



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 30, 2026 – 12:12 PM EDT

PDB ID : 9CG8 / pdb_00009cg8
Title : CRYSTAL STRUCTURE OF THE P285S VARIANT OF SERINE HYDROXYMETHYLTRANSFERASE 8 FROM SOYBEAN CULTIVAR FORREST
Authors : Beamer, L.J.; Samarakoon, V.; Owuocha, L.F.
Deposited on : 2024-06-28
Resolution : 1.90 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

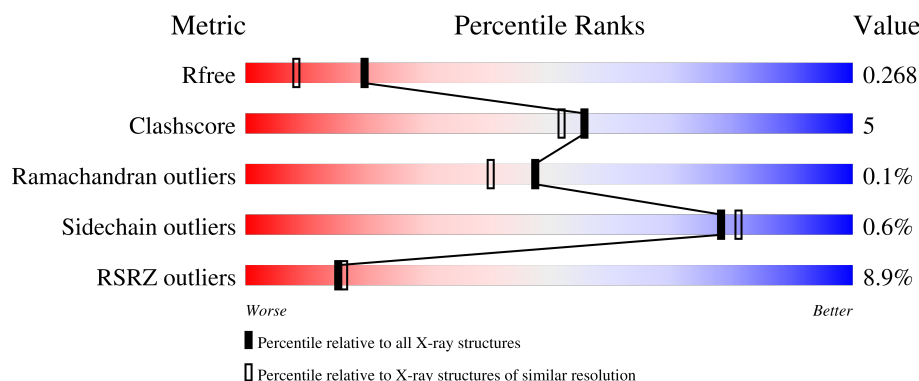
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	496	8% 81% 9% 9%
1	B	496	18% 79% 11% 9%
1	C	496	7% 82% 9% 9%
1	D	496	5% 82% 9% 8%
1	E	496	5% 80% 12% 8%

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Mol	Chain	Length	Quality of chain
1	F	496	5% 84% 6% 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 21383 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 Serine hydroxymethyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	451	Total	C	N	O	P	S	0	1	0
			3397	2155	580	644	1	17			
1	B	451	Total	C	N	O	P	S	0	2	0
			3353	2117	574	643	1	18			
1	C	452	Total	C	N	O	P	S	0	1	0
			3411	2168	585	640	1	17			
1	D	454	Total	C	N	O	P	S	0	1	0
			3476	2209	593	656	1	17			
1	E	458	Total	C	N	O	P	S	0	3	0
			3467	2201	587	659	1	19			
1	F	450	Total	C	N	O	P	S	0	2	0
			3414	2167	581	647	1	18			

There are 156 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	expression tag	UNP K4FW35
A	-23	GLY	-	expression tag	UNP K4FW35
A	-22	MET	-	expression tag	UNP K4FW35
A	-21	HIS	-	expression tag	UNP K4FW35
A	-20	HIS	-	expression tag	UNP K4FW35
A	-19	HIS	-	expression tag	UNP K4FW35
A	-18	HIS	-	expression tag	UNP K4FW35
A	-17	HIS	-	expression tag	UNP K4FW35
A	-16	HIS	-	expression tag	UNP K4FW35
A	-15	SER	-	expression tag	UNP K4FW35
A	-14	SER	-	expression tag	UNP K4FW35
A	-13	GLY	-	expression tag	UNP K4FW35
A	-12	VAL	-	expression tag	UNP K4FW35
A	-11	ASP	-	expression tag	UNP K4FW35
A	-10	LEU	-	expression tag	UNP K4FW35
A	-9	GLY	-	expression tag	UNP K4FW35
A	-8	THR	-	expression tag	UNP K4FW35

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLU	-	expression tag	UNP K4FW35
A	-6	ASN	-	expression tag	UNP K4FW35
A	-5	LEU	-	expression tag	UNP K4FW35
A	-4	TYR	-	expression tag	UNP K4FW35
A	-3	PHE	-	expression tag	UNP K4FW35
A	-2	GLN	-	expression tag	UNP K4FW35
A	-1	SER	-	expression tag	UNP K4FW35
A	0	ASN	-	expression tag	UNP K4FW35
A	285	SER	PRO	engineered mutation	UNP K4FW35
B	-24	MET	-	expression tag	UNP K4FW35
B	-23	GLY	-	expression tag	UNP K4FW35
B	-22	MET	-	expression tag	UNP K4FW35
B	-21	HIS	-	expression tag	UNP K4FW35
B	-20	HIS	-	expression tag	UNP K4FW35
B	-19	HIS	-	expression tag	UNP K4FW35
B	-18	HIS	-	expression tag	UNP K4FW35
B	-17	HIS	-	expression tag	UNP K4FW35
B	-16	HIS	-	expression tag	UNP K4FW35
B	-15	SER	-	expression tag	UNP K4FW35
B	-14	SER	-	expression tag	UNP K4FW35
B	-13	GLY	-	expression tag	UNP K4FW35
B	-12	VAL	-	expression tag	UNP K4FW35
B	-11	ASP	-	expression tag	UNP K4FW35
B	-10	LEU	-	expression tag	UNP K4FW35
B	-9	GLY	-	expression tag	UNP K4FW35
B	-8	THR	-	expression tag	UNP K4FW35
B	-7	GLU	-	expression tag	UNP K4FW35
B	-6	ASN	-	expression tag	UNP K4FW35
B	-5	LEU	-	expression tag	UNP K4FW35
B	-4	TYR	-	expression tag	UNP K4FW35
B	-3	PHE	-	expression tag	UNP K4FW35
B	-2	GLN	-	expression tag	UNP K4FW35
B	-1	SER	-	expression tag	UNP K4FW35
B	0	ASN	-	expression tag	UNP K4FW35
B	285	SER	PRO	engineered mutation	UNP K4FW35
C	-24	MET	-	expression tag	UNP K4FW35
C	-23	GLY	-	expression tag	UNP K4FW35
C	-22	MET	-	expression tag	UNP K4FW35
C	-21	HIS	-	expression tag	UNP K4FW35
C	-20	HIS	-	expression tag	UNP K4FW35
C	-19	HIS	-	expression tag	UNP K4FW35
C	-18	HIS	-	expression tag	UNP K4FW35

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-17	HIS	-	expression tag	UNP K4FW35
C	-16	HIS	-	expression tag	UNP K4FW35
C	-15	SER	-	expression tag	UNP K4FW35
C	-14	SER	-	expression tag	UNP K4FW35
C	-13	GLY	-	expression tag	UNP K4FW35
C	-12	VAL	-	expression tag	UNP K4FW35
C	-11	ASP	-	expression tag	UNP K4FW35
C	-10	LEU	-	expression tag	UNP K4FW35
C	-9	GLY	-	expression tag	UNP K4FW35
C	-8	THR	-	expression tag	UNP K4FW35
C	-7	GLU	-	expression tag	UNP K4FW35
C	-6	ASN	-	expression tag	UNP K4FW35
C	-5	LEU	-	expression tag	UNP K4FW35
C	-4	TYR	-	expression tag	UNP K4FW35
C	-3	PHE	-	expression tag	UNP K4FW35
C	-2	GLN	-	expression tag	UNP K4FW35
C	-1	SER	-	expression tag	UNP K4FW35
C	0	ASN	-	expression tag	UNP K4FW35
C	285	SER	PRO	engineered mutation	UNP K4FW35
D	-24	MET	-	expression tag	UNP K4FW35
D	-23	GLY	-	expression tag	UNP K4FW35
D	-22	MET	-	expression tag	UNP K4FW35
D	-21	HIS	-	expression tag	UNP K4FW35
D	-20	HIS	-	expression tag	UNP K4FW35
D	-19	HIS	-	expression tag	UNP K4FW35
D	-18	HIS	-	expression tag	UNP K4FW35
D	-17	HIS	-	expression tag	UNP K4FW35
D	-16	HIS	-	expression tag	UNP K4FW35
D	-15	SER	-	expression tag	UNP K4FW35
D	-14	SER	-	expression tag	UNP K4FW35
D	-13	GLY	-	expression tag	UNP K4FW35
D	-12	VAL	-	expression tag	UNP K4FW35
D	-11	ASP	-	expression tag	UNP K4FW35
D	-10	LEU	-	expression tag	UNP K4FW35
D	-9	GLY	-	expression tag	UNP K4FW35
D	-8	THR	-	expression tag	UNP K4FW35
D	-7	GLU	-	expression tag	UNP K4FW35
D	-6	ASN	-	expression tag	UNP K4FW35
D	-5	LEU	-	expression tag	UNP K4FW35
D	-4	TYR	-	expression tag	UNP K4FW35
D	-3	PHE	-	expression tag	UNP K4FW35
D	-2	GLN	-	expression tag	UNP K4FW35

