



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:15 PM UTC

PDB ID : 3KTS / pdb_00003kts
Title : CRYSTAL STRUCTURE OF GLYCEROL UPTAKE OPERON ANTITERMINATOR REGULATORY PROTEIN FROM LISTERIA MONOCYTOGENES STR. 4b F2365
Authors : Patskovsky, Y.; Toro, R.; Freeman, J.; Do, J.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYS-GXRC)
Deposited on : 2009-11-25
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:

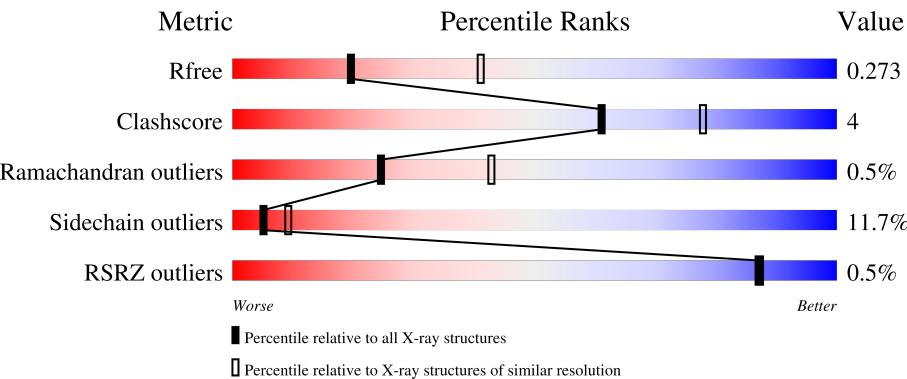
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 2.75 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	192	% 77% 17% ..
1	B	192	2% 81% 14% ..
1	C	192	% 79% 11% 5% 5%
1	D	192	76% 17% .. 5%

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Mol	Chain	Length	Quality of chain
1	E	192	81%15%
1	F	192	% 73%20%
1	G	192	79%16%
1	H	192	% 79%16%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11436 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 Glycerol uptake operon antiterminator regulatory protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	185	Total	C	N	O	S	0	0	0
			1425	909	242	264	10			
1	B	185	Total	C	N	O	S	0	0	0
			1425	909	242	264	10			
1	C	182	Total	C	N	O	S	0	0	0
			1403	895	237	261	10			
1	D	183	Total	C	N	O	S	0	0	0
			1411	901	238	262	10			
1	E	185	Total	C	N	O	S	0	0	0
			1425	909	242	264	10			
1	F	184	Total	C	N	O	S	0	0	0
			1415	903	239	263	10			
1	G	185	Total	C	N	O	S	0	0	0
			1425	909	242	264	10			
1	H	185	Total	C	N	O	S	0	0	0
			1425	909	242	264	10			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q71ZR5
A	1	LEU	-	expression tag	UNP Q71ZR5
A	183	GLU	-	expression tag	UNP Q71ZR5
A	184	GLY	-	expression tag	UNP Q71ZR5
A	185	HIS	-	expression tag	UNP Q71ZR5
A	186	HIS	-	expression tag	UNP Q71ZR5
A	187	HIS	-	expression tag	UNP Q71ZR5
A	188	HIS	-	expression tag	UNP Q71ZR5
A	189	HIS	-	expression tag	UNP Q71ZR5
A	190	HIS	-	expression tag	UNP Q71ZR5
B	0	SER	-	expression tag	UNP Q71ZR5
B	1	LEU	-	expression tag	UNP Q71ZR5
B	183	GLU	-	expression tag	UNP Q71ZR5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	184	GLY	-	expression tag	UNP Q71ZR5
B	185	HIS	-	expression tag	UNP Q71ZR5
B	186	HIS	-	expression tag	UNP Q71ZR5
B	187	HIS	-	expression tag	UNP Q71ZR5
B	188	HIS	-	expression tag	UNP Q71ZR5
B	189	HIS	-	expression tag	UNP Q71ZR5
B	190	HIS	-	expression tag	UNP Q71ZR5
C	0	SER	-	expression tag	UNP Q71ZR5
C	1	LEU	-	expression tag	UNP Q71ZR5
C	183	GLU	-	expression tag	UNP Q71ZR5
C	184	GLY	-	expression tag	UNP Q71ZR5
C	185	HIS	-	expression tag	UNP Q71ZR5
C	186	HIS	-	expression tag	UNP Q71ZR5
C	187	HIS	-	expression tag	UNP Q71ZR5
C	188	HIS	-	expression tag	UNP Q71ZR5
C	189	HIS	-	expression tag	UNP Q71ZR5
C	190	HIS	-	expression tag	UNP Q71ZR5
D	0	SER	-	expression tag	UNP Q71ZR5
D	1	LEU	-	expression tag	UNP Q71ZR5
D	183	GLU	-	expression tag	UNP Q71ZR5
D	184	GLY	-	expression tag	UNP Q71ZR5
D	185	HIS	-	expression tag	UNP Q71ZR5
D	186	HIS	-	expression tag	UNP Q71ZR5
D	187	HIS	-	expression tag	UNP Q71ZR5
D	188	HIS	-	expression tag	UNP Q71ZR5
D	189	HIS	-	expression tag	UNP Q71ZR5
D	190	HIS	-	expression tag	UNP Q71ZR5
E	0	SER	-	expression tag	UNP Q71ZR5
E	1	LEU	-	expression tag	UNP Q71ZR5
E	183	GLU	-	expression tag	UNP Q71ZR5
E	184	GLY	-	expression tag	UNP Q71ZR5
E	185	HIS	-	expression tag	UNP Q71ZR5
E	186	HIS	-	expression tag	UNP Q71ZR5
E	187	HIS	-	expression tag	UNP Q71ZR5
E	188	HIS	-	expression tag	UNP Q71ZR5
E	189	HIS	-	expression tag	UNP Q71ZR5
E	190	HIS	-	expression tag	UNP Q71ZR5
F	0	SER	-	expression tag	UNP Q71ZR5
F	1	LEU	-	expression tag	UNP Q71ZR5
F	183	GLU	-	expression tag	UNP Q71ZR5
F	184	GLY	-	expression tag	UNP Q71ZR5
F	185	HIS	-	expression tag	UNP Q71ZR5

