLY	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
2	SO4	B	1303	5/5	0.49	0.10	131,134,139,147	0
2	SO4	A	1307	5/5	0.67	0.08	121,129,135,139	0
2	SO4	A	1304	5/5	0.69	0.07	119,121,129,134	0
2	SO4	B	1304	5/5	0.70	0.08	132,133,141,145	0
2	SO4	A	1305	5/5	0.87	0.07	99,104,111,120	0
2	SO4	B	1302	5/5	0.89	0.09	91,94,97,98	0
2	SO4	A	1306	5/5	0.91	0.19	92,97,101,120	0
2	SO4	B	1301	5/5	0.95	0.06	65,68,73,75	0
2	SO4	A	1303	5/5	0.95	0.06	58,67,73,73	0
2	SO4	A	1302	5/5	0.96	0.08	67,69,74,75	0
2	SO4	A	1301	5/5	0.98	0.04	68,70,74,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.