



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:00 AM UTC

PDB ID : 3WKR / pdb_00003wkr
Title : Crystal structure of the SepCysS-SepCysE complex from *Methanocaldococcus jannaschii*
Authors : Nakazawa, Y.; Asano, N.; Nakamura, A.; Yao, M.
Deposited on : 2013-10-30
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

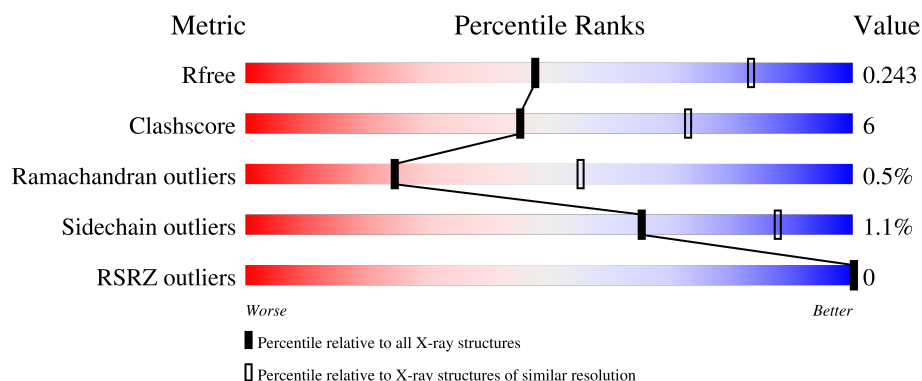
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	416	
1	B	416	
1	E	416	
1	F	416	
2	C	216	

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Mol	Chain	Length	Quality of chain
2	D	216	 28% 69%
2	G	216	 29% 68%
2	H	216	 29% 69%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14030 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 O-phospho-L-seryl-tRNA:Cys-tRNA synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	P	S	0	0	0
			2903	1865	485	539	1	13			
1	B	364	Total	C	N	O	P	S	0	0	0
			2916	1873	487	541	1	14			
1	E	365	Total	C	N	O	P	S	0	0	0
			2923	1877	488	543	1	14			
1	F	363	Total	C	N	O	P	S	0	0	0
			2910	1870	486	540	1	13			

There are 160 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q59072
A	-18	GLY	-	expression tag	UNP Q59072
A	-17	SER	-	expression tag	UNP Q59072
A	-16	SER	-	expression tag	UNP Q59072
A	-15	HIS	-	expression tag	UNP Q59072
A	-14	HIS	-	expression tag	UNP Q59072
A	-13	HIS	-	expression tag	UNP Q59072
A	-12	HIS	-	expression tag	UNP Q59072
A	-11	HIS	-	expression tag	UNP Q59072
A	-10	HIS	-	expression tag	UNP Q59072
A	-9	SER	-	expression tag	UNP Q59072
A	-8	SER	-	expression tag	UNP Q59072
A	-7	GLY	-	expression tag	UNP Q59072
A	-6	LEU	-	expression tag	UNP Q59072
A	-5	VAL	-	expression tag	UNP Q59072
A	-4	PRO	-	expression tag	UNP Q59072
A	-3	ARG	-	expression tag	UNP Q59072
A	-2	GLY	-	expression tag	UNP Q59072
A	-1	SER	-	expression tag	UNP Q59072
A	0	HIS	-	expression tag	UNP Q59072
A	1	MET	-	expression tag	UNP Q59072

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Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLU	-	expression tag	UNP Q59072
A	3	LEU	-	expression tag	UNP Q59072
A	4	GLU	-	expression tag	UNP Q59072
A	5	GLY	-	expression tag	UNP Q59072
A	6	PRO	-	expression tag	UNP Q59072
A	7	TYR	-	expression tag	UNP Q59072
A	8	SER	-	expression tag	UNP Q59072
A	9	LYS	-	expression tag	UNP Q59072
A	10	LYS	-	expression tag	UNP Q59072
A	11	PHE	-	expression tag	UNP Q59072
A	12	GLU	-	expression tag	UNP Q59072
A	13	VAL	-	expression tag	UNP Q59072
A	14	ILE	-	expression tag	UNP Q59072
A	15	THR	-	expression tag	UNP Q59072
A	16	LEU	-	expression tag	UNP Q59072
A	17	ASP	-	expression tag	UNP Q59072
A	18	ILE	-	expression tag	UNP Q59072
A	19	ASN	-	expression tag	UNP Q59072
A	20	LEU	-	expression tag	UNP Q59072
B	-19	MET	-	expression tag	UNP Q59072
B	-18	GLY	-	expression tag	UNP Q59072
B	-17	SER	-	expression tag	UNP Q59072
B	-16	SER	-	expression tag	UNP Q59072
B	-15	HIS	-	expression tag	UNP Q59072
B	-14	HIS	-	expression tag	UNP Q59072
B	-13	HIS	-	expression tag	UNP Q59072
B	-12	HIS	-	expression tag	UNP Q59072
B	-11	HIS	-	expression tag	UNP Q59072
B	-10	HIS	-	expression tag	UNP Q59072
B	-9	SER	-	expression tag	UNP Q59072
B	-8	SER	-	expression tag	UNP Q59072
B	-7	GLY	-	expression tag	UNP Q59072
B	-6	LEU	-	expression tag	UNP Q59072
B	-5	VAL	-	expression tag	UNP Q59072
B	-4	PRO	-	expression tag	UNP Q59072
B	-3	ARG	-	expression tag	UNP Q59072
B	-2	GLY	-	expression tag	UNP Q59072
B	-1	SER	-	expression tag	UNP Q59072
B	0	HIS	-	expression tag	UNP Q59072
B	1	MET	-	expression tag	UNP Q59072
B	2	GLU	-	expression tag	UNP Q59072
B	3	LEU	-	expression tag	UNP Q59072