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Chain	Residue	Modelled	Actual	Comment	Reference
F	186	HIS	-	expression tag	UNP Q71ZR5
F	187	HIS	-	expression tag	UNP Q71ZR5
F	188	HIS	-	expression tag	UNP Q71ZR5
F	189	HIS	-	expression tag	UNP Q71ZR5
F	190	HIS	-	expression tag	UNP Q71ZR5
G	0	SER	-	expression tag	UNP Q71ZR5
G	1	LEU	-	expression tag	UNP Q71ZR5
G	183	GLU	-	expression tag	UNP Q71ZR5
G	184	GLY	-	expression tag	UNP Q71ZR5
G	185	HIS	-	expression tag	UNP Q71ZR5
G	186	HIS	-	expression tag	UNP Q71ZR5
G	187	HIS	-	expression tag	UNP Q71ZR5
G	188	HIS	-	expression tag	UNP Q71ZR5
G	189	HIS	-	expression tag	UNP Q71ZR5
G	190	HIS	-	expression tag	UNP Q71ZR5
H	0	SER	-	expression tag	UNP Q71ZR5
H	1	LEU	-	expression tag	UNP Q71ZR5
H	183	GLU	-	expression tag	UNP Q71ZR5
H	184	GLY	-	expression tag	UNP Q71ZR5
H	185	HIS	-	expression tag	UNP Q71ZR5
H	186	HIS	-	expression tag	UNP Q71ZR5
H	187	HIS	-	expression tag	UNP Q71ZR5
H	188	HIS	-	expression tag	UNP Q71ZR5
H	189	HIS	-	expression tag	UNP Q71ZR5
H	190	HIS	-	expression tag	UNP Q71ZR5

- Molecule 2 is UNKNOWN LIGAND (CCD ID: UNL) (formula:).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			10	3	6	1		
2	B	1	Total	C	O	P	0	0
			10	3	6	1		
2	C	1	Total	C	O	P	0	0
			10	3	6	1		
2	D	1	Total	C	O	P	0	0
			10	3	6	1		
2	E	1	Total	C	O	P	0	0
			10	3	6	1		
2	F	1	Total	C	O	P	0	0
			10	3	6	1		
2	G	1	Total	C	O	P	0	0
			10	3	6	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	H	1	Total	C	O	P	0	0
			10	3	6	1		