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	SER	-	expression tag	UNP K4FW35
D	0	ASN	-	expression tag	UNP K4FW35
D	285	SER	PRO	engineered mutation	UNP K4FW35
E	-24	MET	-	expression tag	UNP K4FW35
E	-23	GLY	-	expression tag	UNP K4FW35
E	-22	MET	-	expression tag	UNP K4FW35
E	-21	HIS	-	expression tag	UNP K4FW35
E	-20	HIS	-	expression tag	UNP K4FW35
E	-19	HIS	-	expression tag	UNP K4FW35
E	-18	HIS	-	expression tag	UNP K4FW35
E	-17	HIS	-	expression tag	UNP K4FW35
E	-16	HIS	-	expression tag	UNP K4FW35
E	-15	SER	-	expression tag	UNP K4FW35
E	-14	SER	-	expression tag	UNP K4FW35
E	-13	GLY	-	expression tag	UNP K4FW35
E	-12	VAL	-	expression tag	UNP K4FW35
E	-11	ASP	-	expression tag	UNP K4FW35
E	-10	LEU	-	expression tag	UNP K4FW35
E	-9	GLY	-	expression tag	UNP K4FW35
E	-8	THR	-	expression tag	UNP K4FW35
E	-7	GLU	-	expression tag	UNP K4FW35
E	-6	ASN	-	expression tag	UNP K4FW35
E	-5	LEU	-	expression tag	UNP K4FW35
E	-4	TYR	-	expression tag	UNP K4FW35
E	-3	PHE	-	expression tag	UNP K4FW35
E	-2	GLN	-	expression tag	UNP K4FW35
E	-1	SER	-	expression tag	UNP K4FW35
E	0	ASN	-	expression tag	UNP K4FW35
E	285	SER	PRO	engineered mutation	UNP K4FW35
F	-24	MET	-	expression tag	UNP K4FW35
F	-23	GLY	-	expression tag	UNP K4FW35
F	-22	MET	-	expression tag	UNP K4FW35
F	-21	HIS	-	expression tag	UNP K4FW35
F	-20	HIS	-	expression tag	UNP K4FW35
F	-19	HIS	-	expression tag	UNP K4FW35
F	-18	HIS	-	expression tag	UNP K4FW35
F	-17	HIS	-	expression tag	UNP K4FW35
F	-16	HIS	-	expression tag	UNP K4FW35
F	-15	SER	-	expression tag	UNP K4FW35
F	-14	SER	-	expression tag	UNP K4FW35
F	-13	GLY	-	expression tag	UNP K4FW35
F	-12	VAL	-	expression tag	UNP K4FW35

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-11	ASP	-	expression tag	UNP K4FW35
F	-10	LEU	-	expression tag	UNP K4FW35
F	-9	GLY	-	expression tag	UNP K4FW35
F	-8	THR	-	expression tag	UNP K4FW35
F	-7	GLU	-	expression tag	UNP K4FW35
F	-6	ASN	-	expression tag	UNP K4FW35
F	-5	LEU	-	expression tag	UNP K4FW35
F	-4	TYR	-	expression tag	UNP K4FW35
F	-3	PHE	-	expression tag	UNP K4FW35
F	-2	GLN	-	expression tag	UNP K4FW35
F	-1	SER	-	expression tag	UNP K4FW35
F	0	ASN	-	expression tag	UNP K4FW35
F	285	SER	PRO	engineered mutation	UNP K4FW35

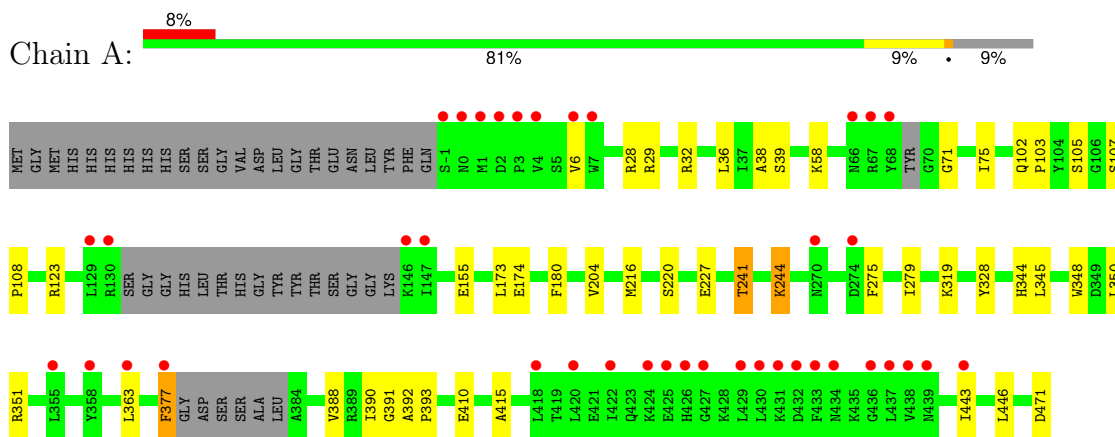
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	154	Total	O	0	0
			154	154		
2	B	82	Total	O	0	0
			82	82		
2	C	140	Total	O	0	0
			140	140		
2	D	164	Total	O	0	0
			164	164		
2	E	166	Total	O	0	0
			166	166		
2	F	159	Total	O	0	0
			159	159		

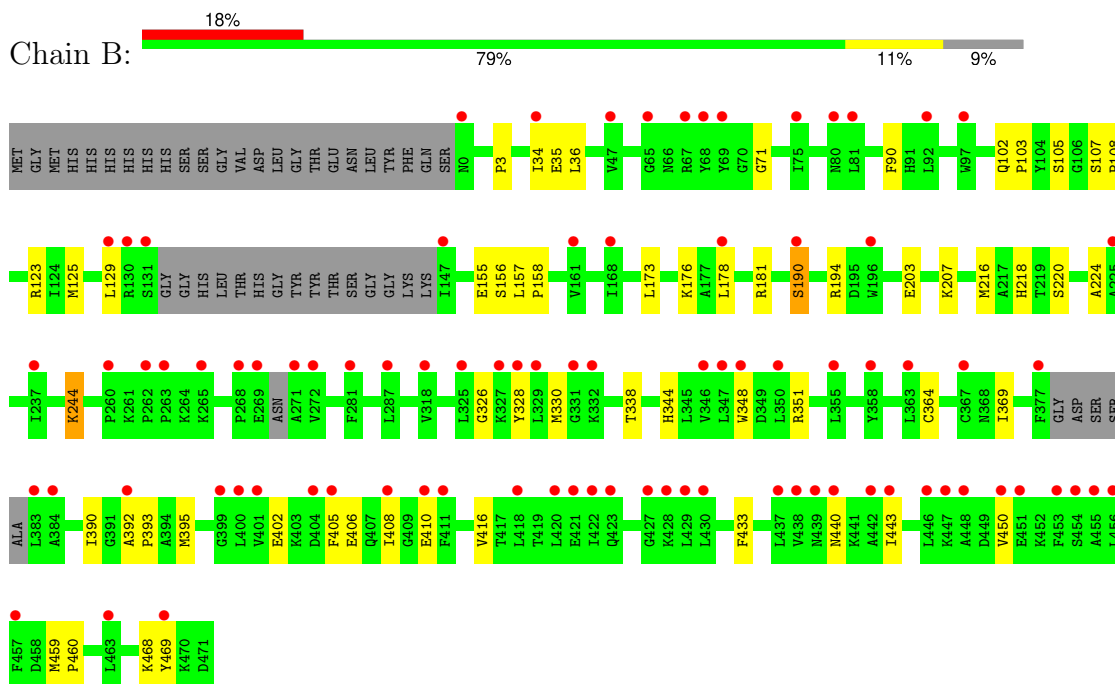
3 Residue-property plots

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.

