



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 4FKB / pdb_00004fkb
Title : An Organic solvent tolerant lipase 42
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Deposited on : 2012-06-13
Resolution : 1.22 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

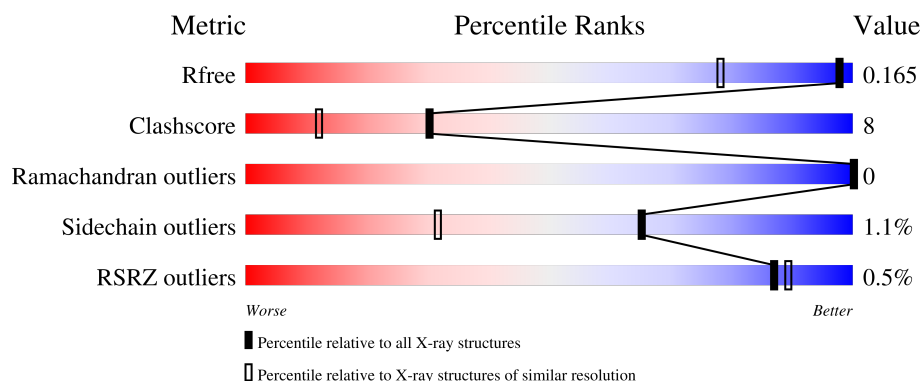
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.22 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	2002 (1.24-1.20)
Clashscore	190562	2061 (1.24-1.20)
Ramachandran outliers	187476	2009 (1.24-1.20)
Sidechain outliers	187428	2008 (1.24-1.20)
RSRZ outliers	180081	2000 (1.24-1.20)

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	388	 88% 11% .
1	B	388	 87% 11% .

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
3	GOL	A	413	-	X	X	-
3	GOL	B	417	-	-	X	-
3	GOL	B	418	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7622 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 Thermostable organic solvent tolerant lipase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	21	0
			3154	2010	549	586	9			
1	B	387	Total	C	N	O	S	0	39	0
			3286	2099	575	601	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	147	LEU	TRP	engineered mutation	UNP Q5U780
A	202	GLU	GLN	engineered mutation	UNP Q5U780
B	147	LEU	TRP	engineered mutation	UNP Q5U780
B	202	GLU	GLN	engineered mutation	UNP Q5U780

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 1 1	0	0
3	A	1	Total C O 12 6 6	0	1
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 5 3 2	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 18 9 9	0	1
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 12 6 6	0	1
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total Na 4 4	0	0
4	B	5	Total Na 5 5	0	0

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	5	Total Cl 5 5	0	0
5	B	5	Total Cl 5 5	0	0

- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total Ca 1 1	0	0
6	B	1	Total Ca 1 1	0	0

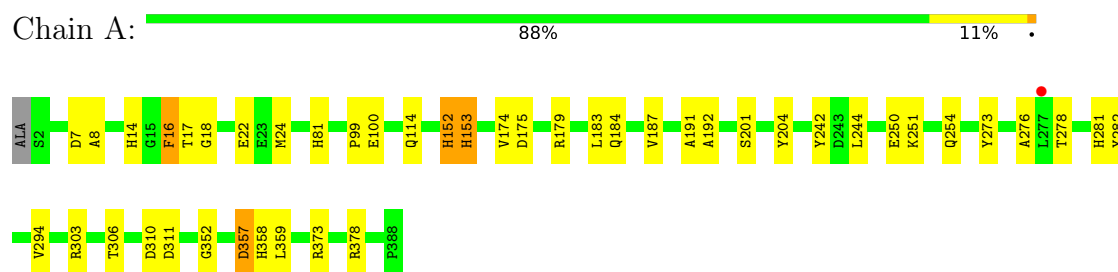
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	524	Total 524	O 524	0	0
7	B	491	Total 491	O 491	0	0