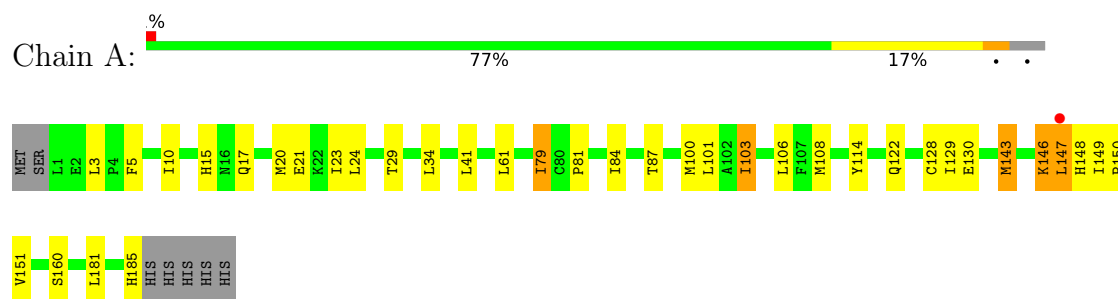
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	2	Total	O	0	0
			2	2		

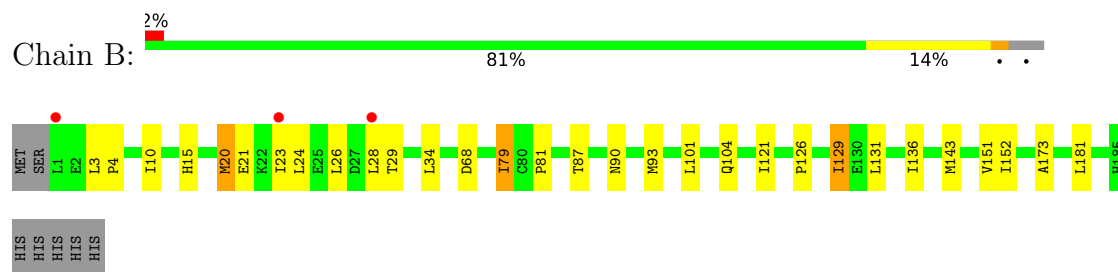
3 Residue-property plots

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.

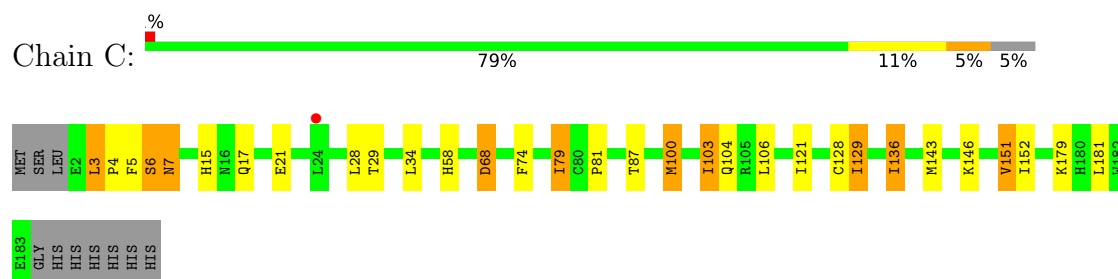
- Molecule 1: Glycerol uptake operon antiterminator regulatory protein



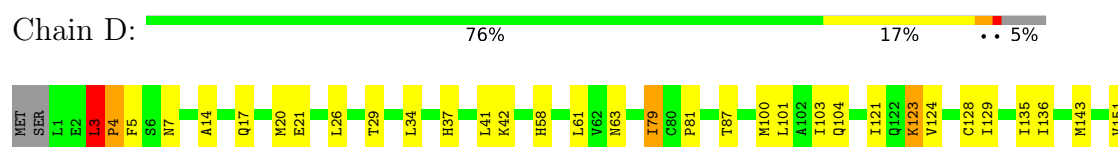
- Molecule 1: Glycerol uptake operon antiterminator regulatory protein



- Molecule 1: Glycerol uptake operon antiterminator regulatory protein



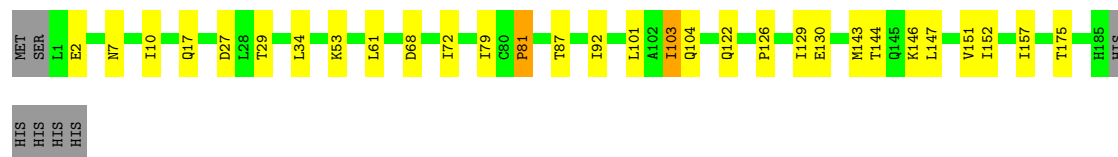
- Molecule 1: Glycerol uptake operon antiterminator regulatory protein





