



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 12:26 AM UTC

PDB ID : 5LOT / pdb_00005lot
Title : Crystal structure of SCP2 thiolase from Leishmania mexicana. Complex of the C123A mutant with acetyl-CoA.
Authors : Harijan, R.K.; Kiema, T.-R.; Wierenga, R.K.
Deposited on : 2016-08-09
Resolution : 2.25 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

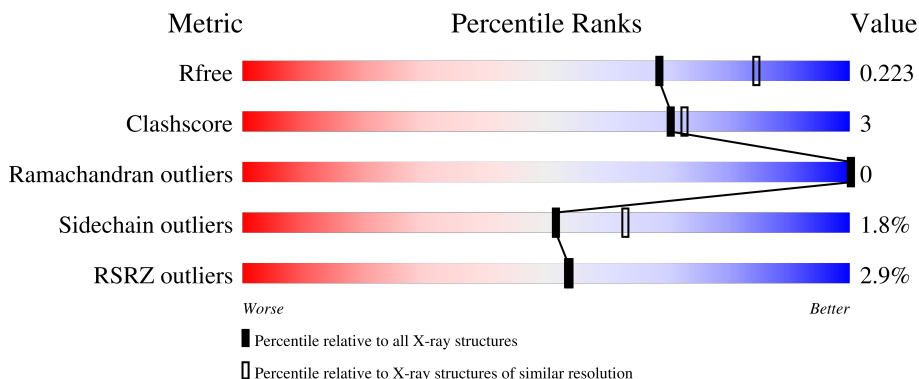
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

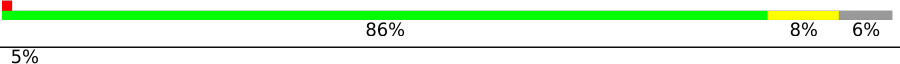
The reported resolution of this entry is 2.25 Å.

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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	457	
1	B	457	
1	C	457	
1	D	457	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13679 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 3-ketoacyl-CoA thiolase-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	0	0
			3223	2010	566	621	26			
1	B	430	Total	C	N	O	S	0	0	0
			3223	2010	566	621	26			
1	C	430	Total	C	N	O	S	0	0	0
			3223	2010	566	621	26			
1	D	430	Total	C	N	O	S	0	0	0
			3223	2010	566	621	26			

There are 180 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	HIS	-	expression tag	UNP E9AW84
A	-14	HIS	-	expression tag	UNP E9AW84
A	-13	HIS	-	expression tag	UNP E9AW84
A	-12	HIS	-	expression tag	UNP E9AW84
A	-11	HIS	-	expression tag	UNP E9AW84
A	-10	HIS	-	expression tag	UNP E9AW84
A	-9	SER	-	expression tag	UNP E9AW84
A	-8	SER	-	expression tag	UNP E9AW84
A	-7	GLY	-	expression tag	UNP E9AW84
A	-6	LEU	-	expression tag	UNP E9AW84
A	-5	VAL	-	expression tag	UNP E9AW84
A	-4	PRO	-	expression tag	UNP E9AW84
A	-3	ARG	-	expression tag	UNP E9AW84
A	-2	GLY	-	expression tag	UNP E9AW84
A	-1	SER	-	expression tag	UNP E9AW84
A	0	HIS	-	expression tag	UNP E9AW84
A	6	LEU	MET	conflict	UNP E9AW84
A	12	ALA	THR	conflict	UNP E9AW84
A	31	PRO	LYS	conflict	UNP E9AW84
A	34	ILE	VAL	conflict	UNP E9AW84
A	45	LYS	GLN	conflict	UNP E9AW84