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Chain	Residue	Modelled	Actual	Comment	Reference
B	4	GLU	-	expression tag	UNP Q59072
B	5	GLY	-	expression tag	UNP Q59072
B	6	PRO	-	expression tag	UNP Q59072
B	7	TYR	-	expression tag	UNP Q59072
B	8	SER	-	expression tag	UNP Q59072
B	9	LYS	-	expression tag	UNP Q59072
B	10	LYS	-	expression tag	UNP Q59072
B	11	PHE	-	expression tag	UNP Q59072
B	12	GLU	-	expression tag	UNP Q59072
B	13	VAL	-	expression tag	UNP Q59072
B	14	ILE	-	expression tag	UNP Q59072
B	15	THR	-	expression tag	UNP Q59072
B	16	LEU	-	expression tag	UNP Q59072
B	17	ASP	-	expression tag	UNP Q59072
B	18	ILE	-	expression tag	UNP Q59072
B	19	ASN	-	expression tag	UNP Q59072
B	20	LEU	-	expression tag	UNP Q59072
E	-19	MET	-	expression tag	UNP Q59072
E	-18	GLY	-	expression tag	UNP Q59072
E	-17	SER	-	expression tag	UNP Q59072
E	-16	SER	-	expression tag	UNP Q59072
E	-15	HIS	-	expression tag	UNP Q59072
E	-14	HIS	-	expression tag	UNP Q59072
E	-13	HIS	-	expression tag	UNP Q59072
E	-12	HIS	-	expression tag	UNP Q59072
E	-11	HIS	-	expression tag	UNP Q59072
E	-10	HIS	-	expression tag	UNP Q59072
E	-9	SER	-	expression tag	UNP Q59072
E	-8	SER	-	expression tag	UNP Q59072
E	-7	GLY	-	expression tag	UNP Q59072
E	-6	LEU	-	expression tag	UNP Q59072
E	-5	VAL	-	expression tag	UNP Q59072
E	-4	PRO	-	expression tag	UNP Q59072
E	-3	ARG	-	expression tag	UNP Q59072
E	-2	GLY	-	expression tag	UNP Q59072
E	-1	SER	-	expression tag	UNP Q59072
E	0	HIS	-	expression tag	UNP Q59072
E	1	MET	-	expression tag	UNP Q59072
E	2	GLU	-	expression tag	UNP Q59072
E	3	LEU	-	expression tag	UNP Q59072
E	4	GLU	-	expression tag	UNP Q59072
E	5	GLY	-	expression tag	UNP Q59072

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Chain	Residue	Modelled	Actual	Comment	Reference
E	6	PRO	-	expression tag	UNP Q59072
E	7	TYR	-	expression tag	UNP Q59072
E	8	SER	-	expression tag	UNP Q59072
E	9	LYS	-	expression tag	UNP Q59072
E	10	LYS	-	expression tag	UNP Q59072
E	11	PHE	-	expression tag	UNP Q59072
E	12	GLU	-	expression tag	UNP Q59072
E	13	VAL	-	expression tag	UNP Q59072
E	14	ILE	-	expression tag	UNP Q59072
E	15	THR	-	expression tag	UNP Q59072
E	16	LEU	-	expression tag	UNP Q59072
E	17	ASP	-	expression tag	UNP Q59072
E	18	ILE	-	expression tag	UNP Q59072
E	19	ASN	-	expression tag	UNP Q59072
E	20	LEU	-	expression tag	UNP Q59072
F	-19	MET	-	expression tag	UNP Q59072
F	-18	GLY	-	expression tag	UNP Q59072
F	-17	SER	-	expression tag	UNP Q59072
F	-16	SER	-	expression tag	UNP Q59072
F	-15	HIS	-	expression tag	UNP Q59072
F	-14	HIS	-	expression tag	UNP Q59072
F	-13	HIS	-	expression tag	UNP Q59072
F	-12	HIS	-	expression tag	UNP Q59072
F	-11	HIS	-	expression tag	UNP Q59072
F	-10	HIS	-	expression tag	UNP Q59072
F	-9	SER	-	expression tag	UNP Q59072
F	-8	SER	-	expression tag	UNP Q59072
F	-7	GLY	-	expression tag	UNP Q59072
F	-6	LEU	-	expression tag	UNP Q59072
F	-5	VAL	-	expression tag	UNP Q59072
F	-4	PRO	-	expression tag	UNP Q59072
F	-3	ARG	-	expression tag	UNP Q59072
F	-2	GLY	-	expression tag	UNP Q59072
F	-1	SER	-	expression tag	UNP Q59072
F	0	HIS	-	expression tag	UNP Q59072
F	1	MET	-	expression tag	UNP Q59072
F	2	GLU	-	expression tag	UNP Q59072
F	3	LEU	-	expression tag	UNP Q59072
F	4	GLU	-	expression tag	UNP Q59072
F	5	GLY	-	expression tag	UNP Q59072
F	6	PRO	-	expression tag	UNP Q59072
F	7	TYR	-	expression tag	UNP Q59072