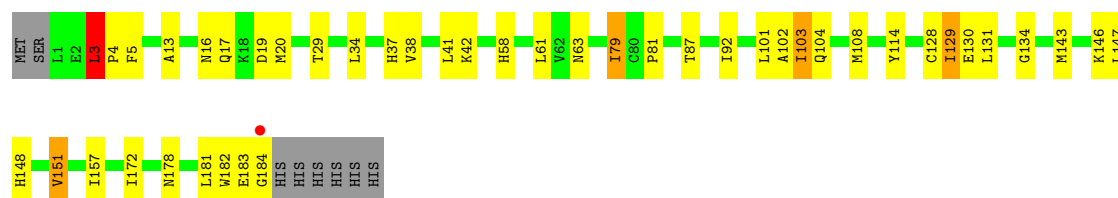
- Molecule 1: Glycerol uptake operon antiterminator regulatory protein

Chain E: 81% 15% . .



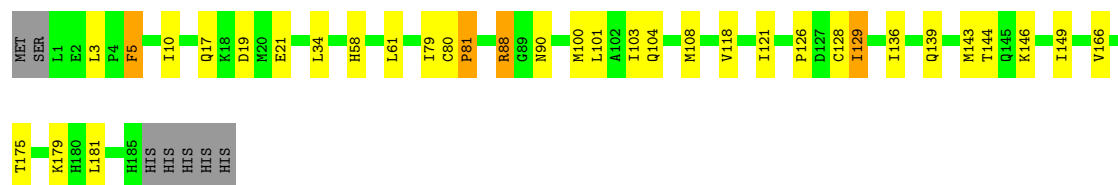
- Molecule 1: Glycerol uptake operon antiterminator regulatory protein

Chain F: 73% 20% . . .



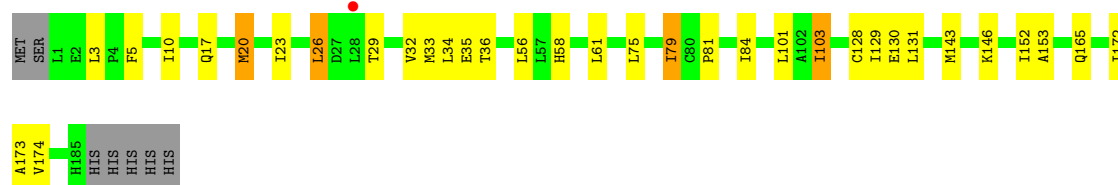
- Molecule 1: Glycerol uptake operon antiterminator regulatory protein

Chain G: 79% 16% . .



- Molecule 1: Glycerol uptake operon antiterminator regulatory protein

Chain H: 79% 16% . .



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	137.60Å 110.60Å 130.97Å 90.00° 116.44° 90.00°	Depositor
Resolution (Å)	20.00 – 2.75 20.00 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.0 (20.00-2.75) 98.7 (20.00-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.64	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.3.0034	Depositor
R, R_{free}	0.218 , 0.272 0.222 , 0.273	Depositor DCC
R_{free} test set	1408 reflections (3.08%)	wwPDB-VP
Wilson B-factor (Å ²)	92.1	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 62.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11436	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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: UNL

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	0.48	0/1447	0.93	1/1957 (0.1%)
1	B	0.50	0/1447	0.97	2/1957 (0.1%)
1	C	0.50	0/1424	0.93	0/1926
1	D	0.49	0/1432	0.97	3/1937 (0.2%)
1	E	0.49	0/1447	0.89	0/1957
1	F	0.50	0/1436	0.95	2/1942 (0.1%)
1	G	0.46	0/1447	0.84	0/1957
1	H	0.47	0/1447	0.90	2/1957 (0.1%)
All	All	0.49	0/11527	0.92	10/15590 (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	2
1	C	0	3
All	All	0	5

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	3	LEU	CA-C-N	9.14	131.26	119.84
1	B	3	LEU	C-N-CA	9.14	131.26	119.84
1	F	3	LEU	CA-C-N	8.88	130.94	119.84
1	F	3	LEU	C-N-CA	8.88	130.94	119.84
1	D	3	LEU	CA-C-N	8.70	130.71	119.84
1	D	3	LEU	C-N-CA	8.70	130.71	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	3	LEU	CA-C-N	6.49	127.95	119.84
1	H	3	LEU	C-N-CA	6.49	127.95	119.84
1	A	41	LEU	N-CA-C	5.59	117.05	111.07
1	D	41	LEU	N-CA-C	5.39	116.84	111.07