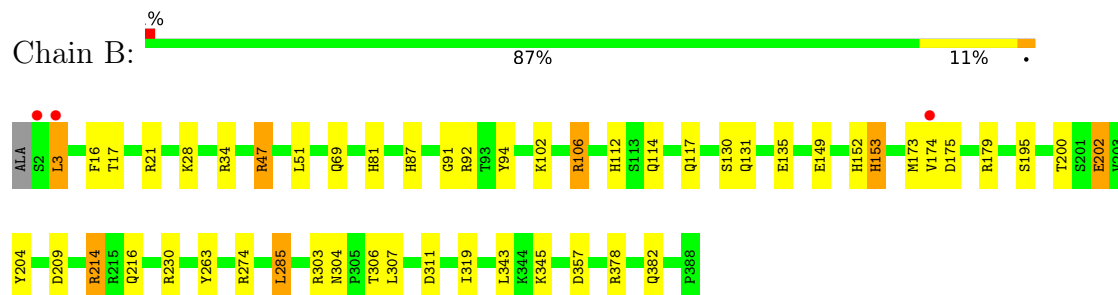
3 Residue-property plots

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: Thermostable organic solvent tolerant lipase



- Molecule 1: Thermostable organic solvent tolerant lipase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	117.82Å 81.00Å 99.53Å 90.00° 96.69° 90.00°	Depositor
Resolution (Å)	50.00 – 1.22 50.00 – 1.22	Depositor EDS
% Data completeness (in resolution range)	93.7 (50.00-1.22) 93.7 (50.00-1.22)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 1.22Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.127 , 0.165 0.127 , 0.165	Depositor DCC
R_{free} test set	13889 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	7.3	Xtriage
Anisotropy	0.136	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 53.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	7622	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.05% 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: CL, ZN, GOL, NA, CA

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	1.46	19/3298 (0.6%)	1.31	6/4483 (0.1%)
1	B	1.44	12/3475 (0.3%)	1.22	7/4711 (0.1%)
All	All	1.45	31/6773 (0.5%)	1.26	13/9194 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	303	ARG	CZ-NH1	10.04	1.46	1.32
1	A	303	ARG	CZ-NH2	8.05	1.44	1.33
1	B	91	GLY	C-O	6.83	1.31	1.24
1	B	274	ARG	CD-NE	-6.75	1.36	1.46
1	B	81	HIS	ND1-CE1	6.60	1.39	1.32
1	B	202	GLU	CD-OE2	5.96	1.36	1.25
1	B	230	ARG	N-CA	5.81	1.53	1.46
1	B	87	HIS	CE1-NE2	5.78	1.38	1.32
1	A	152	HIS	CA-C	-5.75	1.46	1.52
1	A	153	HIS	C-O	-5.74	1.17	1.24
1	A	7	ASP	CA-CB	5.71	1.61	1.53
1	B	202	GLU	N-CA	-5.63	1.39	1.46
1	B	304	ASN	C-O	-5.44	1.18	1.24
1	B	319	ILE	CA-C	5.39	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	152	HIS	ND1-CE1	5.38	1.38	1.32
1	A	152	HIS	N-CA	5.35	1.52	1.46
1	A	281	HIS	CD2-NE2	5.35	1.43	1.37
1	A	175	ASP	C-O	-5.25	1.17	1.24
1	A	175	ASP	N-CA	5.23	1.51	1.46
1	A	174	VAL	C-O	5.22	1.30	1.24
1	B	153	HIS	ND1-CE1	5.21	1.37	1.32
1	A	373	ARG	C-O	5.18	1.30	1.24
1	A	152	HIS	C-O	5.16	1.31	1.24
1	A	153	HIS	CD2-NE2	-5.15	1.32	1.37
1	A	358	HIS	ND1-CE1	5.15	1.37	1.32
1	A	153	HIS	CE1-NE2	5.14	1.37	1.32
1	A	359	LEU	N-CA	5.13	1.52	1.46
1	A	100	GLU	CD-OE1	5.09	1.35	1.25
1	A	81	HIS	ND1-CE1	5.09	1.37	1.32
1	B	112	HIS	CD2-NE2	5.04	1.43	1.37
1	A	352	GLY	N-CA	5.01	1.50	1.45