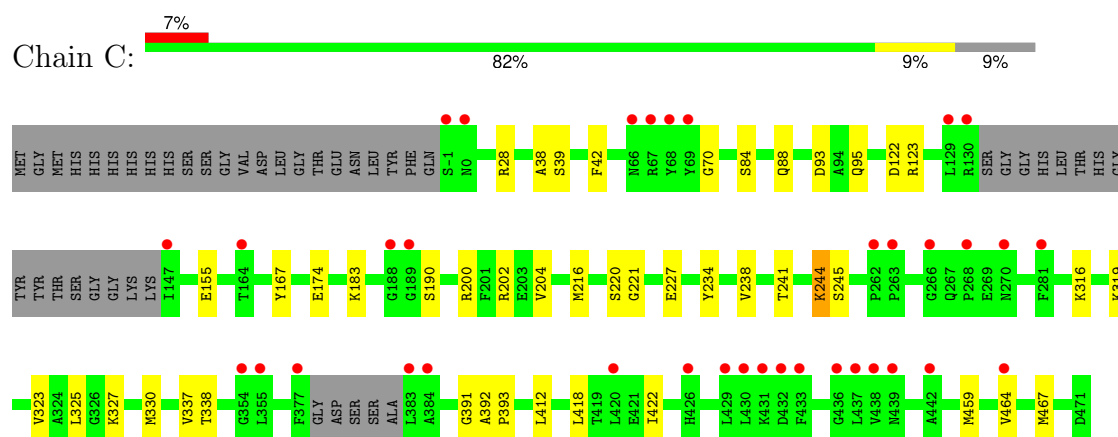
• Molecule 1: Serine hydroxymethyltransferase



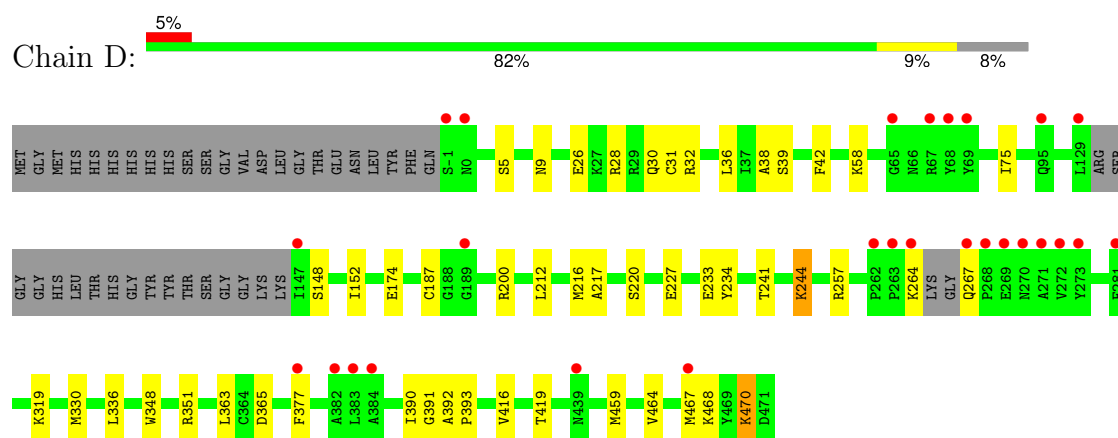
• Molecule 1: Serine hydroxymethyltransferase



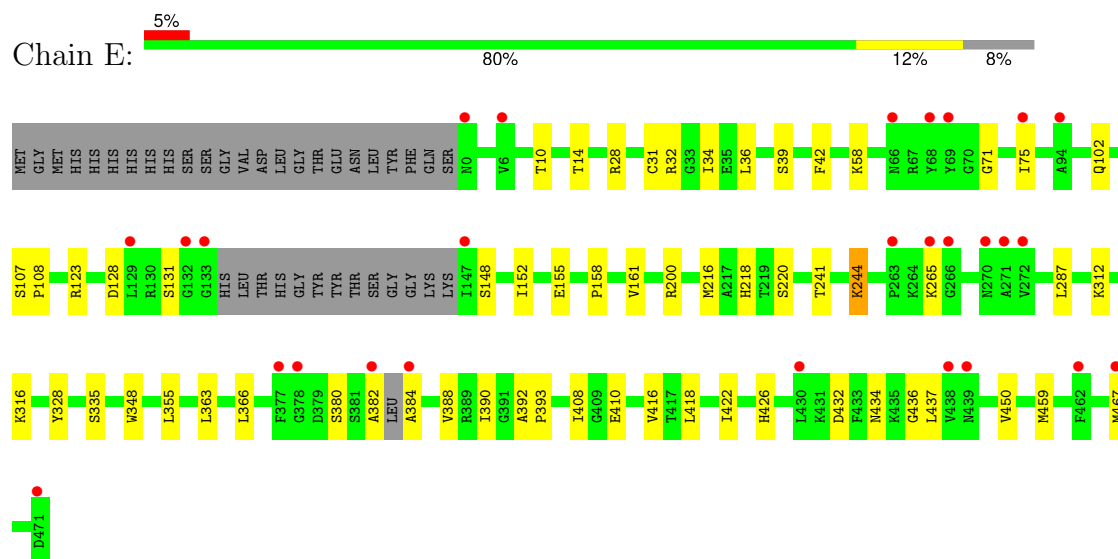
• Molecule 1: Serine hydroxymethyltransferase



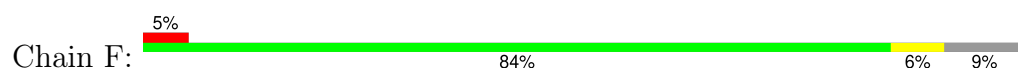
- Molecule 1: Serine hydroxymethyltransferase

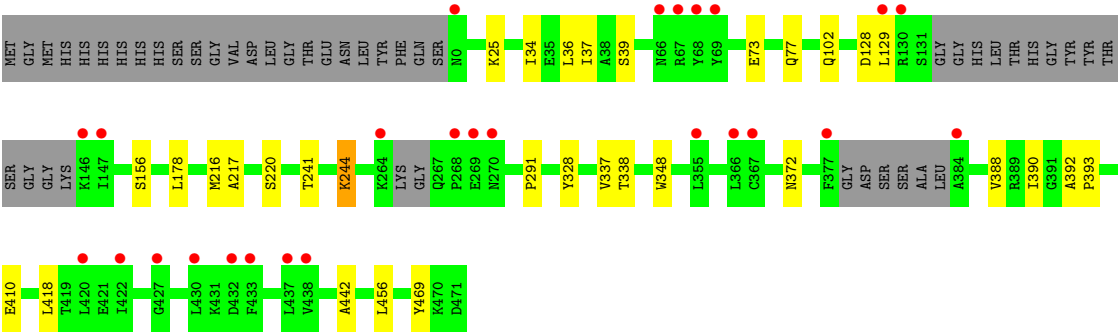


- Molecule 1: Serine hydroxymethyltransferase



- Molecule 1: Serine hydroxymethyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	174.18Å 174.18Å 183.66Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.50 – 1.90 48.50 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.50-1.90) 99.6 (48.50-1.90)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.26 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.227 , 0.266 0.231 , 0.268	Depositor DCC
R_{free} test set	12634 reflections (4.49%)	wwPDB-VP
Wilson B-factor (Å ²)	30.4	Xtriage
Anisotropy	0.422	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.006 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	21383	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 35.98 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.3619e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: LLP

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	0.33	0/3437	0.54	0/4665
1	B	0.30	0/3397	0.52	0/4621
1	C	0.35	0/3453	0.56	0/4685
1	D	0.37	0/3518	0.57	0/4767
1	E	0.38	0/3512	0.58	0/4766
1	F	0.35	0/3458	0.57	0/4693
All	All	0.35	0/20775	0.56	0/28197

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	3397	0	3209	32	0
1	B	3353	0	3059	39	0
1	C	3411	0	3244	29	0
1	D	3476	0	3365	34	0
1	E	3467	0	3306	38	0
1	F	3414	0	3240	19	0
2	A	154	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	82	0	0	0	0
2	C	140	0	0	1	0
2	D	164	0	0	2	0
2	E	166	0	0	1	0
2	F	159	0	0	0	0
All	All	21383	0	19423	187	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 5.