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	147	LEU	Peptide
1	A	148	HIS	Peptide
1	C	3	LEU	Peptide
1	C	5	PHE	Peptide
1	C	7	ASN	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	1425	0	1482	13	0
1	B	1425	0	1482	9	0
1	C	1403	0	1458	10	0
1	D	1411	0	1472	13	0
1	E	1425	0	1482	8	0
1	F	1415	0	1475	20	0
1	G	1425	0	1482	16	0
1	H	1425	0	1482	11	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
2	E	10	0	0	0	0
2	F	10	0	0	0	0
2	G	10	0	0	0	0
2	H	10	0	0	0	0
3	G	2	0	0	0	0
All	All	11436	0	11815	98	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 4.

All (98) 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:149:ILE:HG23	1:A:150:PRO:HD2	1.27	1.09
1:G:129:ILE:HG22	1:G:149:ILE:CG2	1.83	1.09
1:G:129:ILE:HG22	1:G:149:ILE:HG22	1.39	1.05
1:A:149:ILE:CG2	1:A:150:PRO:HD2	1.94	0.98
1:F:181:LEU:O	1:F:181:LEU:HD12	1.67	0.93
1:G:129:ILE:HG22	1:G:149:ILE:HG21	1.63	0.80
1:A:103:ILE:HB	1:A:128:CYS:HB3	1.70	0.74
1:H:34:LEU:O	1:H:61:LEU:HB2	1.87	0.73
1:C:79:ILE:HG22	1:C:81:PRO:HD3	1.73	0.71
1:D:79:ILE:HG22	1:D:81:PRO:HD3	1.76	0.67
1:A:79:ILE:HG22	1:A:81:PRO:HD3	1.76	0.67
1:G:129:ILE:CG2	1:G:149:ILE:CG2	2.70	0.66
1:C:103:ILE:HB	1:C:128:CYS:HB3	1.78	0.65
1:D:7:ASN:O	1:D:7:ASN:ND2	2.31	0.64
1:G:3:LEU:HD13	1:G:5:PHE:CE1	2.35	0.62
1:A:149:ILE:CG2	1:A:150:PRO:CD	2.73	0.61
1:E:79:ILE:HG22	1:E:81:PRO:HD3	1.81	0.61
1:B:79:ILE:HG22	1:B:81:PRO:HD3	1.82	0.60
1:G:34:LEU:O	1:G:61:LEU:HB2	2.02	0.59
1:F:34:LEU:O	1:F:61:LEU:HB2	2.02	0.59
1:H:103:ILE:HB	1:H:128:CYS:HB3	1.86	0.57
1:D:34:LEU:O	1:D:61:LEU:HB2	2.05	0.57
1:G:3:LEU:HD13	1:G:5:PHE:HE1	1.69	0.56
1:A:103:ILE:HD11	1:A:130:GLU:HB2	1.88	0.56
1:F:181:LEU:O	1:F:181:LEU:CD1	2.50	0.55
1:C:15:HIS:CE1	1:C:34:LEU:HD13	2.40	0.55
1:A:34:LEU:O	1:A:61:LEU:HB2	2.07	0.55
1:G:129:ILE:CG2	1:G:149:ILE:HG21	2.34	0.55
1:F:103:ILE:HB	1:F:128:CYS:HB3	1.88	0.55
1:F:129:ILE:HG13	1:F:151:VAL:HG12	1.90	0.54
1:E:34:LEU:O	1:E:61:LEU:HB2	2.08	0.53
1:B:20:MET:HA	1:B:23:ILE:HG12	1.91	0.52
1:D:14:ALA:O	1:D:34:LEU:HB2	2.10	0.52
1:F:16:ASN:H	1:F:19:ASP:HB2	1.75	0.52
1:H:79:ILE:HG22	1:H:81:PRO:HD3	1.91	0.52
1:F:183:GLU:HG2	1:F:184:GLY:H	1.75	0.51
1:E:103:ILE:HD11	1:E:130:GLU:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:104:GLN:HE22	1:G:121:ILE:HA	1.76	0.51
1:D:81:PRO:O	1:D:100:MET:HE1	2.12	0.50
1:C:81:PRO:HD2	1:C:100:MET:HE1	1.93	0.50
1:H:103:ILE:HD11	1:H:130:GLU:HB2	1.93	0.50
1:G:79:ILE:HG22	1:G:81:PRO:HD3	1.92	0.50
1:A:122:GLN:OE1	1:A:147:LEU:HD21	2.12	0.49
1:D:104:GLN:HE22	1:D:121:ILE:HA	1.77	0.48
1:H:153:ALA:HB3	1:H:174:VAL:HG22	1.96	0.48
1:F:92:ILE:HG13	1:F:104:GLN:HG3	1.96	0.48
1:D:152:ILE:HD12	1:D:173:ALA:HB3	1.96	0.48
1:C:129:ILE:HG13	1:C:151:VAL:HG12	1.95	0.48