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	278	THR	CA-C-N	-11.18	104.96	122.51
1	A	278	THR	C-N-CA	-11.18	104.96	122.51
1	B	106	ARG	NE-CZ-NH1	-6.78	114.72	121.50
1	A	357	ASP	CA-CB-CG	5.83	118.43	112.60
1	B	303	ARG	NE-CZ-NH2	5.75	124.37	119.20
1	B	200	THR	CA-C-N	-5.59	109.58	121.32
1	B	200	THR	C-N-CA	-5.59	109.58	121.32
1	A	99	PRO	N-CA-C	-5.32	107.49	114.03
1	B	3	LEU	CA-CB-CG	-5.28	97.81	116.30
1	B	274	ARG	CB-CG-CD	5.26	123.39	111.30
1	B	47	ARG	NE-CZ-NH2	-5.06	114.64	119.20
1	A	8	ALA	CA-C-N	-5.01	115.07	120.03
1	A	8	ALA	C-N-CA	-5.01	115.07	120.03

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	16[A]	PHE	Peptide
1	B	16	PHE	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	3154	0	3068	40	0
1	B	3286	0	3258	61	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	54	0	68	15	0
3	B	90	0	120	21	0
4	A	4	0	0	0	0
4	B	5	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	1	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	524	0	0	15	2
7	B	491	0	0	17	3
All	All	7622	0	6514	110	3

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 8.