All (187) 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:B:34:ILE:HA	1:B:395:MET:HE2	1.47	0.95
1:B:34:ILE:HG13	1:B:450:VAL:HG13	1.67	0.75
1:F:216:MET:HG3	1:F:220:SER:HB3	1.68	0.74
1:E:34:ILE:HG13	1:E:450:VAL:HG13	1.69	0.72
1:B:459:MET:HG3	1:B:460:PRO:N	2.04	0.72
1:D:31:CYS:HB3	1:D:459:MET:HG2	1.71	0.70
1:D:58:LYS:HG2	1:D:75:ILE:HG13	1.73	0.70
1:E:58:LYS:HE2	1:E:71:GLY:O	1.98	0.63
1:D:351:ARG:NH1	1:D:377:PHE:O	2.30	0.63
1:E:36:LEU:HB3	1:E:390:ILE:HG23	1.80	0.63
1:E:382:ALA:O	1:E:384:ALA:N	2.32	0.62
1:B:203:GLU:O	1:B:207:LYS:HG3	2.01	0.61
1:A:28:ARG:NE	1:A:471:ASP:O	2.33	0.60
1:D:39:SER:HB2	1:D:244[B]:LLP:HE3	1.83	0.60
1:B:34:ILE:HA	1:B:395:MET:CE	2.25	0.60
1:D:363:LEU:HD13	1:D:419:THR:OG1	2.01	0.60
1:E:434:ASN:HA	1:E:437:LEU:HD12	1.82	0.60
1:D:39:SER:HB2	1:D:244[A]:LLP:HE3	1.84	0.60
1:B:351:ARG:HG3	1:B:351:ARG:HH11	1.67	0.60
1:D:330:MET:HE2	1:D:336:LEU:HD12	1.84	0.60
1:E:131:SER:HB3	1:E:161:VAL:HG13	1.83	0.59
1:D:264:LYS:O	1:D:267:GLN:N	2.35	0.59
1:D:233:GLU:HG2	1:D:234:TYR:CE2	2.37	0.59
1:F:217:ALA:HB1	1:F:241:THR:HG22	1.84	0.58
1:A:28:ARG:HG2	1:A:28:ARG:HH11	1.67	0.58
1:F:36:LEU:HB3	1:F:390:ILE:HG23	1.86	0.58
1:B:328:TYR:OH	1:B:410:GLU:OE1	2.21	0.58
1:E:328:TYR:OH	1:E:410:GLU:OE2	2.21	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:42:PHE:CD1	1:E:459:MET:HE1	2.38	0.58
1:E:58:LYS:HG2	1:E:75:ILE:HG12	1.84	0.58
1:C:70:GLY:HA3	1:D:365:ASP:OD2	2.04	0.57
1:A:363:LEU:HD12	1:A:415:ALA:HB1	1.85	0.57
1:D:5:SER:HA	1:D:9:ASN:HD22	1.70	0.56
1:E:128:ASP:HB2	1:E:158:PRO:HB2	1.87	0.56
1:C:122:ASP:OD2	1:C:183:LYS:HE2	2.06	0.56
1:B:36:LEU:HB3	1:B:390:ILE:HG23	1.87	0.55
1:D:216:MET:HE3	1:D:220:SER:HB2	1.88	0.55
1:D:174:GLU:OE1	1:D:200:ARG:NH2	2.38	0.55
1:B:178:LEU:O	1:B:181:ARG:HD3	2.06	0.55
1:D:464:VAL:HA	1:D:467:MET:HG3	1.90	0.53
1:A:328:TYR:OH	1:A:410:GLU:OE1	2.27	0.53
1:D:227:GLU:HB3	1:D:319:LYS:HD3	1.90	0.53
1:C:84:SER:O	1:C:88:GLN:HG3	2.08	0.52
1:E:148:SER:O	1:E:152:ILE:HG13	2.08	0.52
1:C:337:VAL:HG12	1:C:338:THR:HG23	1.91	0.52
1:B:129:LEU:HD23	1:B:156:SER:HB2	1.90	0.52
1:D:42:PHE:CD1	1:D:459:MET:HE1	2.45	0.52
1:E:363:LEU:HD11	1:E:418:LEU:HG	1.92	0.51
1:E:102:GLN:HB3	1:E:287:LEU:HD11	1.91	0.51
1:F:39:SER:HB2	1:F:244[A]:LLP:HE3	1.93	0.51
1:B:405:PHE:HA	1:B:408:ILE:HD12	1.92	0.50
1:F:39:SER:HB2	1:F:244[B]:LLP:HE3	1.92	0.50
1:D:348:TRP:CH2	1:D:416:VAL:HG21	2.46	0.50
1:C:216:MET:HG3	1:C:220:SER:HB3	1.92	0.50
1:A:363:LEU:HD21	1:A:443:ILE:HD11	1.93	0.50
1:B:216:MET:HG3	1:B:220:SER:HB3	1.94	0.50
1:B:218:HIS:HA	1:B:244[A]:LLP:HD2	1.94	0.49
1:A:377:PHE:N	1:A:377:PHE:CD1	2.79	0.49
1:C:28:ARG:HA	1:C:467:MET:HE1	1.95	0.49
1:A:241:THR:HG21	1:A:244[A]:LLP:HE2	1.95	0.48
1:F:102:GLN:OE1	1:F:291:PRO:HG3	2.14	0.48
1:A:28:ARG:CZ	1:A:32:ARG:HD2	2.43	0.48
1:B:125:MET:HE1	1:B:176:LYS:HB3	1.96	0.48
1:C:325:LEU:HD11	1:C:412:LEU:HD12	1.96	0.48
1:A:275:PHE:HB3	1:A:279:ILE:HD12	1.95	0.48
1:C:39:SER:HB2	1:C:244[B]:LLP:HE3	1.96	0.48
1:D:148:SER:O	1:D:152:ILE:HG13	2.12	0.48
1:F:392:ALA:N	1:F:393:PRO:CD	2.77	0.47
1:A:39:SER:HB2	1:A:244[B]:LLP:HE3	1.94	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:257:ARG:HG2	2:D:587:HOH:O	2.14	0.47
1:C:323:VAL:O	1:C:327:LYS:HG3	2.15	0.47
1:B:468:LYS:HD3	1:B:469:TYR:CE2	2.49	0.47
1:D:58:LYS:CG	1:D:75:ILE:HG13	2.44	0.47
1:F:25:LYS:HG3	1:F:469:TYR:CE2	2.50	0.47
1:B:402:GLU:O	1:B:406:GLU:HG3	2.14	0.46
1:F:418:LEU:HD21	1:F:442:ALA:HB1	1.98	0.46
1:C:392:ALA:N	1:C:393:PRO:HD3	2.30	0.46
1:E:58:LYS:NZ	2:E:513:HOH:O	2.47	0.46
1:E:366:LEU:HD12	1:E:437:LEU:HD22	1.97	0.46
1:E:392:ALA:N	1:E:393:PRO:CD	2.79	0.46
1:D:233:GLU:HG2	1:D:234:TYR:CD2	2.50	0.46
1:B:90:PHE:CE1	1:B:224:ALA:HB2	2.51	0.46
1:A:58:LYS:HG2	1:A:75:ILE:HG12	1.97	0.46
1:B:190:SER:HB2	1:B:244[A]:LLP:H5'2	1.98	0.46
1:D:58:LYS:HG2	1:D:75:ILE:CG1	2.44	0.46
1:C:174:GLU:HG3	1:C:204:VAL:HG22	1.98	0.46
1:E:123:ARG:HA	1:E:155:GLU:O	2.15	0.46
1:E:107:SER:HB3	1:E:108:PRO:HD3	1.98	0.46
1:D:470:LYS:HB3	1:D:470:LYS:HE3	1.45	0.45
1:A:58:LYS:HE2	1:A:71:GLY:O	2.16	0.45
1:F:392:ALA:N	1:F:393:PRO:HD3	2.31	0.45
1:A:227:GLU:HB3	1:A:319:LYS:HD3	1.99	0.45
1:E:312:LYS:HE3	1:E:316:LYS:HE2	1.98	0.45
1:A:348:TRP:HB3	1:A:388:VAL:HG23	1.98	0.45
1:C:93:ASP:OD2	1:C:95:GLN:HB2	2.16	0.45
1:A:415:ALA:HB2	1:A:446:LEU:HD21	1.97	0.45