1:A:108:MET:HE1	1:A:114:TYR:HD1	1.79	0.47
1:F:13:ALA:HB1	1:F:34:LEU:HD11	1.97	0.47
1:B:26:LEU:HD23	1:B:28:LEU:HD12	1.97	0.47
1:G:118:VAL:HA	1:G:121:ILE:HD12	1.97	0.47
1:E:104:GLN:HB3	1:E:129:ILE:HG22	1.97	0.47
1:F:79:ILE:HG22	1:F:81:PRO:HD3	1.97	0.46
1:F:134:GLY:HA3	1:F:157:ILE:HD11	1.98	0.46
1:A:34:LEU:O	1:A:61:LEU:CB	2.64	0.46
1:H:23:ILE:HA	1:H:26:LEU:HD13	1.98	0.46
1:G:103:ILE:HG12	1:G:128:CYS:HB3	1.98	0.45
1:A:15:HIS:CE1	1:A:34:LEU:HD13	2.51	0.45
1:E:104:GLN:HE21	1:E:126:PRO:HG3	1.81	0.45
1:E:152:ILE:HG21	1:E:175:THR:HG23	1.99	0.45
1:B:152:ILE:HD12	1:B:173:ALA:HB3	1.98	0.45
1:A:143:MET:O	1:A:146:LYS:HB2	2.17	0.45
1:B:104:GLN:HE22	1:B:121:ILE:HA	1.83	0.44
1:D:103:ILE:HG12	1:D:128:CYS:HB3	1.99	0.44
1:H:20:MET:HA	1:H:23:ILE:HG12	1.99	0.44
1:B:90:ASN:HA	1:B:93:MET:HE2	1.99	0.43
1:B:129:ILE:HG13	1:B:151:VAL:HG12	1.99	0.43
1:E:68:ASP:O	1:E:72:ILE:HG12	2.18	0.43
1:F:3:LEU:HD11	1:F:128:CYS:HB2	1.99	0.43
1:C:74:PHE:HB2	1:F:38:VAL:HG13	1.99	0.43
1:C:104:GLN:HE22	1:C:121:ILE:HA	1.84	0.43
1:G:126:PRO:O	1:G:149:ILE:HD12	2.19	0.43
1:D:26:LEU:HD21	1:D:182:TRP:HB3	2.00	0.42
1:F:103:ILE:HD11	1:F:130:GLU:HB2	2.01	0.42
1:G:108:MET:HE2	1:G:139:GLN:HB3	2.01	0.42
1:D:37:HIS:HD2	1:D:63:ASN:HD22	1.66	0.42
1:B:15:HIS:CE1	1:B:34:LEU:HD13	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:6:SER:O	1:C:7:ASN:HB2	2.18	0.42
1:H:34:LEU:O	1:H:61:LEU:CB	2.62	0.42
1:F:92:ILE:HG23	1:F:102:ALA:HB1	2.02	0.42
1:F:37:HIS:CD2	1:F:63:ASN:HD22	2.38	0.41
1:F:37:HIS:HD2	1:F:63:ASN:HD22	1.69	0.41
1:G:81:PRO:O	1:G:100:MET:HE1	2.21	0.41
1:E:92:ILE:HG13	1:E:104:GLN:HG3	2.03	0.41
1:H:32:VAL:HG22	1:H:56:LEU:HB2	2.03	0.41
1:F:41:LEU:HD23	1:F:79:ILE:HG12	2.03	0.41
1:H:75:LEU:HA	1:H:79:ILE:HB	2.03	0.41
1:C:3:LEU:HA	1:C:4:PRO:HD3	1.42	0.41
1:D:123:LYS:HG2	1:D:124:VAL:HG13	2.02	0.41
1:H:152:ILE:HD12	1:H:173:ALA:HB3	2.02	0.41
1:A:20:MET:HA	1:A:23:ILE:HG12	2.03	0.41
1:B:104:GLN:HE21	1:B:126:PRO:HG3	1.86	0.41
1:F:178:ASN:HB3	1:F:181:LEU:CD2	2.51	0.41
1:C:136:ILE:HD13	1:D:135:ILE:HD11	2.03	0.40
1:F:108:MET:HE1	1:F:114:TYR:CD1	2.56	0.40
1:G:88:ARG:HH21	1:G:90:ASN:HB2	1.86	0.40
1:D:3:LEU:HA	1:D:4:PRO:HD2	1.75	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	183/192 (95%)	177 (97%)	6 (3%)	0	100	100
1	B	183/192 (95%)	169 (92%)	12 (7%)	2 (1%)	11	22
1	C	180/192 (94%)	171 (95%)	8 (4%)	1 (1%)	21	38
1	D	181/192 (94%)	171 (94%)	9 (5%)	1 (1%)	21	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	183/192 (95%)	172 (94%)	10 (6%)	1 (0%)	24	43
1	F	182/192 (95%)	168 (92%)	13 (7%)	1 (0%)	24	43
1	G	183/192 (95%)	174 (95%)	8 (4%)	1 (0%)	24	43
1	H	183/192 (95%)	173 (94%)	10 (6%)	0	100	100
All	All	1458/1536 (95%)	1375 (94%)	76 (5%)	7 (0%)	24	43