All (110) 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:16[B]:PHE:CE2	7:A:710:HOH:O	1.83	1.29
1:B:214[A]:ARG:HH21	1:B:214[A]:ARG:CG	1.35	1.29
1:B:174[B]:VAL:HG23	7:B:968:HOH:O	1.30	1.26
1:A:306[B]:THR:HG22	7:A:683:HOH:O	1.04	1.22
1:B:173[B]:MET:CE	1:B:179[B]:ARG:HD2	1.71	1.20
3:B:418:GOL:H31	7:B:627:HOH:O	1.41	1.20
1:B:131[B]:GLN:NE2	1:B:135[B]:GLU:OE2	1.74	1.19
1:B:174[B]:VAL:O	1:B:175[B]:ASP:HB2	1.45	1.12
1:B:378:ARG:HB3	3:B:417:GOL:O1	1.55	1.05
1:B:106:ARG:HH21	1:B:153:HIS:HE1	1.07	1.02
1:B:214[A]:ARG:NH2	1:B:214[A]:ARG:HG2	1.36	1.00
1:B:173[B]:MET:HE3	1:B:179[B]:ARG:CD	1.92	0.98
1:A:254[A]:GLN:NE2	7:A:935:HOH:O	1.96	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:HIS:HB3	3:A:412:GOL:H12	1.45	0.97
1:B:173[B]:MET:HE3	1:B:179[B]:ARG:HD2	0.97	0.96
1:B:149[B]:GLU:OE2	7:B:938:HOH:O	1.85	0.94
1:B:214[A]:ARG:NH2	1:B:216:GLN:HE22	1.66	0.94
1:B:378:ARG:HG3	7:B:804:HOH:O	1.76	0.86
1:A:242:TYR:OH	1:A:251:LYS:HD3	1.76	0.85
1:B:174[B]:VAL:O	1:B:175[B]:ASP:CB	2.20	0.85
1:B:131[B]:GLN:OE1	7:B:977:HOH:O	1.95	0.84
1:A:16[B]:PHE:CZ	7:A:710:HOH:O	2.17	0.82
1:B:378:ARG:HE	3:B:417:GOL:C1	1.92	0.81
1:B:214[A]:ARG:NH2	1:B:216:GLN:NE2	2.29	0.81
1:A:179:ARG:HD3	1:A:294[B]:VAL:HG21	1.67	0.77
1:A:114:GLN:NE2	1:A:244[B]:LEU:HG	2.00	0.76
1:B:214[A]:ARG:NH2	7:B:749:HOH:O	2.19	0.76
1:B:214[A]:ARG:CG	1:B:214[A]:ARG:NH2	2.09	0.74
1:B:378:ARG:HE	3:B:417:GOL:H11	1.51	0.73
1:B:106:ARG:HH21	1:B:153:HIS:CE1	1.99	0.71
1:A:250[A]:GLU:OE1	7:A:947:HOH:O	2.11	0.69
1:B:173[B]:MET:CE	1:B:179[B]:ARG:CD	2.58	0.69
1:B:195:SER:OG	3:B:418:GOL:H32	1.94	0.68
1:B:214[A]:ARG:HH21	1:B:214[A]:ARG:HG2	0.55	0.67
1:A:114:GLN:HE22	1:A:244[B]:LEU:HG	1.57	0.66
5:B:423:CL:CL	7:B:725:HOH:O	2.50	0.66
1:B:195:SER:OG	3:B:418:GOL:H12	1.96	0.65
1:B:214[A]:ARG:CZ	1:B:216:GLN:HE22	2.11	0.64
1:A:311[A]:ASP:H	3:A:413:GOL:H31	1.63	0.63
1:B:92[A]:ARG:NH2	7:B:801:HOH:O	2.31	0.63
1:A:153:HIS:CG	7:A:839:HOH:O	2.52	0.62
1:B:131[B]:GLN:NE2	1:B:135[B]:GLU:CD	2.56	0.62
1:B:378:ARG:CB	3:B:417:GOL:O1	2.38	0.61
1:B:106:ARG:NH2	1:B:153:HIS:HE1	1.89	0.61
3:A:413:GOL:H12	7:A:617:HOH:O	2.01	0.60
1:B:214[A]:ARG:HH22	1:B:216:GLN:NE2	1.99	0.60
1:B:378:ARG:NE	3:B:417:GOL:H11	2.17	0.60
1:B:131[B]:GLN:HE21	1:B:135[B]:GLU:CD	2.10	0.59
1:A:311[A]:ASP:OD2	3:A:413:GOL:H32	2.03	0.58
1:B:51:LEU:HD22	1:B:69[B]:GLN:CG	2.33	0.58
1:B:51:LEU:HD22	1:B:69[B]:GLN:HG3	1.86	0.58
1:B:378:ARG:HH21	3:B:417:GOL:H32	1.69	0.57
1:A:311[B]:ASP:H	3:A:413:GOL:H31	1.68	0.57