1:C:183:LYS:HB3	1:C:183:LYS:HE3	1.63	0.45
1:E:31[B]:CYS:O	1:E:32:ARG:HD3	2.17	0.45
1:E:216:MET:HG3	1:E:220:SER:HB3	1.99	0.45
1:A:174:GLU:HG3	1:A:204:VAL:HG22	1.99	0.45
1:D:36:LEU:HB3	1:D:390:ILE:HG23	1.99	0.45
1:B:348:TRP:CH2	1:B:416:VAL:HG21	2.52	0.44
1:B:218:HIS:HA	1:B:244[B]:LLP:HD2	1.99	0.44
1:E:426:HIS:O	1:E:432:ASP:HB3	2.17	0.44
1:A:36:LEU:HB3	1:A:390:ILE:HG23	1.98	0.44
1:A:344:HIS:CE1	1:A:345:LEU:HD23	2.53	0.44
1:B:364:CYS:HB3	1:B:369:ILE:HB	2.00	0.44
1:C:123:ARG:HA	1:C:155:GLU:O	2.18	0.44
1:E:34:ILE:HD13	1:E:408:ILE:HG12	2.00	0.44
1:B:157:LEU:HD12	1:B:158:PRO:HD2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:ALA:N	1:A:393:PRO:CD	2.81	0.44
1:F:73:GLU:O	1:F:77:GLN:HG3	2.18	0.44
1:F:129:LEU:HG	1:F:156:SER:HB2	2.00	0.44
1:A:102:GLN:N	1:A:103:PRO:CD	2.81	0.43
1:E:31[B]:CYS:SG	1:E:459:MET:HB2	2.58	0.43
1:D:38:ALA:HA	1:D:391:GLY:HA3	2.00	0.43
1:A:39:SER:HB2	1:A:244[A]:LLP:HE3	2.01	0.43
1:C:39:SER:HB2	1:C:244[A]:LLP:HE3	2.00	0.43
1:D:392:ALA:N	1:D:393:PRO:CD	2.82	0.43
1:B:105:SER:C	1:B:108:PRO:HD2	2.44	0.43
1:B:433:PHE:CD2	1:B:433:PHE:C	2.97	0.43
1:C:38:ALA:HA	1:C:391:GLY:HA3	2.00	0.43
1:C:42:PHE:CD1	1:C:459:MET:HE1	2.54	0.43
1:C:221:GLY:N	1:C:245:SER:OG	2.48	0.43
1:E:218:HIS:HA	1:E:244[A]:LLP:HD2	2.01	0.43
1:E:348:TRP:HB3	1:E:388:VAL:HG23	2.01	0.42
1:F:216:MET:HG3	1:F:216:MET:O	2.18	0.42
1:D:187:CYS:SG	1:D:212:LEU:HD11	2.59	0.42
1:A:351:ARG:HE	1:A:351:ARG:HB2	1.51	0.42
1:D:28:ARG:CZ	1:D:32:ARG:HG3	2.50	0.42
1:E:10:THR:HG22	1:E:14:THR:OG1	2.20	0.42
1:F:178:LEU:HD23	1:F:178:LEU:HA	1.80	0.42
1:B:3:PRO:HG2	1:E:335:SER:OG	2.19	0.42
1:E:28:ARG:HA	1:E:467:MET:HE1	2.02	0.42
1:B:102:GLN:N	1:B:103:PRO:CD	2.83	0.42
1:B:440:ASN:HD22	1:B:443:ILE:H	1.67	0.42
1:F:34:ILE:HG22	1:F:36:LEU:HG	2.00	0.42
1:A:6:VAL:HG22	1:C:330:MET:HE3	2.01	0.42
1:D:26:GLU:O	1:D:30:GLN:HG3	2.20	0.42
1:E:392:ALA:N	1:E:393:PRO:HD3	2.35	0.42
1:E:392:ALA:H	1:E:393:PRO:HD3	1.85	0.42
1:A:105:SER:C	1:A:108:PRO:HD2	2.45	0.42
1:C:190:SER:HB3	1:C:244[A]:LLP:H5'2	2.02	0.42
1:A:38:ALA:HA	1:A:391:GLY:HA3	2.02	0.41
1:A:392:ALA:N	1:A:393:PRO:HD3	2.35	0.41
1:E:42:PHE:CG	1:E:459:MET:HE1	2.55	0.41
1:C:227:GLU:HB3	1:C:319:LYS:HD3	2.02	0.41
1:B:459:MET:HG3	1:B:460:PRO:CD	2.49	0.41
1:E:218:HIS:HA	1:E:244[B]:LLP:HD2	2.02	0.41
1:A:350:LEU:HD11	1:A:388:VAL:HG13	2.03	0.41
1:B:344:HIS:HD1	1:B:344:HIS:H	1.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:58:LYS:HD2	1:D:58:LYS:HA	1.83	0.41
1:B:194:ARG:NH2	1:B:338:THR:O	2.53	0.41
1:C:167:TYR:HB3	2:C:607:HOH:O	2.20	0.41
1:F:37:ILE:HG23	1:F:372:ASN:OD1	2.19	0.41
1:F:348:TRP:HB3	1:F:388:VAL:HG23	2.02	0.41
1:A:123:ARG:HB3	1:A:180:PHE:CE2	2.55	0.41
1:A:244[A]:LLP:H4'1	1:A:244[A]:LLP:OP3	2.20	0.41
1:B:105:SER:O	1:B:108:PRO:HD2	2.20	0.41
1:C:216:MET:HE2	1:C:238:VAL:HG12	2.01	0.41
1:C:227:GLU:CD	1:C:316:LYS:HG2	2.45	0.41
1:D:468:LYS:NZ	2:D:512:HOH:O	2.52	0.41
1:B:107:SER:HB3	1:B:108:PRO:HD3	2.03	0.41
1:B:326:GLY:O	1:B:330:MET:HB2	2.21	0.41
1:C:216:MET:HE2	1:C:238:VAL:CG1	2.50	0.41
1:E:200:ARG:HD2	1:E:200:ARG:HA	1.87	0.41
1:A:29:ARG:HD2	1:B:71:GLY:HA2	2.03	0.41
1:F:328:TYR:OH	1:F:410:GLU:OE1	2.34	0.41
1:F:337:VAL:HG12	1:F:338:THR:HG23	2.03	0.41
1:A:123:ARG:HA	1:A:155:GLU:O	2.21	0.41
1:B:123:ARG:HA	1:B:155:GLU:O	2.20	0.41
1:B:190:SER:HB2	1:B:244[A]:LLP:C5'	2.51	0.41
1:D:363:LEU:HD23	1:D:363:LEU:C	2.46	0.41
1:B:392:ALA:N	1:B:393:PRO:CD	2.84	0.41
1:E:355:LEU:HD23	1:E:355:LEU:HA	1.93	0.41
1:C:200:ARG:HA	1:C:200:ARG:HD2	1.86	0.40
1:E:39:SER:HB2	1:E:244[B]:LLP:HE3	2.03	0.40
1:D:392:ALA:N	1:D:393:PRO:HD3	2.36	0.40
1:A:216:MET:HG3	1:A:220:SER:HB3	2.03	0.40
1:C:392:ALA:N	1:C:393:PRO:CD	2.84	0.40
1:B:35:GLU:H	1:B:395:MET:HE2	1.86	0.40
1:C:202:ARG:HB2	1:C:234:TYR:HB3	2.04	0.40
1:D:216:MET:O	1:D:217:ALA:C	2.64	0.40
1:E:348:TRP:CH2	1:E:416:VAL:HG21	2.56	0.40
1:C:418:LEU:O	1:C:422:ILE:HG13	2.21	0.40
1:E:422:ILE:HG23	1:E:436:GLY:HA3	2.03	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	442/496 (89%)	433 (98%)	9 (2%)	0	100	100
1	B	443/496 (89%)	423 (96%)	20 (4%)	0	100	100
1	C	445/496 (90%)	437 (98%)	8 (2%)	0	100	100
1	D	447/496 (90%)	431 (96%)	16 (4%)	0	100	100
1	E	453/496 (91%)	439 (97%)	13 (3%)	1 (0%)	43	36
1	F	442/496 (89%)	429 (97%)	12 (3%)	1 (0%)	43	36
All	All	2672/2976 (90%)	2592 (97%)	78 (3%)	2 (0%)	48	40