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Chain	Residue	Modelled	Actual	Comment	Reference
A	59	GLY	HIS	conflict	UNP E9AW84
A	61	MET	ILE	conflict	UNP E9AW84
A	63	HIS	ASN	conflict	UNP E9AW84
A	64	SER	VAL	conflict	UNP E9AW84
A	65	GLY	SER	conflict	UNP E9AW84
A	67	SER	LYS	conflict	UNP E9AW84
A	69	ARG	LYS	conflict	UNP E9AW84
A	71	GLY	ASP	conflict	UNP E9AW84
A	107	SER	ALA	conflict	UNP E9AW84
A	112	MET	ILE	conflict	UNP E9AW84
A	123	ALA	CYS	conflict	UNP E9AW84
A	134	MET	THR	conflict	UNP E9AW84
A	139	SER	ALA	conflict	UNP E9AW84
A	145	THR	VAL	conflict	UNP E9AW84
A	157	SER	THR	conflict	UNP E9AW84
A	176	GLN	LYS	conflict	UNP E9AW84
A	229	THR	SER	conflict	UNP E9AW84
A	268	VAL	ALA	conflict	UNP E9AW84
A	277	ARG	LYS	conflict	UNP E9AW84
A	327	ILE	VAL	conflict	UNP E9AW84
A	356	GLU	ASP	conflict	UNP E9AW84
A	368	ASP	GLU	conflict	UNP E9AW84
A	398	ILE	VAL	conflict	UNP E9AW84
A	410	GLU	GLY	conflict	UNP E9AW84
B	-15	HIS	-	expression tag	UNP E9AW84
B	-14	HIS	-	expression tag	UNP E9AW84
B	-13	HIS	-	expression tag	UNP E9AW84
B	-12	HIS	-	expression tag	UNP E9AW84
B	-11	HIS	-	expression tag	UNP E9AW84
B	-10	HIS	-	expression tag	UNP E9AW84
B	-9	SER	-	expression tag	UNP E9AW84
B	-8	SER	-	expression tag	UNP E9AW84
B	-7	GLY	-	expression tag	UNP E9AW84
B	-6	LEU	-	expression tag	UNP E9AW84
B	-5	VAL	-	expression tag	UNP E9AW84
B	-4	PRO	-	expression tag	UNP E9AW84
B	-3	ARG	-	expression tag	UNP E9AW84
B	-2	GLY	-	expression tag	UNP E9AW84
B	-1	SER	-	expression tag	UNP E9AW84
B	0	HIS	-	expression tag	UNP E9AW84
B	6	LEU	MET	conflict	UNP E9AW84
B	12	ALA	THR	conflict	UNP E9AW84

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Chain	Residue	Modelled	Actual	Comment	Reference
B	31	PRO	LYS	conflict	UNP E9AW84
B	34	ILE	VAL	conflict	UNP E9AW84
B	45	LYS	GLN	conflict	UNP E9AW84
B	59	GLY	HIS	conflict	UNP E9AW84
B	61	MET	ILE	conflict	UNP E9AW84
B	63	HIS	ASN	conflict	UNP E9AW84
B	64	SER	VAL	conflict	UNP E9AW84
B	65	GLY	SER	conflict	UNP E9AW84
B	67	SER	LYS	conflict	UNP E9AW84
B	69	ARG	LYS	conflict	UNP E9AW84
B	71	GLY	ASP	conflict	UNP E9AW84
B	107	SER	ALA	conflict	UNP E9AW84
B	112	MET	ILE	conflict	UNP E9AW84
B	123	ALA	CYS	conflict	UNP E9AW84
B	134	MET	THR	conflict	UNP E9AW84
B	139	SER	ALA	conflict	UNP E9AW84
B	145	THR	VAL	conflict	UNP E9AW84
B	157	SER	THR	conflict	UNP E9AW84
B	176	GLN	LYS	conflict	UNP E9AW84
B	229	THR	SER	conflict	UNP E9AW84
B	268	VAL	ALA	conflict	UNP E9AW84
B	277	ARG	LYS	conflict	UNP E9AW84
B	327	ILE	VAL	conflict	UNP E9AW84
B	356	GLU	ASP	conflict	UNP E9AW84
B	368	ASP	GLU	conflict	UNP E9AW84
B	398	ILE	VAL	conflict	UNP E9AW84
B	410	GLU	GLY	conflict	UNP E9AW84
C	-15	HIS	-	expression tag	UNP E9AW84
C	-14	HIS	-	expression tag	UNP E9AW84
C	-13	HIS	-	expression tag	UNP E9AW84
C	-12	HIS	-	expression tag	UNP E9AW84
C	-11	HIS	-	expression tag	UNP E9AW84
C	-10	HIS	-	expression tag	UNP E9AW84
C	-9	SER	-	expression tag	UNP E9AW84
C	-8	SER	-	expression tag	UNP E9AW84
C	-7	GLY	-	expression tag	UNP E9AW84
C	-6	LEU	-	expression tag	UNP E9AW84
C	-5	VAL	-	expression tag	UNP E9AW84
C	-4	PRO	-	expression tag	UNP E9AW84
C	-3	ARG	-	expression tag	UNP E9AW84
C	-2	GLY	-	expression tag	UNP E9AW84
C	-1	SER	-	expression tag	UNP E9AW84