3:B:418:GOL:H32	7:B:608:HOH:O	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:SER:OG	3:B:418:GOL:C3	2.54	0.56
1:A:17[A]:THR:OG1	1:A:204:TYR:OH	2.23	0.56
1:A:310:ASP:CA	3:A:413:GOL:H31	2.36	0.55
1:A:378[A]:ARG:HG3	7:A:1023:HOH:O	2.07	0.55
1:A:378[B]:ARG:NH1	7:A:1010:HOH:O	1.81	0.54
3:B:407:GOL:C3	7:B:916:HOH:O	2.51	0.54
1:A:152:HIS:HD2	7:A:561:HOH:O	1.91	0.54
1:A:153:HIS:CD2	7:A:839:HOH:O	2.61	0.53
1:B:34:ARG:O	3:B:415:GOL:H31	2.09	0.53
1:A:184:GLN:O	1:A:187[B]:VAL:HG12	2.09	0.52
1:B:92[B]:ARG:HD2	1:B:94:TYR:OH	2.12	0.50
1:A:310:ASP:HB2	3:A:413:GOL:H31	1.93	0.50
1:B:263:TYR:CZ	1:B:345[B]:LYS:HE2	2.46	0.50
1:B:306:THR:HG23	7:B:747:HOH:O	2.12	0.50
1:A:276:ALA:O	7:A:996:HOH:O	2.20	0.49
1:B:378:ARG:HB2	3:B:417:GOL:H31	1.93	0.49
1:B:311:ASP:HB2	7:B:965:HOH:O	2.12	0.49
1:A:114:GLN:CD	1:A:244[B]:LEU:HG	2.37	0.49
1:B:378:ARG:HB3	3:B:417:GOL:HO1	1.74	0.48
1:A:152:HIS:HE1	7:A:845:HOH:O	1.95	0.48
1:B:102[A]:LYS:NZ	7:B:867:HOH:O	2.45	0.48
1:B:378:ARG:NH2	3:B:417:GOL:H32	2.28	0.48
1:A:311[A]:ASP:H	3:A:413:GOL:C3	2.24	0.48
1:B:47:ARG:HG2	7:B:824:HOH:O	2.13	0.48
1:B:28[B]:LYS:NZ	7:B:663:HOH:O	2.24	0.47
1:A:192:ALA:O	3:A:414:GOL:H2	2.15	0.47
1:A:14:HIS:CG	1:A:18:GLY:HA3	2.50	0.47
1:A:24[B]:MET:HG3	1:A:191:ALA:HB1	1.98	0.46
1:A:310:ASP:C	3:A:413:GOL:H31	2.40	0.46
1:B:130:SER:OG	3:B:416:GOL:H12	2.17	0.45
1:B:92[A]:ARG:HE	1:B:209[A]:ASP:CG	2.25	0.45
1:B:21:ARG:HD3	3:B:413[A]:GOL:H2	1.99	0.45
1:B:130:SER:OG	3:B:416:GOL:C1	2.65	0.45
1:B:3:LEU:H	1:B:3:LEU:HD12	1.81	0.44
1:A:311[B]:ASP:H	3:A:413:GOL:C3	2.28	0.44
1:A:273:TYR:CZ	1:A:282:TYR:HB2	2.53	0.44
1:A:201[A]:SER:HB2	3:A:411:GOL:HO2	1.82	0.44
1:A:201[B]:SER:OG	3:A:411:GOL:O2	1.81	0.43
3:B:418:GOL:C3	7:B:608:HOH:O	2.65	0.43
1:B:92[B]:ARG:HD2	1:B:94:TYR:CZ	2.53	0.43
1:A:251:LYS:HE2	1:A:251:LYS:HB3	1.65	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:TYR:CE1	1:B:345[B]:LYS:HE2	2.54	0.43
1:B:343:LEU:HD11	1:B:382[B]:GLN:HG2	1.99	0.43
1:A:24[A]:MET:CE	1:A:191:ALA:CB	2.97	0.42
1:A:153:HIS:HA	7:A:839:HOH:O	2.19	0.42
1:B:114[B]:GLN:HG3	1:B:117:GLN:OE1	2.19	0.42
1:B:17[B]:THR:OG1	1:B:204:TYR:OH	2.23	0.42
1:A:24[A]:MET:HE3	1:A:24[A]:MET:HB3	1.73	0.41
1:A:378[A]:ARG:NH2	7:A:776:HOH:O	2.54	0.41
1:B:214[A]:ARG:HH22	1:B:216:GLN:HE22	1.49	0.41
1:B:3:LEU:HD12	1:B:3:LEU:N	2.36	0.40
3:B:407:GOL:H32	7:B:916:HOH:O	2.19	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:683:HOH:O	7:B:718:HOH:O[4_545]	2.03	0.17
7:B:736:HOH:O	7:B:761:HOH:O[4_545]	2.09	0.11
7:A:684:HOH:O	7:B:682:HOH:O[3_445]	2.16	0.04