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	128	ASP
1	E	265	LYS

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	334/406 (82%)	330 (99%)	4 (1%)	63	63
1	B	319/406 (79%)	317 (99%)	2 (1%)	78	81
1	C	336/406 (83%)	334 (99%)	2 (1%)	78	81
1	D	354/406 (87%)	352 (99%)	2 (1%)	78	81
1	E	347/406 (86%)	345 (99%)	2 (1%)	78	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	340/406 (84%)	339 (100%)	1 (0%)	86	88
All	All	2030/2436 (83%)	2017 (99%)	13 (1%)	78	81

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	107	SER
1	A	173	LEU
1	A	241	THR
1	A	377	PHE
1	B	173	LEU
1	B	190	SER
1	C	241	THR
1	C	464	VAL
1	D	241	THR
1	D	470	LYS
1	E	241	THR
1	E	380	SER
1	F	456	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	80	ASN
1	A	95	GLN
1	A	374	ASN
1	B	293	ASN
1	B	440	ASN
1	C	88	GLN
1	C	267	GLN
1	C	426	HIS
1	D	9	ASN
1	D	20	HIS
1	D	77	GLN
1	D	80	ASN
1	D	317	GLN
1	D	423	GLN
1	E	80	ASN
1	E	267	GLN
1	F	20	HIS

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Mol	Chain	Res	Type
1	F	80	ASN
1	F	226	GLN
1	F	305	GLN
1	F	407	GLN
1	F	434	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
1	LLP	E	244[A]	1	23,24,25	2.50	6 (26%)	25,32,34	1.57	6 (24%)
1	LLP	C	244[A]	1	23,24,25	2.54	6 (26%)	25,32,34	1.29	3 (12%)
1	LLP	B	244[B]	1	7,8,25	0.69	0	3,8,34	0.62	0
1	LLP	D	244[A]	1	23,24,25	2.57	7 (30%)	25,32,34	1.48	6 (24%)
1	LLP	F	244[B]	1	7,8,25	0.66	0	3,8,34	0.58	0
1	LLP	F	244[A]	1	23,24,25	2.63	6 (26%)	25,32,34	1.40	4 (16%)
1	LLP	B	244[A]	1	23,24,25	2.59	6 (26%)	25,32,34	1.60	6 (24%)
1	LLP	A	244[B]	1	7,8,25	0.73	0	3,8,34	0.56	0
1	LLP	E	244[B]	1	7,8,25	0.69	0	3,8,34	0.66	0
1	LLP	C	244[B]	1	7,8,25	0.65	0	3,8,34	0.60	0
1	LLP	D	244[B]	1	7,8,25	0.72	0	3,8,34	0.50	0
1	LLP	A	244[A]	1	23,24,25	2.55	6 (26%)	25,32,34	1.55	5 (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
1	LLP	E	244[A]	1	-	7/16/17/19	0/1/1/1
1	LLP	C	244[A]	1	-	4/16/17/19	0/1/1/1
1	LLP	B	244[B]	1	-	2/6/7/19	-
1	LLP	D	244[A]	1	-	4/16/17/19	0/1/1/1
1	LLP	F	244[B]	1	-	1/6/7/19	-
1	LLP	F	244[A]	1	-	2/16/17/19	0/1/1/1
1	LLP	B	244[A]	1	-	6/16/17/19	0/1/1/1
1	LLP	A	244[B]	1	-	1/6/7/19	-
1	LLP	E	244[B]	1	-	2/6/7/19	-
1	LLP	C	244[B]	1	-	2/6/7/19	-
1	LLP	D	244[B]	1	-	1/6/7/19	-
1	LLP	A	244[A]	1	-	7/16/17/19	0/1/1/1

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	244[A]	LLP	C4-C4'	7.77	1.63	1.46
1	D	244[A]	LLP	C4-C4'	7.52	1.62	1.46
1	C	244[A]	LLP	C4-C4'	7.52	1.62	1.46
1	E	244[A]	LLP	C4-C4'	7.39	1.62	1.46
1	B	244[A]	LLP	C4-C4'	7.14	1.61	1.46
1	A	244[A]	LLP	C4-C4'	7.06	1.61	1.46
1	B	244[A]	LLP	C4'-NZ	5.05	1.44	1.27
1	F	244[A]	LLP	C4'-NZ	5.04	1.44	1.27
1	E	244[A]	LLP	C4'-NZ	4.96	1.43	1.27
1	C	244[A]	LLP	C4'-NZ	4.89	1.43	1.27
1	A	244[A]	LLP	C4'-NZ	4.86	1.43	1.27
1	D	244[A]	LLP	C4'-NZ	4.71	1.42	1.27
1	B	244[A]	LLP	C4-C5	-4.22	1.36	1.42
1	D	244[A]	LLP	C4-C5	-4.15	1.36	1.42
1	A	244[A]	LLP	C4-C5	-4.08	1.36	1.42
1	B	244[A]	LLP	C2'-C2	4.08	1.56	1.50
1	D	244[A]	LLP	C2'-C2	3.88	1.56	1.50
1	F	244[A]	LLP	C4-C5	-3.83	1.36	1.42
1	F	244[A]	LLP	C2'-C2	3.80	1.56	1.50
1	A	244[A]	LLP	C2'-C2	3.79	1.56	1.50
1	E	244[A]	LLP	C4-C5	-3.67	1.36	1.42
1	C	244[A]	LLP	C4-C5	-3.65	1.36	1.42
1	C	244[A]	LLP	C2'-C2	3.65	1.56	1.50
1	E	244[A]	LLP	C2'-C2	3.59	1.56	1.50
1	D	244[A]	LLP	C6-N1	3.56	1.41	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	244[A]	LLP	C6-N1	3.55	1.41	1.34
1	F	244[A]	LLP	C6-N1	3.50	1.41	1.34
1	A	244[A]	LLP	C6-N1	3.40	1.41	1.34
1	E	244[A]	LLP	C6-N1	3.32	1.41	1.34
1	C	244[A]	LLP	C6-N1	3.26	1.41	1.34
1	C	244[A]	LLP	C3-C2	2.37	1.43	1.41
1	F	244[A]	LLP	C3-C2	2.21	1.43	1.41
1	E	244[A]	LLP	C6-C5	2.08	1.41	1.37
1	A	244[A]	LLP	C5'-C5	2.07	1.56	1.50
1	B	244[A]	LLP	O3-C3	2.06	1.41	1.36
1	D	244[A]	LLP	C6-C5	2.06	1.41	1.37
1	D	244[A]	LLP	C5'-C5	2.03	1.56	1.50