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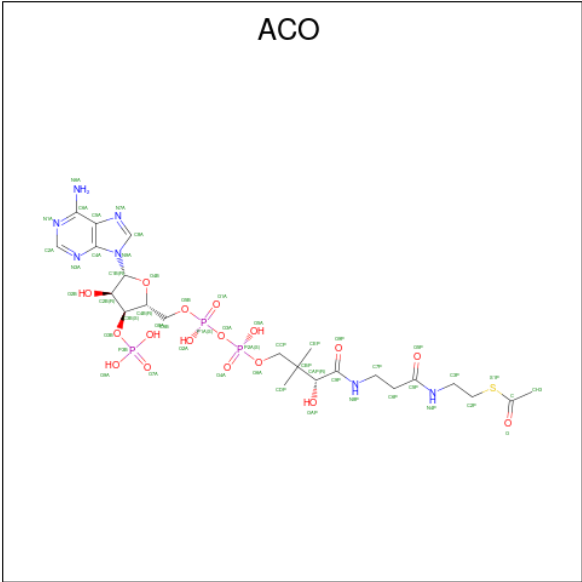
Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP E9AW84
C	6	LEU	MET	conflict	UNP E9AW84
C	12	ALA	THR	conflict	UNP E9AW84
C	31	PRO	LYS	conflict	UNP E9AW84
C	34	ILE	VAL	conflict	UNP E9AW84
C	45	LYS	GLN	conflict	UNP E9AW84
C	59	GLY	HIS	conflict	UNP E9AW84
C	61	MET	ILE	conflict	UNP E9AW84
C	63	HIS	ASN	conflict	UNP E9AW84
C	64	SER	VAL	conflict	UNP E9AW84
C	65	GLY	SER	conflict	UNP E9AW84
C	67	SER	LYS	conflict	UNP E9AW84
C	69	ARG	LYS	conflict	UNP E9AW84
C	71	GLY	ASP	conflict	UNP E9AW84
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C	112	MET	ILE	conflict	UNP E9AW84
C	123	ALA	CYS	conflict	UNP E9AW84
C	134	MET	THR	conflict	UNP E9AW84
C	139	SER	ALA	conflict	UNP E9AW84
C	145	THR	VAL	conflict	UNP E9AW84
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C	327	ILE	VAL	conflict	UNP E9AW84
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D	-15	HIS	-	expression tag	UNP E9AW84
D	-14	HIS	-	expression tag	UNP E9AW84
D	-13	HIS	-	expression tag	UNP E9AW84
D	-12	HIS	-	expression tag	UNP E9AW84
D	-11	HIS	-	expression tag	UNP E9AW84
D	-10	HIS	-	expression tag	UNP E9AW84
D	-9	SER	-	expression tag	UNP E9AW84
D	-8	SER	-	expression tag	UNP E9AW84
D	-7	GLY	-	expression tag	UNP E9AW84
D	-6	LEU	-	expression tag	UNP E9AW84
D	-5	VAL	-	expression tag	UNP E9AW84
D	-4	PRO	-	expression tag	UNP E9AW84