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	406/388 (105%)	397 (98%)	9 (2%)	0	100	100
1	B	425/388 (110%)	414 (97%)	11 (3%)	0	100	100
All	All	831/776 (107%)	811 (98%)	20 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	337/318 (106%)	335 (99%)	2 (1%)	78	52
1	B	357/318 (112%)	350 (98%)	7 (2%)	48	12
All	All	694/636 (109%)	685 (99%)	9 (1%)	65	25

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	357	ASP
1	B	202	GLU
1	B	214[A]	ARG
1	B	214[B]	ARG
1	B	285[A]	LEU
1	B	285[B]	LEU
1	B	307	LEU
1	B	357	ASP

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	59	ASN
1	A	152	HIS
1	B	44	ASN
1	B	128	ASN
1	B	153	HIS
1	B	216	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 48 ligands modelled in this entry, 23 are monoatomic and 1 is modelled with single atom - leaving 24 for Mogul analysis.

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	GOL	A	410	-	5,5,5	0.83	0	5,5,5	1.36	1 (20%)
3	GOL	B	413[A]	-	5,5,5	1.22	0	5,5,5	1.28	1 (20%)
3	GOL	B	417	-	5,5,5	0.79	0	5,5,5	0.68	0
3	GOL	B	410	-	5,5,5	0.99	0	5,5,5	0.66	0
3	GOL	A	408	-	5,5,5	0.22	0	5,5,5	0.83	0
3	GOL	A	403[A]	-	5,5,5	0.71	0	5,5,5	0.89	0
3	GOL	A	412	-	5,5,5	0.78	0	5,5,5	1.21	1 (20%)
3	GOL	A	414	-	5,5,5	0.68	0	5,5,5	0.87	0
3	GOL	B	407	-	5,5,5	0.92	0	5,5,5	0.98	0
3	GOL	B	411	-	5,5,5	1.24	1 (20%)	5,5,5	0.88	0
3	GOL	B	408[B]	-	5,5,5	0.53	0	5,5,5	0.95	0
3	GOL	B	414	-	5,5,5	0.33	0	5,5,5	0.68	0
3	GOL	B	408[C]	-	5,5,5	0.66	0	5,5,5	0.68	0
3	GOL	B	416	-	5,5,5	0.83	0	5,5,5	1.26	0
3	GOL	B	413[B]	-	5,5,5	1.15	1 (20%)	5,5,5	1.47	1 (20%)
3	GOL	A	411	-	4,4,5	0.56	0	4,4,5	1.93	1 (25%)
3	GOL	A	403[B]	-	5,5,5	0.46	0	5,5,5	1.16	0
3	GOL	A	409	-	5,5,5	1.01	0	5,5,5	1.17	1 (20%)
3	GOL	B	418	-	5,5,5	0.67	0	5,5,5	1.67	1 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	A	413	-	5,5,5	1.06	0	5,5,5	2.83	2 (40%)
3	GOL	B	408[A]	-	5,5,5	0.71	0	5,5,5	0.93	0
3	GOL	B	409	-	5,5,5	0.56	0	5,5,5	1.91	3 (60%)
3	GOL	B	412	-	5,5,5	0.67	0	5,5,5	1.02	0
3	GOL	B	415	-	5,5,5	0.82	0	5,5,5	1.01	0

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
3	GOL	A	410	-	-	0/4/4/4	-
3	GOL	B	413[A]	-	-	0/4/4/4	-
3	GOL	B	417	-	-	0/4/4/4	-
3	GOL	B	410	-	-	0/4/4/4	-
3	GOL	A	408	-	-	4/4/4/4	-
3	GOL	A	403[A]	-	-	3/4/4/4	-
3	GOL	A	412	-	-	2/4/4/4	-
3	GOL	A	414	-	-	4/4/4/4	-
3	GOL	B	407	-	-	0/4/4/4	-
3	GOL	B	411	-	-	0/4/4/4	-
3	GOL	B	408[B]	-	-	0/4/4/4	-
3	GOL	B	414	-	-	1/4/4/4	-
3	GOL	B	408[C]	-	-	1/4/4/4	-
3	GOL	B	416	-	-	2/4/4/4	-
3	GOL	B	413[B]	-	-	0/4/4/4	-
3	GOL	A	411	-	-	2/2/2/4	-
3	GOL	A	403[B]	-	-	2/4/4/4	-
3	GOL	A	409	-	-	0/4/4/4	-
3	GOL	B	418	-	-	4/4/4/4	-
3	GOL	A	413	-	-	4/4/4/4	-
3	GOL	B	408[A]	-	-	0/4/4/4	-
3	GOL	B	409	-	-	1/4/4/4	-
3	GOL	B	412	-	-	1/4/4/4	-
3	GOL	B	415	-	-	2/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	411	GOL	O3-C3	2.45	1.52	1.42
3	B	413[B]	GOL	O2-C2	-2.12	1.37	1.43