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Chain	Residue	Modelled	Actual	Comment	Reference
F	8	SER	-	expression tag	UNP Q59072
F	9	LYS	-	expression tag	UNP Q59072
F	10	LYS	-	expression tag	UNP Q59072
F	11	PHE	-	expression tag	UNP Q59072
F	12	GLU	-	expression tag	UNP Q59072
F	13	VAL	-	expression tag	UNP Q59072
F	14	ILE	-	expression tag	UNP Q59072
F	15	THR	-	expression tag	UNP Q59072
F	16	LEU	-	expression tag	UNP Q59072
F	17	ASP	-	expression tag	UNP Q59072
F	18	ILE	-	expression tag	UNP Q59072
F	19	ASN	-	expression tag	UNP Q59072
F	20	LEU	-	expression tag	UNP Q59072

- Molecule 2 is a protein called Uncharacterized protein MJ1481.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	68	Total	C	N	O	S	0	0	0
			562	368	89	103	2			
2	D	68	Total	C	N	O	S	0	0	0
			565	371	89	103	2			
2	G	69	Total	C	N	O	S	0	0	0
			573	377	90	104	2			
2	H	68	Total	C	N	O	S	0	0	0
			565	371	89	103	2			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	MET	-	expression tag	UNP Q58876
C	-1	ASN	-	expression tag	UNP Q58876
C	0	HIS	-	expression tag	UNP Q58876
D	-2	MET	-	expression tag	UNP Q58876
D	-1	ASN	-	expression tag	UNP Q58876
D	0	HIS	-	expression tag	UNP Q58876
G	-2	MET	-	expression tag	UNP Q58876
G	-1	ASN	-	expression tag	UNP Q58876
G	0	HIS	-	expression tag	UNP Q58876
H	-2	MET	-	expression tag	UNP Q58876
H	-1	ASN	-	expression tag	UNP Q58876
H	0	HIS	-	expression tag	UNP Q58876

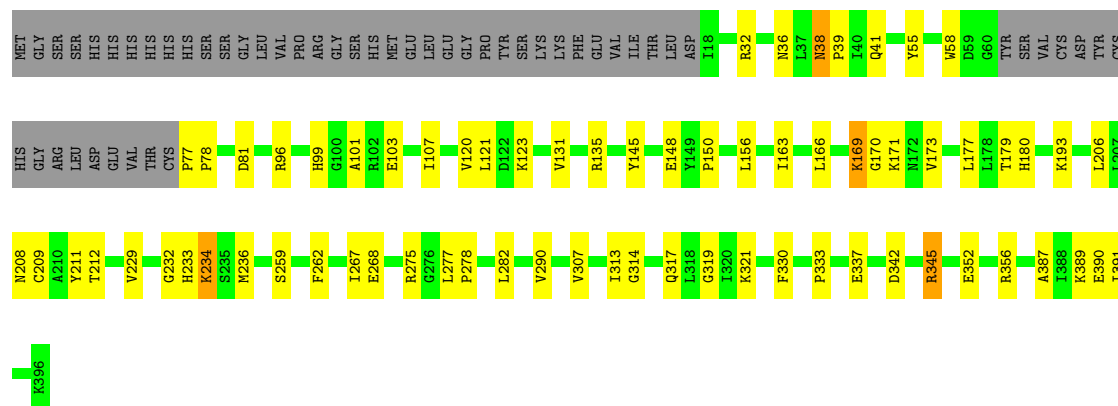
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	24	Total 24	O 24	0	0
3	B	26	Total 26	O 26	0	0
3	C	4	Total 4	O 4	0	0
3	D	4	Total 4	O 4	0	0
3	E	22	Total 22	O 22	0	0
3	F	20	Total 20	O 20	0	0
3	G	6	Total 6	O 6	0	0
3	H	7	Total 7	O 7	0	0



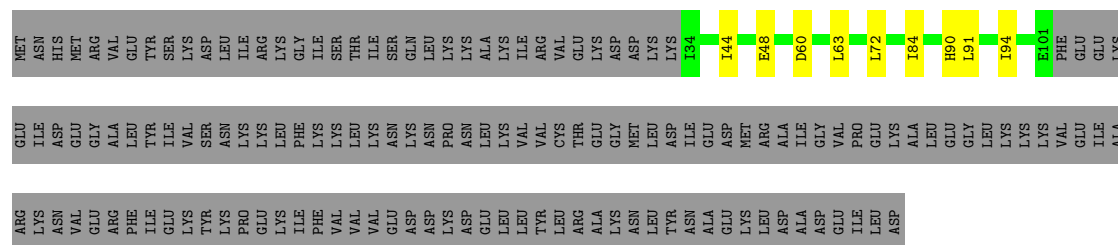
- Molecule 1: O-phospho-L-seryl-tRNA:Cys-tRNA synthase

Chain F: 70% 16% 13%



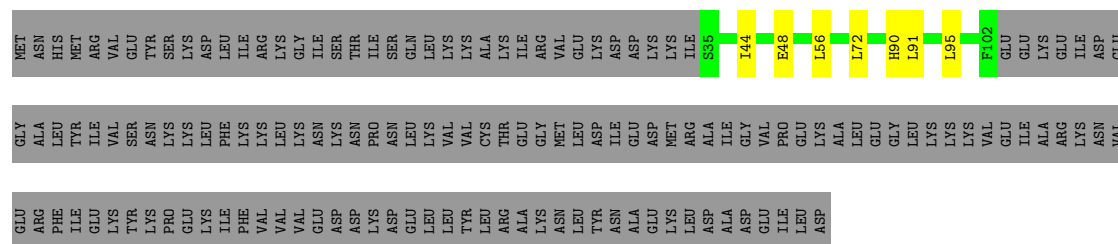
- Molecule 2: Uncharacterized protein MJ1481