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-3	ARG	-	expression tag	UNP E9AW84
D	-2	GLY	-	expression tag	UNP E9AW84
D	-1	SER	-	expression tag	UNP E9AW84
D	0	HIS	-	expression tag	UNP E9AW84
D	6	LEU	MET	conflict	UNP E9AW84
D	12	ALA	THR	conflict	UNP E9AW84
D	31	PRO	LYS	conflict	UNP E9AW84
D	34	ILE	VAL	conflict	UNP E9AW84
D	45	LYS	GLN	conflict	UNP E9AW84
D	59	GLY	HIS	conflict	UNP E9AW84
D	61	MET	ILE	conflict	UNP E9AW84
D	63	HIS	ASN	conflict	UNP E9AW84
D	64	SER	VAL	conflict	UNP E9AW84
D	65	GLY	SER	conflict	UNP E9AW84
D	67	SER	LYS	conflict	UNP E9AW84
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D	107	SER	ALA	conflict	UNP E9AW84
D	112	MET	ILE	conflict	UNP E9AW84
D	123	ALA	CYS	conflict	UNP E9AW84
D	134	MET	THR	conflict	UNP E9AW84
D	139	SER	ALA	conflict	UNP E9AW84
D	145	THR	VAL	conflict	UNP E9AW84
D	157	SER	THR	conflict	UNP E9AW84
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D	327	ILE	VAL	conflict	UNP E9AW84
D	356	GLU	ASP	conflict	UNP E9AW84
D	368	ASP	GLU	conflict	UNP E9AW84
D	398	ILE	VAL	conflict	UNP E9AW84
D	410	GLU	GLY	conflict	UNP E9AW84

- Molecule 2 is ACETYL COENZYME *A (CCD ID: ACO) (formula: C₂₃H₃₈N₇O₁₇P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		
2	D	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		

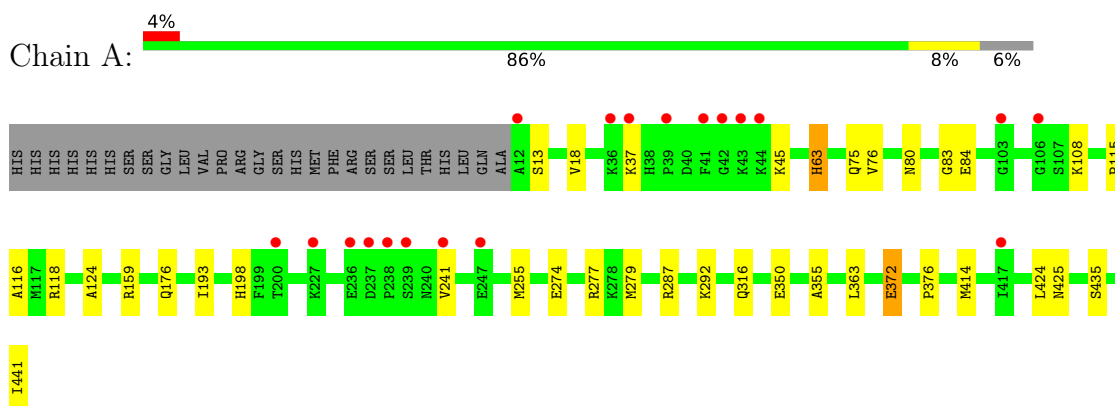
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	122	Total	O	0	0
			122	122		
3	B	197	Total	O	0	0
			197	197		
3	C	151	Total	O	0	0
			151	151		
3	D	164	Total	O	0	0
			164	164		

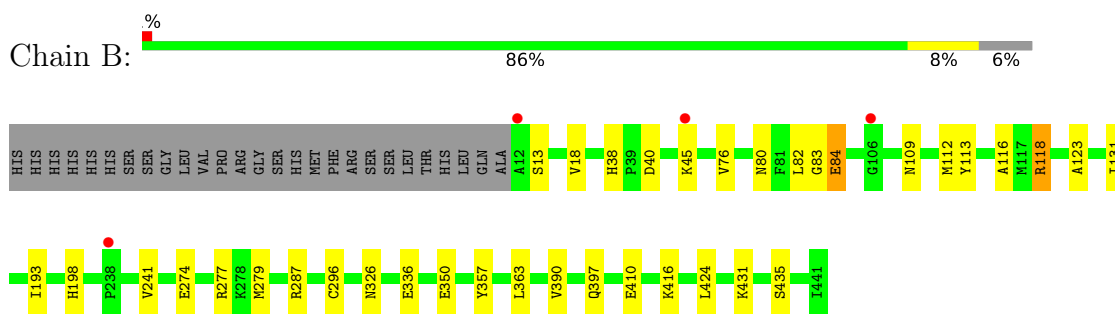
3 Residue-property plots [i](#)

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.