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	413	GOL	O2-C2-C3	4.11	126.21	109.18
3	A	413	GOL	O3-C3-C2	4.01	128.45	110.38
3	A	411	GOL	O2-C2-C3	-3.62	97.25	114.99
3	B	413[B]	GOL	O2-C2-C3	-2.50	98.83	109.18
3	A	412	GOL	O2-C2-C1	-2.44	99.09	109.18
3	B	409	GOL	C3-C2-C1	-2.31	103.31	111.80
3	B	418	GOL	C3-C2-C1	-2.30	103.36	111.80
3	B	409	GOL	O1-C1-C2	-2.27	100.16	110.38
3	B	409	GOL	O2-C2-C3	-2.22	99.97	109.18
3	A	409	GOL	C3-C2-C1	-2.09	104.11	111.80
3	B	413[A]	GOL	O2-C2-C3	-2.07	100.59	109.18
3	A	410	GOL	O1-C1-C2	-2.00	101.37	110.38

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	412	GOL	C1-C2-C3-O3
3	A	412	GOL	O2-C2-C3-O3
3	A	414	GOL	O1-C1-C2-C3
3	B	416	GOL	C1-C2-C3-O3
3	B	416	GOL	O2-C2-C3-O3
3	B	418	GOL	C1-C2-C3-O3
3	A	414	GOL	O1-C1-C2-O2
3	B	418	GOL	O1-C1-C2-O2
3	A	403[A]	GOL	O1-C1-C2-C3
3	A	403[B]	GOL	O1-C1-C2-C3
3	A	408	GOL	O1-C1-C2-C3
3	A	408	GOL	C1-C2-C3-O3
3	A	413	GOL	O1-C1-C2-C3
3	A	413	GOL	C1-C2-C3-O3
3	A	414	GOL	C1-C2-C3-O3
3	B	412	GOL	O1-C1-C2-C3
3	B	415	GOL	C1-C2-C3-O3
3	B	418	GOL	O1-C1-C2-C3
3	A	403[A]	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	A	408	GOL	O1-C1-C2-O2
3	A	413	GOL	O1-C1-C2-O2
3	A	413	GOL	O2-C2-C3-O3
3	B	418	GOL	O2-C2-C3-O3
3	A	414	GOL	O2-C2-C3-O3
3	A	411	GOL	O2-C2-C3-O3
3	A	408	GOL	O2-C2-C3-O3
3	B	409	GOL	O2-C2-C3-O3
3	A	411	GOL	C1-C2-C3-O3
3	A	403[A]	GOL	O2-C2-C3-O3
3	B	415	GOL	O2-C2-C3-O3
3	A	403[B]	GOL	O1-C1-C2-O2
3	B	408[C]	GOL	O1-C1-C2-C3
3	B	414	GOL	O1-C1-C2-C3

There are no ring outliers.

10 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	413[A]	GOL	1	0
3	B	417	GOL	9	0
3	A	412	GOL	1	0
3	A	414	GOL	1	0
3	B	407	GOL	2	0
3	B	416	GOL	2	0
3	A	411	GOL	2	0
3	B	418	GOL	6	0
3	A	413	GOL	11	0
3	B	415	GOL	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	387/388 (99%)	-0.75	1 (0%) 90 92	5, 11, 23, 33	21 (5%)
1	B	387/388 (99%)	-0.85	3 (0%) 82 85	4, 9, 22, 49	39 (10%)
All	All	774/776 (99%)	-0.80	4 (0%) 87 89	4, 10, 22, 49	60 (7%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	174[A]	VAL	5.0
1	B	2	SER	2.7
1	B	3	LEU	2.5
1	A	277	LEU	2.1