Chain C: 27% 69%



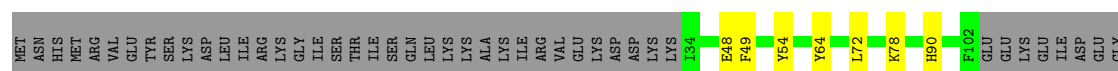
- Molecule 2: Uncharacterized protein MJ1481

Chain D: 28% 69%



- Molecule 2: Uncharacterized protein MJ1481

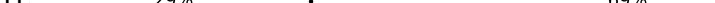
Chain G: 29% 68%



ALA LEU LEU TYR ILE VAL SER ASN LYS LYS LYS LYS PHE LYS LYS LYS LYS ASN ASN ASN ASN PRO THR GLU GLY MET MET LEU ASP THR ILE GLY GLY VAL VAL GLU LYS ALA ALA LEU GLU GLY LEU LYS LYS LYS VAL VAL ILE ALA ARG LYS LYS VAL VAL THR

ARG	PHE	ILE	GLU	LYS	TYR	LYS	PRO	GLU	LYS	ILE	PHE	VAL	VAL	VAL	GLU	ASP	ASP	LYS	LYS	ASP	GLU	LEU	LEU	TYR	LEU	ARG	ALA	ALA	LYS	ASN	LEU	LEU	TYR	ASN	ASN	GLU	LYS	LEU	ASP	ALA	ALA	ASP	GLU	ILE	LEU	LEU	ASP
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- Molecule 2: Uncharacterized protein MJ1481

Chain H:  29% . 69%

MET	ASIN	HIS	MET	ARG	VAL	GLU	TYR	SER	SER	LYS	ASP	LEU	LEU	ILE	ARG	LYS	GLY	ILE	ILE	SER	THR	ILE	SER	GLN	LYS	LYS	ALA	LYS	ILE	ILE	ARG	VAL	GLU	GLU	ASP	ASP	LYS	LYS	ILE	S35	Y36	K37		I44	E48		L63		H90	F102	GLU	GLU	LYS	LYS	GLU	ILE	ASP	GLU	GLY	ALA	LEU
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TYR	ILE	VAL	VAL	ASN	SER	ASN	LYS	LYS	LEU	LEU	LEU	LEU	ASN	ASN	ASN	ASN	ASN	LEU	LYS	VAL	VAL	CYS	THR	THR	GLU	GLY	MET	MET	LEU	ASP	ILE	GLU	ASP	MET	ARG	ALA	ILE	GLY	VAL	PRO	PRO	GLU	LYS	LYS	ALA	ALA	LEU	GLU	GLY	LEU	LEU	LYS	LYS	LYS	VAL	GLU	ILE	ILE	ALA	ARG	LYS	ASN	ASN	VAL	GLU	GLU	ARG	THR
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ILE	GLU	LYS	TYR	LYS	PRO	GLU	LYS	ILE	PHE	VAL	VAL	VAL	GLU	ASP	ASP	LYS	ASP	GLU	LEU	LEU	TYR	LEU	ARG	ALA	ALA	LYS	ASN	LEU	TYR	ASN	ALA	GLU	LYS	LEU	ASP	ALA	ASP	GLU	ILE	LEU	ASP
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4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	106.88Å 106.88Å 495.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.30 – 2.80 47.30 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.9 (47.30-2.80) 95.0 (47.30-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.7.2_869	Depositor
R, R_{free}	0.224 , 0.242 0.227 , 0.243	Depositor DCC
R_{free} test set	4279 reflections (5.46%)	wwPDB-VP
Wilson B-factor (Å ²)	71.2	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 100.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	0.467 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.498 for h,-h-k,-l	Depositor
Outliers	0 of 74017 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14030	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.17% 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: 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	0.25	0/2933	0.67	0/3939
1	B	0.25	0/2947	0.68	2/3960 (0.1%)
1	E	0.25	0/2954	0.67	0/3970
1	F	0.25	0/2941	0.68	4/3951 (0.1%)
2	C	0.22	0/573	0.60	0/765
2	D	0.22	0/577	0.57	0/770
2	G	0.22	0/585	0.59	0/781
2	H	0.22	0/577	0.57	0/770
All	All	0.25	0/14087	0.66	6/18906 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	169	LYS	N-CA-C	5.72	117.31	111.14
1	B	345	ARG	CB-CA-C	-5.50	110.22	116.54
1	F	345	ARG	CB-CA-C	-5.45	110.27	116.54
1	B	169	LYS	N-CA-C	5.38	117.96	111.40
1	F	38	ASN	CA-C-N	5.04	124.83	119.28
1	F	38	ASN	C-N-CA	5.04	124.83	119.28

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	2903	0	2951	37	0
1	B	2916	0	2962	43	0
1	E	2923	0	2969	46	0
1	F	2910	0	2958	45	0
2	C	562	0	571	7	0
2	D	565	0	569	5	0
2	G	573	0	580	6	0
2	H	565	0	569	4	0
3	A	24	0	0	1	0
3	B	26	0	0	0	0
3	C	4	0	0	0	0
3	D	4	0	0	0	0
3	E	22	0	0	0	0
3	F	20	0	0	0	0
3	G	6	0	0	0	0
3	H	7	0	0	0	0
All	All	14030	0	14129	165	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 6.