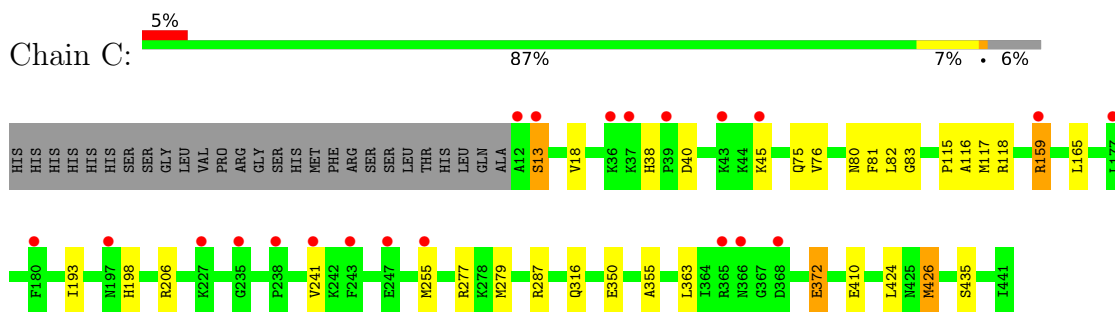
- Molecule 1: 3-ketoacyl-CoA thiolase-like protein



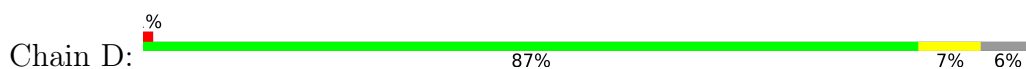
- Molecule 1: 3-ketoacyl-CoA thiolase-like protein

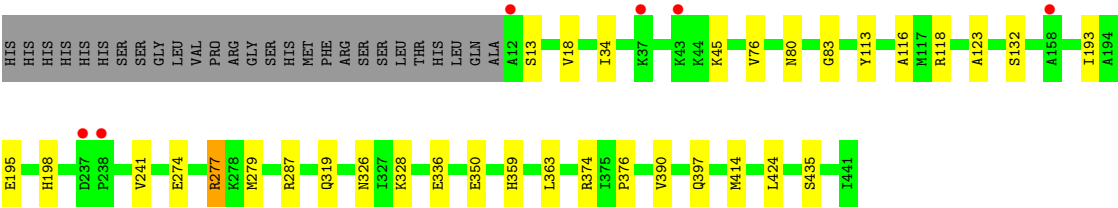


- Molecule 1: 3-ketoacyl-CoA thiolase-like protein



- Molecule 1: 3-ketoacyl-CoA thiolase-like protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.70Å 89.89Å 126.33Å 90.00° 105.78° 90.00°	Depositor
Resolution (Å)	47.30 – 2.25 47.30 – 2.25	Depositor EDS
% Data completeness (in resolution range)	98.2 (47.30-2.25) 98.2 (47.30-2.25)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.213 , 0.255 0.221 , 0.223	Depositor DCC
R_{free} test set	3952 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.012 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13679	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.28% 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: ACO

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	0.82	0/3276	0.88	1/4417 (0.0%)
1	B	0.91	3/3276 (0.1%)	0.93	1/4417 (0.0%)
1	C	0.87	0/3276	0.92	2/4417 (0.0%)
1	D	0.86	1/3276 (0.0%)	0.89	1/4417 (0.0%)
All	All	0.87	4/13104 (0.0%)	0.91	5/17668 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	109	ASN	C-O	-6.00	1.20	1.25
1	B	241	VAL	CA-CB	5.72	1.60	1.54
1	B	84	GLU	CD-OE1	5.62	1.36	1.25
1	D	241	VAL	CA-CB	5.57	1.60	1.54

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	426	MET	CB-CG-SD	-5.90	95.01	112.70
1	C	159	ARG	N-CA-CB	-5.54	102.03	110.07
1	B	287	ARG	CG-CD-NE	-5.36	100.21	112.00
1	A	63	HIS	CB-CA-C	5.17	121.13	110.31
1	D	287	ARG	CG-CD-NE	-5.17	100.62	112.00

There are no chirality outliers.

There are no planarity outliers.