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	244[A]	LLP	C4-C4'-NZ	-3.54	107.70	124.04
1	B	244[A]	LLP	C4-C4'-NZ	-3.32	108.74	124.04
1	D	244[A]	LLP	C4-C4'-NZ	-3.16	109.46	124.04
1	D	244[A]	LLP	CE-NZ-C4'	-3.16	108.61	118.72
1	C	244[A]	LLP	C4-C4'-NZ	-3.15	109.51	124.04
1	E	244[A]	LLP	C2'-C2-C3	-3.05	117.23	120.80
1	A	244[A]	LLP	CE-NZ-C4'	-2.95	109.27	118.72
1	F	244[A]	LLP	C4-C4'-NZ	-2.95	110.44	124.04
1	E	244[A]	LLP	C4-C4'-NZ	-2.92	110.59	124.04
1	F	244[A]	LLP	C5-C6-N1	-2.89	119.12	123.83
1	B	244[A]	LLP	CE-NZ-C4'	-2.88	109.48	118.72
1	C	244[A]	LLP	CE-NZ-C4'	-2.86	109.56	118.72
1	B	244[A]	LLP	C5-C6-N1	-2.80	119.28	123.83
1	B	244[A]	LLP	C3-C4-C5	2.79	120.52	118.28
1	A	244[A]	LLP	OP4-P-OP1	2.72	113.79	106.44
1	E	244[A]	LLP	CE-NZ-C4'	-2.68	110.12	118.72
1	E	244[A]	LLP	C2'-C2-N1	2.68	122.69	117.64
1	F	244[A]	LLP	CE-NZ-C4'	-2.63	110.31	118.72
1	A	244[A]	LLP	C3-C4-C5	2.57	120.34	118.28
1	A	244[A]	LLP	C5-C6-N1	-2.52	119.74	123.83
1	E	244[A]	LLP	C5-C6-N1	-2.49	119.79	123.83
1	C	244[A]	LLP	C5-C6-N1	-2.46	119.83	123.83
1	D	244[A]	LLP	C5-C6-N1	-2.42	119.90	123.83
1	B	244[A]	LLP	C2'-C2-N1	2.36	122.08	117.64
1	D	244[A]	LLP	C2'-C2-N1	2.28	121.95	117.64
1	B	244[A]	LLP	C2'-C2-C3	-2.23	118.18	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	244[A]	LLP	C4-C3-C2	2.23	121.40	120.14
1	D	244[A]	LLP	C2'-C2-C3	-2.20	118.22	120.80
1	D	244[A]	LLP	OP3-P-OP4	2.16	112.30	106.67
1	F	244[A]	LLP	C2'-C2-N1	2.01	121.42	117.64

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	244[A]	LLP	C5'-OP4-P-OP2
1	A	244[A]	LLP	C5'-OP4-P-OP3
1	A	244[A]	LLP	O-C-CA-CB
1	A	244[B]	LLP	O-C-CA-CB
1	B	244[A]	LLP	C5'-OP4-P-OP1
1	B	244[A]	LLP	C5'-OP4-P-OP2
1	B	244[A]	LLP	C5'-OP4-P-OP3
1	B	244[A]	LLP	C-CA-CB-CG
1	B	244[A]	LLP	O-C-CA-CB
1	B	244[B]	LLP	O-C-CA-CB
1	C	244[A]	LLP	O-C-CA-CB
1	C	244[B]	LLP	O-C-CA-CB
1	D	244[A]	LLP	O-C-CA-CB
1	D	244[B]	LLP	O-C-CA-CB
1	E	244[A]	LLP	C5'-OP4-P-OP1
1	E	244[A]	LLP	C5'-OP4-P-OP2
1	E	244[A]	LLP	C5'-OP4-P-OP3
1	E	244[A]	LLP	O-C-CA-CB
1	E	244[A]	LLP	CG-CD-CE-NZ
1	E	244[B]	LLP	O-C-CA-CB
1	F	244[A]	LLP	O-C-CA-CB
1	F	244[B]	LLP	O-C-CA-CB
1	A	244[A]	LLP	C4-C4'-NZ-CE
1	C	244[A]	LLP	CG-CD-CE-NZ
1	D	244[A]	LLP	CG-CD-CE-NZ
1	A	244[A]	LLP	C3-C4-C4'-NZ
1	A	244[A]	LLP	C5'-OP4-P-OP1
1	C	244[A]	LLP	C3-C4-C4'-NZ
1	B	244[B]	LLP	C-CA-CB-CG
1	D	244[A]	LLP	C3-C4-C4'-NZ
1	F	244[A]	LLP	C3-C4-C4'-NZ
1	E	244[A]	LLP	C3-C4-C4'-NZ
1	A	244[A]	LLP	C5-C4-C4'-NZ

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Mol	Chain	Res	Type	Atoms
1	E	244[A]	LLP	CD-CE-NZ-C4'
1	D	244[A]	LLP	C5'-OP4-P-OP3
1	C	244[B]	LLP	CG-CD-CE-NZ
1	C	244[A]	LLP	C5-C4-C4'-NZ
1	B	244[A]	LLP	C3-C4-C4'-NZ
1	E	244[B]	LLP	C-CA-CB-CG

There are no ring outliers.

12 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	E	244[A]	LLP	1	0
1	C	244[A]	LLP	2	0
1	B	244[B]	LLP	1	0
1	D	244[A]	LLP	1	0
1	F	244[B]	LLP	1	0
1	F	244[A]	LLP	1	0
1	B	244[A]	LLP	4	0
1	A	244[B]	LLP	1	0
1	E	244[B]	LLP	2	0
1	C	244[B]	LLP	1	0
1	D	244[B]	LLP	1	0
1	A	244[A]	LLP	3	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	450/496 (90%)	0.68	39 (8%) 16 17	23, 38, 63, 82	0
1	B	450/496 (90%)	1.27	87 (19%) 3 3	25, 47, 72, 81	1 (0%)
1	C	451/496 (90%)	0.75	36 (7%) 18 19	24, 38, 67, 82	0
1	D	453/496 (91%)	0.50	27 (5%) 27 29	22, 35, 56, 80	0
1	E	457/496 (92%)	0.63	27 (5%) 28 30	18, 34, 56, 83	2 (0%)
1	F	449/496 (90%)	0.62	26 (5%) 29 30	21, 37, 62, 75	1 (0%)
All	All	2710/2976 (91%)	0.74	242 (8%) 15 16	18, 38, 65, 83	4 (0%)

All (242) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	383	LEU	6.6
1	A	422	ILE	4.9
1	F	377	PHE	4.7
1	D	147	ILE	4.6
1	E	129	LEU	4.5
1	B	383	LEU	4.4
1	B	129	LEU	4.4
1	F	69	TYR	4.4
1	C	69	TYR	4.3
1	B	438	VAL	4.3
1	D	264	LYS	4.3
1	B	422	ILE	4.2
1	A	-1	SER	4.2
1	C	147	ILE	4.2
1	B	329	LEU	4.2
1	C	68	TYR	4.1
1	C	438	VAL	4.1
1	A	68	TYR	4.1
1	B	272	VAL	4.0