All (165) 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:101:ALA:HB3	1:B:234:LLP:H5'2	1.73	0.71
1:A:268:GLU:HA	1:B:135:ARG:HD2	1.75	0.69
1:B:232:GLY:HA2	1:B:236:MET:HB2	1.75	0.68
1:E:268:GLU:HA	1:F:135:ARG:HD2	1.77	0.66
1:E:55:TYR:O	1:F:32:ARG:NH1	2.26	0.66
1:E:135:ARG:HD2	1:F:268:GLU:HA	1.79	0.65
1:E:107:ILE:HG23	1:E:268:GLU:HG2	1.79	0.65
1:E:234:LLP:NZ	1:E:234:LLP:O3	2.27	0.65
1:A:352:GLU:HB3	1:A:391:ILE:HD12	1.79	0.65
1:A:55:TYR:O	1:B:32:ARG:NH1	2.28	0.64
1:E:340:LYS:HG2	1:E:341:LYS:HG3	1.79	0.63
1:A:318:LEU:HD11	1:A:331:GLU:HB2	1.80	0.63
1:A:99:HIS:HB2	1:A:103:GLU:HG3	1.79	0.63
1:A:232:GLY:HA2	1:A:236:MET:HB2	1.80	0.63
1:B:121:LEU:HB3	1:B:177:LEU:HB3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:352:GLU:HB3	1:F:391:ILE:HD12	1.81	0.62
1:A:340:LYS:HG2	1:A:341:LYS:HG3	1.81	0.62
1:B:179:THR:HA	1:B:208:ASN:HB3	1.81	0.62
1:B:234:LLP:O3	1:B:234:LLP:NZ	2.29	0.62
1:B:367:LYS:HG3	1:B:368:GLU:HG2	1.82	0.62
1:E:179:THR:HA	1:E:208:ASN:HB3	1.82	0.61
1:A:233:HIS:NE2	1:A:234:LLP:OP2	2.31	0.61
1:A:135:ARG:HD2	1:B:268:GLU:HA	1.83	0.60
1:A:234:LLP:NZ	1:A:234:LLP:O3	2.33	0.60
1:B:352:GLU:HB3	1:B:391:ILE:HD12	1.82	0.60
1:E:307:VAL:HA	1:E:330:PHE:HZ	1.67	0.60
1:B:242:CYS:HB2	1:B:278:PRO:HB2	1.84	0.59
1:A:40:ILE:HG13	1:A:41:GLN:HG3	1.84	0.59
1:A:179:THR:HA	1:A:208:ASN:HB3	1.84	0.59
1:E:205:PHE:HD2	1:E:225:ALA:HA	1.67	0.58
1:F:234:LLP:O3	1:F:234:LLP:NZ	2.37	0.58
1:E:179:THR:HG22	1:E:208:ASN:HD22	1.70	0.57
1:A:205:PHE:HD2	1:A:225:ALA:HA	1.69	0.57
1:B:169:LYS:N	1:B:170:GLY:HA2	2.18	0.57
1:F:356:ARG:NH2	1:F:390:GLU:OE1	2.38	0.57
1:A:114:LYS:NZ	1:F:342:ASP:O	2.30	0.56
1:E:123:LYS:NZ	1:E:142:GLU:OE1	2.38	0.56
1:B:166:LEU:O	1:B:171:LYS:N	2.37	0.56
1:E:105:LYS:HG2	1:E:206:LEU:HD21	1.89	0.55
1:F:101:ALA:HB3	1:F:234:LLP:H5'2	1.89	0.55
1:F:179:THR:HA	1:F:208:ASN:HB3	1.88	0.55
1:A:276:GLY:HA2	1:B:240:ALA:HB3	1.89	0.55
1:F:99:HIS:HB2	1:F:103:GLU:HG3	1.89	0.54
1:E:352:GLU:HB3	1:E:391:ILE:HD12	1.89	0.54
1:A:102:ARG:NH2	1:A:127:TYR:OH	2.35	0.53
1:E:318:LEU:HD11	1:E:331:GLU:HB2	1.89	0.53
1:A:307:VAL:HA	1:A:330:PHE:HZ	1.73	0.53
1:B:233:HIS:NE2	1:B:234:LLP:OP2	2.34	0.53
1:E:93:ASP:HB2	1:E:248:SER:HA	1.90	0.53
1:E:332:THR:OG1	1:E:366:THR:O	2.26	0.52
1:F:313:ILE:HD11	1:F:389:LYS:HG2	1.91	0.52
1:E:26:LEU:HD11	1:F:77:PRO:HG3	1.91	0.52
1:B:107:ILE:HG23	1:B:268:GLU:HG2	1.92	0.52
1:B:99:HIS:HB2	1:B:103:GLU:HG3	1.92	0.52
1:F:78:PRO:HB2	1:F:81:ASP:HB2	1.93	0.51
1:F:150:PRO:O	1:F:319:GLY:N	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:125:ALA:HA	1:E:179:THR:HG21	1.93	0.51
1:A:125:ALA:HA	1:A:179:THR:HG21	1.92	0.50
1:F:166:LEU:O	1:F:171:LYS:N	2.44	0.50
2:C:72:LEU:HD11	2:D:44:ILE:HD11	1.93	0.50
1:B:40:ILE:HG23	1:B:234:LLP:HE3	1.92	0.50
1:E:156:LEU:HD12	1:E:193:LYS:HD2	1.94	0.50
1:A:114:LYS:HG2	1:F:345:ARG:HH11	1.77	0.49
1:B:107:ILE:HD12	1:B:268:GLU:HG3	1.94	0.49
1:B:34:PHE:HE2	1:B:354:LYS:HA	1.77	0.49
2:G:72:LEU:HD11	2:H:44:ILE:HD11	1.95	0.49
1:A:192:LYS:NZ	3:A:402:HOH:O	2.43	0.49
1:E:239:SER:HB2	1:F:277:LEU:HD13	1.95	0.49
2:C:84:ILE:HB	2:D:95:LEU:HD21	1.95	0.48
1:E:99:HIS:HB2	1:E:103:GLU:HG3	1.96	0.48
1:B:342:ASP:O	1:E:114:LYS:NZ	2.39	0.48
1:F:232:GLY:HA2	1:F:236:MET:HB2	1.93	0.48
1:F:121:LEU:HB3	1:F:177:LEU:HB3	1.95	0.48
1:F:169:LYS:N	1:F:170:GLY:HA2	2.28	0.48
1:A:341:LYS:HA	1:A:342:ASP:HA	1.58	0.47
1:E:56:GLU:OE2	2:G:54:TYR:OH	2.31	0.47
1:A:277:LEU:N	1:A:278:PRO:HD2	2.30	0.47
1:E:32:ARG:NH2	2:G:64:TYR:OH	2.47	0.47
2:H:48:GLU:OE2	2:H:90:HIS:ND1	2.37	0.47
1:B:307:VAL:HA	1:B:330:PHE:HZ	1.79	0.47
1:A:123:LYS:NZ	1:A:142:GLU:OE1	2.48	0.47
1:B:110:HIS:ND1	1:B:268:GLU:OE2	2.48	0.47
1:A:306:VAL:HG21	1:A:373:VAL:HG11	1.97	0.46
2:C:60:ASP:HA	2:C:63:LEU:HD13	1.96	0.46
1:E:259:SER:OG	1:E:262:PHE:O	2.31	0.46
1:A:22:LYS:HB3	1:B:78:PRO:HG3	1.96	0.46
1:A:338:ILE:HG21	1:A:395:CYS:HB3	1.96	0.46
1:B:196:LYS:HD2	1:B:224:LYS:HE3	1.98	0.46