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	4	PRO
1	B	68	ASP
1	D	4	PRO
1	G	81	PRO
1	C	68	ASP
1	E	81	PRO
1	F	4	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	157/164 (96%)	136 (87%)	21 (13%)	4	7
1	B	157/164 (96%)	144 (92%)	13 (8%)	10	19
1	C	155/164 (94%)	135 (87%)	20 (13%)	4	7
1	D	156/164 (95%)	139 (89%)	17 (11%)	6	12
1	E	157/164 (96%)	140 (89%)	17 (11%)	6	12
1	F	156/164 (95%)	136 (87%)	20 (13%)	4	7
1	G	157/164 (96%)	139 (88%)	18 (12%)	5	10
1	H	157/164 (96%)	137 (87%)	20 (13%)	4	8
All	All	1252/1312 (95%)	1106 (88%)	146 (12%)	5	9

All (146) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	5	PHE
1	A	10	ILE
1	A	17	GLN
1	A	21	GLU
1	A	24	LEU
1	A	29	THR
1	A	79	ILE
1	A	84	ILE
1	A	87	THR
1	A	100	MET
1	A	101	LEU
1	A	103	ILE
1	A	106	LEU
1	A	129	ILE
1	A	143	MET
1	A	146	LYS
1	A	151	VAL
1	A	160	SER
1	A	181	LEU
1	A	185	HIS
1	B	10	ILE
1	B	20	MET
1	B	21	GLU
1	B	24	LEU
1	B	29	THR
1	B	79	ILE
1	B	87	THR
1	B	101	LEU
1	B	129	ILE
1	B	131	LEU
1	B	136	ILE
1	B	143	MET
1	B	181	LEU
1	C	6	SER
1	C	17	GLN
1	C	21	GLU
1	C	28	LEU
1	C	29	THR
1	C	58	HIS
1	C	68	ASP
1	C	79	ILE
1	C	87	THR

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Mol	Chain	Res	Type
1	C	100	MET
1	C	103	ILE
1	C	106	LEU
1	C	129	ILE
1	C	136	ILE
1	C	143	MET
1	C	146	LYS
1	C	151	VAL
1	C	152	ILE
1	C	179	LYS
1	C	181	LEU
1	D	3	LEU
1	D	5	PHE
1	D	17	GLN
1	D	20	MET
1	D	21	GLU
1	D	29	THR
1	D	42	LYS
1	D	58	HIS
1	D	79	ILE
1	D	87	THR
1	D	101	LEU
1	D	123	LYS
1	D	129	ILE
1	D	136	ILE
1	D	143	MET
1	D	151	VAL
1	D	181	LEU
1	E	2	GLU
1	E	7	ASN
1	E	10	ILE
1	E	17	GLN
1	E	27	ASP
1	E	29	THR
1	E	53	LYS
1	E	87	THR
1	E	101	LEU
1	E	103	ILE
1	E	122	GLN
1	E	143	MET
1	E	144	THR
1	E	146	LYS