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(Å ²)	Q<0.9
4	NA	A	406	1/1	0.82	0.16	64,64,64,64	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	A	403[B]	6/6	0.85	0.15	21,25,27,34	6
3	GOL	A	403[A]	6/6	0.85	0.15	21,34,37,39	6
3	GOL	B	409	6/6	0.87	0.13	34,39,49,60	0
3	GOL	A	414	6/6	0.87	0.16	43,51,82,94	0
3	GOL	B	417	6/6	0.89	0.13	20,31,43,51	0
3	GOL	B	416	6/6	0.89	0.13	20,35,50,56	0
3	GOL	B	418	6/6	0.91	0.10	24,35,44,46	0
3	GOL	A	410	6/6	0.92	0.09	31,35,38,41	0
3	GOL	A	411	5/6	0.92	0.10	30,33,43,44	0
3	GOL	B	414	6/6	0.92	0.10	44,50,60,71	0
3	GOL	B	415	6/6	0.92	0.10	27,33,35,53	0
3	GOL	A	408	6/6	0.94	0.09	41,43,57,72	0
3	GOL	A	412	6/6	0.95	0.12	19,42,51,61	0
3	GOL	B	407	6/6	0.95	0.07	16,22,25,28	0
3	GOL	A	413	6/6	0.95	0.08	16,22,32,44	0
4	NA	A	407	1/1	0.95	0.38	37,37,37,37	0
4	NA	B	405	1/1	0.95	0.17	42,42,42,42	0
3	GOL	B	412	6/6	0.96	0.07	15,21,25,25	0
3	GOL	B	410	6/6	0.97	0.07	14,17,24,32	0
3	GOL	B	408[B]	6/6	0.97	0.06	8,12,13,14	6
3	GOL	B	413[A]	6/6	0.97	0.06	13,17,21,23	6
3	GOL	B	413[B]	6/6	0.97	0.06	8,17,24,30	6
3	GOL	B	408[C]	6/6	0.97	0.06	6,14,18,18	6
4	NA	B	403	1/1	0.97	0.11	40,40,40,40	0
4	NA	B	404	1/1	0.97	0.09	31,31,31,31	0
3	GOL	B	408[A]	6/6	0.97	0.06	7,14,18,24	6
4	NA	A	404	1/1	0.98	0.11	26,26,26,26	0
4	NA	B	402	1/1	0.99	0.07	26,26,26,26	0
4	NA	A	405	1/1	0.99	0.10	25,25,25,25	0
3	GOL	B	411	6/6	0.99	0.04	11,14,20,23	0
3	GOL	A	409	6/6	0.99	0.04	12,14,15,19	0
5	CL	A	415	1/1	0.99	0.13	20,20,20,20	0
2	ZN	A	401	1/1	1.00	0.01	7,7,7,7	0
2	ZN	B	401	1/1	1.00	0.01	6,6,6,6	0
4	NA	B	406	1/1	1.00	0.01	12,12,12,12	0
3	GOL	A	402	1/6	1.00	0.07	15,15,15,15	0
5	CL	A	416	1/1	1.00	0.02	15,15,15,15	1
5	CL	A	417	1/1	1.00	0.06	14,14,14,14	0
5	CL	A	418	1/1	1.00	0.01	18,18,18,18	0
5	CL	A	419	1/1	1.00	0.07	18,18,18,18	0
5	CL	B	419	1/1	1.00	0.05	12,12,12,12	0
5	CL	B	420	1/1	1.00	0.03	14,14,14,14	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	CL	B	421	1/1	1.00	0.01	13,13,13,13	0
5	CL	B	422	1/1	1.00	0.05	15,15,15,15	0
5	CL	B	423	1/1	1.00	0.12	16,16,16,16	0
6	CA	A	420	1/1	1.00	0.03	8,8,8,8	1
6	CA	B	424	1/1	1.00	0.03	7,7,7,7	1

6.5 Other polymers [i](#)

There are no such residues in this entry.