1:A:179:THR:HG22	1:A:208:ASN:HD22	1.81	0.45
1:A:186:GLY:O	1:A:323:LYS:NZ	2.49	0.45
1:B:148:GLU:O	1:B:152:TYR:N	2.37	0.45
1:F:156:LEU:HD12	1:F:193:LYS:HD2	1.98	0.45
1:F:233:HIS:NE2	1:F:234:LLP:OP2	2.43	0.45
1:F:208:ASN:HD22	1:F:229:VAL:HG23	1.82	0.45
1:E:289:VAL:HA	1:E:292:ARG:HG2	1.98	0.45
1:F:236:MET:HE2	1:F:282:LEU:HD11	1.98	0.45
1:A:156:LEU:HD12	1:A:193:LYS:HD2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:356:ARG:HE	1:E:391:ILE:HD11	1.82	0.45
1:E:107:ILE:HD12	1:E:268:GLU:HG3	1.99	0.44
1:F:36:ASN:ND2	1:F:41:GLN:OE1	2.50	0.44
1:B:235:SER:O	1:B:292:ARG:NH2	2.50	0.44
1:F:209:CYS:HB3	1:F:212:THR:HB	2.00	0.44
1:E:96:ARG:HG2	1:E:269:MET:HE3	1.99	0.44
1:E:96:ARG:NH2	1:E:266:GLU:OE2	2.51	0.44
2:G:48:GLU:OE2	2:G:90:HIS:ND1	2.37	0.44
1:E:277:LEU:N	1:E:278:PRO:HD2	2.33	0.44
1:A:121:LEU:HB3	1:A:177:LEU:HB3	1.99	0.43
1:E:103:GLU:O	1:E:107:ILE:HG12	2.18	0.43
1:E:180:HIS:ND1	1:E:212:THR:HG21	2.33	0.43
1:F:103:GLU:O	1:F:107:ILE:HG12	2.18	0.43
1:E:40:ILE:HG13	1:E:41:GLN:HG3	1.99	0.43
1:E:32:ARG:NH1	1:F:55:TYR:O	2.48	0.43
1:B:103:GLU:O	1:B:107:ILE:HG12	2.17	0.43
1:B:252:SER:O	1:B:256:THR:OG1	2.25	0.43
2:G:49:PHE:CD2	2:H:63:LEU:HD11	2.53	0.43
1:A:239:SER:HB2	1:B:277:LEU:HD13	1.99	0.43
1:B:356:ARG:HE	1:B:391:ILE:HD11	1.83	0.43
1:E:131:VAL:HG13	1:F:267:ILE:HA	2.01	0.43
1:B:259:SER:OG	1:B:262:PHE:O	2.32	0.43
1:E:177:LEU:HD12	1:E:206:LEU:HD23	2.01	0.43
1:E:267:ILE:HA	1:F:131:VAL:HG13	2.00	0.43
1:F:123:LYS:HD3	1:F:145:TYR:CE2	2.53	0.43
1:A:239:SER:O	1:A:278:PRO:HB3	2.19	0.43
1:B:205:PHE:HD2	1:B:225:ALA:HA	1.84	0.42
1:A:180:HIS:ND1	1:A:212:THR:HG21	2.32	0.42
1:E:122:ASP:HB2	1:E:154:ILE:HG13	2.01	0.42
1:B:317:GLN:NE2	1:B:321:LYS:HG2	2.34	0.42
2:C:91:LEU:HD23	2:C:94:ILE:HD12	2.02	0.42
1:F:107:ILE:HD12	1:F:268:GLU:HG3	2.00	0.42
1:F:307:VAL:HA	1:F:330:PHE:HZ	1.85	0.42
1:F:387:ALA:O	1:F:391:ILE:HG12	2.19	0.42
1:B:151:THR:O	1:B:319:GLY:HA2	2.20	0.42
2:C:84:ILE:HG23	2:D:91:LEU:HD22	2.01	0.42
1:E:58:TRP:HE1	1:F:32:ARG:HH22	1.68	0.42
1:F:38:ASN:HA	1:F:39:PRO:HD3	1.87	0.42
1:B:209:CYS:HB3	1:B:212:THR:HB	2.02	0.42
1:B:387:ALA:O	1:B:391:ILE:HG12	2.19	0.42
1:E:26:LEU:O	1:E:28:ARG:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:HIS:ND1	1:B:212:THR:HG21	2.34	0.42
2:C:48:GLU:OE2	2:C:90:HIS:ND1	2.42	0.42
1:E:130:TYR:CE2	1:E:140:VAL:HG11	2.55	0.42
1:E:341:LYS:HA	1:E:342:ASP:HA	1.65	0.42
1:F:180:HIS:ND1	1:F:212:THR:HG21	2.35	0.42
1:F:277:LEU:N	1:F:278:PRO:HD2	2.35	0.42
1:B:169:LYS:H	1:B:170:GLY:HA2	1.83	0.41
1:B:277:LEU:N	1:B:278:PRO:HD2	2.35	0.41
1:A:38:ASN:HA	1:A:39:PRO:HD3	1.89	0.41
1:E:303:THR:HG23	1:E:328:ILE:HD11	2.02	0.41
1:B:125:ALA:HA	1:B:179:THR:HG21	2.02	0.41
1:F:166:LEU:HA	1:F:169:LYS:HD2	2.02	0.41
1:B:303:THR:HG23	1:B:328:ILE:HD11	2.03	0.41
2:C:44:ILE:HD11	2:D:72:LEU:HD11	2.02	0.41
2:D:48:GLU:OE2	2:D:90:HIS:ND1	2.43	0.41
1:A:343:LYS:HB3	1:A:344:ARG:H	1.68	0.41
1:F:177:LEU:HD12	1:F:206:LEU:HD23	2.02	0.41
1:F:333:PRO:O	1:F:337:GLU:HG2	2.20	0.41
2:G:78:LYS:HE3	2:H:37:LYS:HA	2.03	0.41
1:E:92:MET:HG2	1:E:246:ALA:HB1	2.04	0.40
1:F:163:ILE:HG23	1:F:173:VAL:HG21	2.03	0.40
1:F:259:SER:OG	1:F:262:PHE:O	2.37	0.40
1:A:303:THR:HG23	1:A:328:ILE:HD11	2.02	0.40
1:A:332:THR:OG1	1:A:366:THR:O	2.38	0.40
1:E:20:LEU:HD21	1:F:290:VAL:HG11	2.04	0.40
1:B:332:THR:OG1	1:B:366:THR:O	2.39	0.40
1:F:317:GLN:NE2	1:F:321:LYS:HG2	2.36	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	357/416 (86%)	333 (93%)	23 (6%)	1 (0%)	36	66
1	B	359/416 (86%)	342 (95%)	15 (4%)	2 (1%)	21	51
1	E	360/416 (86%)	335 (93%)	22 (6%)	3 (1%)	16	44
1	F	358/416 (86%)	340 (95%)	16 (4%)	2 (1%)	21	51
2	C	66/216 (31%)	65 (98%)	1 (2%)	0	100	100
2	D	66/216 (31%)	65 (98%)	1 (2%)	0	100	100
2	G	67/216 (31%)	66 (98%)	1 (2%)	0	100	100
2	H	66/216 (31%)	65 (98%)	1 (2%)	0	100	100
All	All	1699/2528 (67%)	1611 (95%)	80 (5%)	8 (0%)	24	55