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Mol	Chain	Res	Type
1	E	147	LEU
1	E	151	VAL
1	E	157	ILE
1	F	3	LEU
1	F	5	PHE
1	F	17	GLN
1	F	20	MET
1	F	29	THR
1	F	42	LYS
1	F	58	HIS
1	F	79	ILE
1	F	87	THR
1	F	101	LEU
1	F	103	ILE
1	F	129	ILE
1	F	131	LEU
1	F	143	MET
1	F	146	LYS
1	F	147	LEU
1	F	148	HIS
1	F	151	VAL
1	F	172	ILE
1	F	182	TRP
1	G	5	PHE
1	G	10	ILE
1	G	17	GLN
1	G	19	ASP
1	G	21	GLU
1	G	58	HIS
1	G	80	CYS
1	G	88	ARG
1	G	101	LEU
1	G	129	ILE
1	G	136	ILE
1	G	143	MET
1	G	144	THR
1	G	146	LYS
1	G	166	VAL
1	G	175	THR
1	G	179	LYS
1	G	181	LEU
1	H	5	PHE

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Mol	Chain	Res	Type
1	H	10	ILE
1	H	17	GLN
1	H	20	MET
1	H	26	LEU
1	H	29	THR
1	H	33	MET
1	H	35	GLU
1	H	36	THR
1	H	58	HIS
1	H	79	ILE
1	H	84	ILE
1	H	101	LEU
1	H	103	ILE
1	H	129	ILE
1	H	131	LEU
1	H	143	MET
1	H	146	LYS
1	H	165	GLN
1	H	172	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (47) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	15	HIS
1	A	97	GLN
1	A	165	GLN
1	B	7	ASN
1	B	15	HIS
1	B	97	GLN
1	B	104	GLN
1	B	139	GLN
1	C	16	ASN
1	C	104	GLN
1	C	122	GLN
1	C	178	ASN
1	D	15	HIS
1	D	63	ASN
1	D	90	ASN
1	D	97	GLN
1	D	98	HIS
1	D	104	GLN

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Mol	Chain	Res	Type
1	D	139	GLN
1	D	145	GLN
1	D	178	ASN
1	E	7	ASN
1	E	15	HIS
1	E	17	GLN
1	E	37	HIS
1	E	40	GLN
1	E	104	GLN
1	E	122	GLN
1	E	139	GLN
1	E	178	ASN
1	F	15	HIS
1	F	58	HIS
1	F	63	ASN
1	F	180	HIS
1	G	58	HIS
1	G	90	ASN
1	G	104	GLN
1	G	139	GLN
1	G	145	GLN
1	G	165	GLN
1	G	178	ASN
1	G	180	HIS
1	H	15	HIS
1	H	58	HIS
1	H	139	GLN
1	H	180	HIS

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 8 ligands modelled in this entry, 8 are unknown - 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 [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	185/192 (96%)	-0.38	1 (0%) 87 87	69, 87, 121, 149	0
1	B	185/192 (96%)	0.01	3 (1%) 70 69	70, 94, 126, 157	0
1	C	182/192 (94%)	-0.29	1 (0%) 87 87	67, 91, 124, 150	0
1	D	183/192 (95%)	-0.24	0 100 100	66, 90, 120, 150	0
1	E	185/192 (96%)	-0.23	0 100 100	64, 94, 131, 148	0
1	F	184/192 (95%)	0.03	1 (0%) 87 87	73, 100, 128, 142	0
1	G	185/192 (96%)	-0.09	0 100 100	73, 104, 140, 149	0
1	H	185/192 (96%)	-0.01	1 (0%) 87 87	80, 106, 132, 153	0
All	All	1474/1536 (95%)	-0.15	7 (0%) 87 87	64, 96, 131, 157	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	184	GLY	3.1
1	A	147	LEU	2.5
1	B	1	LEU	2.5
1	H	28	LEU	2.4
1	B	28	LEU	2.3
1	C	24	LEU	2.2
1	B	23	ILE	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
2	UNL	H	191	10/-	0.81	0.10	92,105,117,119	0
2	UNL	G	191	10/-	0.87	0.10	106,113,120,123	0
2	UNL	D	191	10/-	0.89	0.10	78,100,105,119	0
2	UNL	E	191	10/-	0.90	0.12	105,123,132,137	0
2	UNL	F	191	10/-	0.90	0.09	100,115,125,132	0
2	UNL	B	191	10/-	0.90	0.09	68,92,96,106	0
2	UNL	A	191	10/-	0.90	0.10	77,96,105,105	0
2	UNL	C	191	10/-	0.93	0.11	71,99,107,111	0

6.5 Other polymers [i](#)

There are no such residues in this entry.