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Mol	Chain	Res	Type	RSRZ
1	F	0	ASN	4.0
1	B	377	PHE	4.0
1	B	68	TYR	3.9
1	A	130	ARG	3.9
1	A	429	LEU	3.8
1	B	437	LEU	3.8
1	E	0	ASN	3.7
1	B	0	ASN	3.6
1	B	439	ASN	3.6
1	A	67	ARG	3.6
1	F	420	LEU	3.6
1	F	422	ILE	3.6
1	B	420	LEU	3.6
1	C	129	LEU	3.6
1	D	384	ALA	3.5
1	B	350	LEU	3.5
1	C	437	LEU	3.5
1	E	270	ASN	3.5
1	A	433	PHE	3.5
1	A	355	LEU	3.5
1	F	129	LEU	3.5
1	B	69	TYR	3.4
1	E	147	ILE	3.4
1	D	272	VAL	3.4
1	B	399	GLY	3.4
1	F	130	ARG	3.4
1	B	443	ILE	3.4
1	F	264	LYS	3.4
1	A	437	LEU	3.4
1	B	263	PRO	3.4
1	C	432	ASP	3.3
1	E	69	TYR	3.3
1	A	0	ASN	3.3
1	C	164	THR	3.3
1	A	436	GLY	3.3
1	B	67	ARG	3.3
1	B	271	ALA	3.2
1	C	429	LEU	3.2
1	D	383	LEU	3.2
1	A	66	ASN	3.2
1	E	377	PHE	3.2
1	A	438	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	325	LEU	3.1
1	C	436	GLY	3.1
1	F	147	ILE	3.1
1	B	161	VAL	3.1
1	D	270	ASN	3.1
1	F	367	CYS	3.1
1	B	147	ILE	3.0
1	B	418	LEU	3.0
1	F	68	TYR	3.0
1	B	348	TRP	3.0
1	A	377	PHE	3.0
1	C	377	PHE	3.0
1	C	67	ARG	3.0
1	A	6	VAL	3.0
1	D	69	TYR	3.0
1	A	7	TRP	3.0
1	E	265	LYS	3.0
1	D	439	ASN	3.0
1	B	265	LYS	2.9
1	F	438	VAL	2.9
1	F	66	ASN	2.9
1	B	411	PHE	2.9
1	C	384	ALA	2.9
1	A	424	LYS	2.9
1	B	408	ILE	2.9
1	B	363	LEU	2.9
1	B	400	LEU	2.9
1	D	271	ALA	2.9
1	B	196	TRP	2.9
1	C	439	ASN	2.8
1	A	432	ASP	2.8
1	E	462	PHE	2.8
1	C	130	ARG	2.8
1	C	430	LEU	2.8
1	E	132	GLY	2.8
1	E	378	GLY	2.8
1	F	268	PRO	2.8
1	B	225	ALA	2.7
1	A	439	ASN	2.7
1	D	269	GLU	2.7
1	D	382	ALA	2.7
1	B	427	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	266	GLY	2.7
1	A	363	LEU	2.7
1	E	439	ASN	2.7
1	E	271	ALA	2.7
1	D	95	GLN	2.7
1	B	450	VAL	2.7
1	F	384	ALA	2.6
1	B	454	SER	2.6
1	D	268	PRO	2.6
1	B	75	ILE	2.6
1	C	66	ASN	2.6
1	E	430	LEU	2.6
1	A	147	ILE	2.6
1	B	65	GLY	2.6
1	C	426	HIS	2.6
1	A	3	PRO	2.6
1	D	281	PHE	2.6
1	B	130	ARG	2.6
1	B	47	VAL	2.6
1	E	382	ALA	2.6
1	B	428	LYS	2.5
1	E	266	GLY	2.5
1	E	263	PRO	2.5
1	C	433	PHE	2.5
1	F	146	LYS	2.5
1	B	331	GLY	2.5
1	E	94	ALA	2.5
1	A	1	MET	2.5
1	C	262	PRO	2.5
1	C	263	PRO	2.5
1	A	2	ASP	2.5
1	A	129	LEU	2.5
1	B	347	LEU	2.5
1	B	355	LEU	2.5
1	B	453	PHE	2.5
1	A	434	ASN	2.5
1	E	75	ILE	2.5
1	E	467	MET	2.5
1	B	262	PRO	2.5
1	F	67	ARG	2.4
1	A	420	LEU	2.4
1	B	405	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	457	PHE	2.4
1	A	425	GLU	2.4
1	B	269	GLU	2.4
1	D	-1	SER	2.4
1	C	0	ASN	2.4
1	D	65	GLY	2.4
1	B	178	LEU	2.4
1	B	430	LEU	2.4
1	B	168	ILE	2.4
1	D	267	GLN	2.4
1	F	427	GLY	2.4
1	B	429	LEU	2.4
1	B	237	ILE	2.4
1	C	354	GLY	2.4
1	B	92	LEU	2.3
1	F	437	LEU	2.3
1	A	358	TYR	2.3
1	D	467	MET	2.3
1	A	431	LYS	2.3
1	B	346	VAL	2.3
1	F	269	GLU	2.3
1	A	418	LEU	2.3
1	F	366	LEU	2.3
1	F	430	LEU	2.3
1	D	68	TYR	2.3
1	B	80	ASN	2.3
1	C	442	ALA	2.3
1	B	34	ILE	2.3
1	C	189	GLY	2.3
1	B	131	SER	2.3
1	D	67	ARG	2.3
1	B	463	LEU	2.3
1	D	0	ASN	2.3
1	E	66	ASN	2.3
1	B	384	ALA	2.3
1	F	432	ASP	2.3
1	E	6	VAL	2.3
1	F	270	ASN	2.3
1	B	327	LYS	2.3
1	B	446	LEU	2.3
1	C	355	LEU	2.3
1	F	433	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	404	ASP	2.2
1	A	426	HIS	2.2
1	B	367	CYS	2.2
1	B	456	LEU	2.2
1	B	190	SER	2.2
1	B	268	PRO	2.2
1	E	133	GLY	2.2
1	E	471	ASP	2.2
1	B	97	TRP	2.2
1	B	358	TYR	2.2
1	C	281	PHE	2.2
1	A	146	LYS	2.2
1	C	431	LYS	2.2
1	B	401	VAL	2.2
1	E	272	VAL	2.2
1	A	274	ASP	2.2
1	B	421	GLU	2.2
1	A	4	VAL	2.2
1	E	438	VAL	2.2
1	A	430	LEU	2.2
1	C	420	LEU	2.2
1	B	442	ALA	2.2
1	B	455	ALA	2.2
1	E	384	ALA	2.2
1	D	377	PHE	2.1
1	D	189	GLY	2.1
1	B	260	PRO	2.1
1	D	262	PRO	2.1
1	A	270	ASN	2.1
1	B	281	PHE	2.1
1	B	328	TYR	2.1
1	D	273	TYR	2.1
1	A	427	GLY	2.1
1	B	451	GLU	2.1
1	B	81	LEU	2.1
1	C	268	PRO	2.1
1	D	129	LEU	2.1
1	B	392	ALA	2.1
1	B	332	LYS	2.1
1	B	410	GLU	2.1
1	B	318	VAL	2.1
1	C	270	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	423	GLN	2.1
1	D	263	PRO	2.1
1	B	287	LEU	2.1
1	B	448	ALA	2.1
1	A	443	ILE	2.0
1	B	447	LYS	2.0
1	E	68	TYR	2.0
1	C	188	GLY	2.0
1	B	440	ASN	2.0
1	F	355	LEU	2.0
1	C	-1	SER	2.0
1	B	469	TYR	2.0
1	C	464	VAL	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	LLP	B	244[A]	24/25	0.80	0.16	37,51,59,66	24
1	LLP	B	244[B]	9/25	0.80	0.16	37,39,46,50	9
1	LLP	F	244[A]	24/25	0.87	0.15	28,44,57,60	24
1	LLP	F	244[B]	9/25	0.87	0.15	28,29,37,42	9
1	LLP	C	244[A]	24/25	0.89	0.13	28,40,51,52	24
1	LLP	C	244[B]	9/25	0.89	0.13	28,30,36,39	9
1	LLP	E	244[A]	24/25	0.89	0.12	25,38,47,53	24
1	LLP	E	244[B]	9/25	0.89	0.12	24,28,35,37	9
1	LLP	A	244[A]	24/25	0.89	0.13	27,42,51,53	24
1	LLP	A	244[B]	9/25	0.89	0.13	29,32,36,39	9
1	LLP	D	244[A]	24/25	0.91	0.11	25,37,45,48	24
1	LLP	D	244[B]	9/25	0.91	0.11	24,28,34,35	9

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.