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	148	GLU
1	E	115	GLU
1	E	148	GLU
1	F	314	GLY
1	A	148	GLU
1	B	148	GLU
1	E	27	THR
1	B	314	GLY

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	309/358 (86%)	306 (99%)	3 (1%)	68	88
1	B	311/358 (87%)	307 (99%)	4 (1%)	61	86
1	E	312/358 (87%)	308 (99%)	4 (1%)	61	86
1	F	310/358 (87%)	305 (98%)	5 (2%)	55	83
2	C	60/195 (31%)	60 (100%)	0	100	100
2	D	60/195 (31%)	59 (98%)	1 (2%)	53	83

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	G	61/195 (31%)	61 (100%)	0	100	100
2	H	60/195 (31%)	60 (100%)	0	100	100
All	All	1483/2212 (67%)	1466 (99%)	17 (1%)	65	87

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	VAL
1	A	211	TYR
1	A	344	ARG
1	B	58	TRP
1	B	120	VAL
1	B	211	TYR
1	B	368	GLU
2	D	56	LEU
1	E	120	VAL
1	E	121	LEU
1	E	211	TYR
1	E	344	ARG
1	F	58	TRP
1	F	96	ARG
1	F	120	VAL
1	F	211	TYR
1	F	275	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	208	ASN
1	A	288	HIS
1	B	126	HIS
2	D	97	ASN
1	E	172	ASN
1	E	208	ASN
1	F	288	HIS
2	G	88	GLN
2	H	97	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	B	234	1	23,24,25	1.71	2 (8%)	25,32,34	1.77	4 (16%)
1	LLP	F	234	1	23,24,25	1.74	3 (13%)	25,32,34	1.68	4 (16%)
1	LLP	A	234	1	23,24,25	1.71	2 (8%)	25,32,34	1.69	4 (16%)
1	LLP	E	234	1	23,24,25	1.72	2 (8%)	25,32,34	1.81	3 (12%)

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	B	234	1	-	7/16/17/19	0/1/1/1
1	LLP	F	234	1	-	7/16/17/19	0/1/1/1
1	LLP	A	234	1	-	7/16/17/19	0/1/1/1
1	LLP	E	234	1	-	7/16/17/19	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	234	LLP	O3-C3	-5.89	1.23	1.36
1	F	234	LLP	O3-C3	-5.88	1.23	1.36
1	B	234	LLP	O3-C3	-5.86	1.23	1.36
1	A	234	LLP	O3-C3	-5.83	1.23	1.36
1	F	234	LLP	C2-N1	2.67	1.38	1.33
1	A	234	LLP	C2-N1	2.58	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	234	LLP	C2-N1	2.55	1.38	1.33
1	B	234	LLP	C2-N1	2.49	1.38	1.33
1	F	234	LLP	C6-N1	2.02	1.38	1.34

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	234	LLP	OP4-C5'-C5	6.71	121.94	109.36
1	B	234	LLP	OP4-C5'-C5	6.41	121.37	109.36
1	A	234	LLP	OP4-C5'-C5	5.83	120.29	109.36
1	F	234	LLP	OP4-C5'-C5	5.63	119.92	109.36
1	E	234	LLP	C4-C4'-NZ	-3.16	109.45	124.04
1	B	234	LLP	C4-C4'-NZ	-3.12	109.67	124.04
1	F	234	LLP	C4-C4'-NZ	-3.09	109.78	124.04
1	A	234	LLP	C4-C4'-NZ	-2.93	110.51	124.04
1	A	234	LLP	CE-NZ-C4'	-2.64	110.28	118.72
1	F	234	LLP	CE-NZ-C4'	-2.61	110.35	118.72
1	E	234	LLP	CE-NZ-C4'	-2.58	110.45	118.72
1	B	234	LLP	CE-NZ-C4'	-2.56	110.51	118.72
1	A	234	LLP	C5-C6-N1	-2.22	120.22	123.83
1	B	234	LLP	C5-C6-N1	-2.14	120.35	123.83
1	F	234	LLP	C5-C6-N1	-2.14	120.36	123.83

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	234	LLP	C4-C4'-NZ-CE
1	A	234	LLP	C4-C5-C5'-OP4
1	A	234	LLP	C6-C5-C5'-OP4
1	A	234	LLP	C5'-OP4-P-OP1
1	A	234	LLP	C5'-OP4-P-OP2
1	A	234	LLP	C5'-OP4-P-OP3
1	B	234	LLP	C4-C4'-NZ-CE
1	B	234	LLP	C4-C5-C5'-OP4
1	B	234	LLP	C6-C5-C5'-OP4
1	B	234	LLP	C5'-OP4-P-OP1
1	B	234	LLP	C5'-OP4-P-OP2
1	B	234	LLP	C5'-OP4-P-OP3
1	B	234	LLP	O-C-CA-CB
1	E	234	LLP	C4-C5-C5'-OP4
1	E	234	LLP	C6-C5-C5'-OP4

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Mol	Chain	Res	Type	Atoms
1	E	234	LLP	C5'-OP4-P-OP1
1	E	234	LLP	C5'-OP4-P-OP2
1	E	234	LLP	C5'-OP4-P-OP3
1	E	234	LLP	C-CA-CB-CG
1	F	234	LLP	C4-C4'-NZ-CE
1	F	234	LLP	C4-C5-C5'-OP4
1	F	234	LLP	C5'-OP4-P-OP1
1	F	234	LLP	C5'-OP4-P-OP2
1	F	234	LLP	C5'-OP4-P-OP3
1	F	234	LLP	O-C-CA-CB
1	E	234	LLP	C4-C4'-NZ-CE
1	F	234	LLP	C6-C5-C5'-OP4
1	A	234	LLP	CD-CE-NZ-C4'

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	234	LLP	4	0
1	F	234	LLP	3	0
1	A	234	LLP	2	0
1	E	234	LLP	1	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 [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	361/416 (86%)	-1.23	0 100 100	36, 102, 154, 181	0
1	B	363/416 (87%)	-1.32	0 100 100	39, 74, 168, 203	0
1	E	364/416 (87%)	-1.19	0 100 100	42, 102, 163, 201	0
1	F	362/416 (87%)	-1.11	0 100 100	34, 114, 182, 213	0
2	C	68/216 (31%)	-1.33	0 100 100	43, 79, 146, 173	0
2	D	68/216 (31%)	-1.34	0 100 100	39, 71, 169, 182	0
2	G	69/216 (31%)	-1.30	0 100 100	45, 94, 154, 196	0
2	H	68/216 (31%)	-1.08	0 100 100	35, 116, 178, 193	0
All	All	1723/2528 (68%)	-1.22	0 100 100	34, 97, 171, 213	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	LLP	E	234	24/25	0.99	0.05	79,128,147,154	0
1	LLP	B	234	24/25	1.00	0.04	59,80,91,95	0
1	LLP	A	234	24/25	1.00	0.04	29,65,85,92	0
1	LLP	F	234	24/25	1.00	0.03	63,92,100,105	0

6.3 Carbohydrates [i](#)

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.