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	3223	0	3202	27	0
1	B	3223	0	3202	21	0
1	C	3223	0	3202	23	0
1	D	3223	0	3202	18	0
2	A	51	0	34	3	0
2	B	51	0	34	1	0
2	D	51	0	34	3	0
3	A	122	0	0	5	0
3	B	197	0	0	6	0
3	C	151	0	0	6	0
3	D	164	0	0	6	0
All	All	13679	0	12910	88	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 3.

All (88) 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:441:ILE:C	3:A:632:HOH:O	1.83	1.22
1:C:410:GLU:HG3	3:C:614:HOH:O	1.80	0.82
1:D:18:VAL:HG21	1:D:279:MET:HE1	1.62	0.81
2:A:501:ACO:H8A	2:A:501:ACO:O5B	1.82	0.79
1:A:18:VAL:O	1:A:287:ARG:NH1	2.17	0.78
1:D:374:ARG:HD2	3:D:722:HOH:O	1.85	0.77
1:D:336:GLU:OE2	3:D:601:HOH:O	2.13	0.67
1:C:75:GLN:HG2	1:C:115:PRO:HB2	1.78	0.66
1:A:75:GLN:HG2	1:A:115:PRO:HB2	1.81	0.63
1:D:80:ASN:HD21	1:D:83:GLY:H	1.49	0.61
1:A:80:ASN:HD21	1:A:83:GLY:H	1.48	0.61
1:B:123:ALA:HA	1:B:390:VAL:HG13	1.83	0.61
1:D:123:ALA:HA	1:D:390:VAL:HG13	1.82	0.60
1:B:84:GLU:CD	1:B:118:ARG:HH22	2.09	0.60
1:B:18:VAL:HG11	1:B:279:MET:HE1	1.84	0.59
1:B:80:ASN:HD21	1:B:83:GLY:H	1.50	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:ASN:HD21	1:C:83:GLY:H	1.50	0.58
1:D:195:GLU:HG3	3:D:677:HOH:O	2.04	0.58
1:A:18:VAL:HG11	1:A:279:MET:HE1	1.84	0.57
1:A:274:GLU:OE2	1:A:277:ARG:NH2	2.37	0.57
1:B:410:GLU:HG3	3:B:742:HOH:O	2.04	0.57
1:C:18:VAL:HG11	1:C:279:MET:HE1	1.86	0.57
2:D:501:ACO:HH32	3:D:713:HOH:O	2.05	0.57
2:D:501:ACO:H131	2:D:501:ACO:O9P	2.04	0.57
1:A:37:LYS:HE3	3:A:702:HOH:O	2.05	0.56
1:A:37:LYS:HG3	3:A:702:HOH:O	2.08	0.54
1:B:193:ILE:HG23	1:B:198:HIS:HB3	1.89	0.54
1:B:431:LYS:HE3	3:B:747:HOH:O	2.07	0.54
1:A:193:ILE:HG23	1:A:198:HIS:HB3	1.90	0.54
2:A:501:ACO:H141	2:A:501:ACO:O9P	2.08	0.53
1:A:176:GLN:HE21	1:A:176:GLN:HA	1.74	0.53
1:A:84:GLU:HG3	3:B:602:HOH:O	2.09	0.53
1:C:38:HIS:CE1	3:C:529:HOH:O	2.61	0.52
1:C:193:ILE:HG23	1:C:198:HIS:HB3	1.91	0.52
1:D:193:ILE:HG23	1:D:198:HIS:HB3	1.91	0.52
1:A:424:LEU:HD13	1:A:435:SER:HB2	1.92	0.51
1:B:274:GLU:OE2	1:B:277:ARG:NH2	2.44	0.51
1:C:372:GLU:OE2	1:C:372:GLU:HA	2.11	0.51
1:B:198:HIS:HD2	3:B:757:HOH:O	1.94	0.51
1:B:38:HIS:HD2	1:B:40:ASP:H	1.58	0.51
1:C:316:GLN:HG3	1:D:113:TYR:CG	2.46	0.50
1:A:18:VAL:HG12	1:A:287:ARG:NH1	2.27	0.49
2:B:501:ACO:HH32	3:B:721:HOH:O	2.10	0.49
1:D:274:GLU:OE2	1:D:277:ARG:NH2	2.45	0.49
1:A:159:ARG:HB2	3:A:686:HOH:O	2.10	0.49
2:D:501:ACO:O9P	2:D:501:ACO:CDP	2.61	0.49
1:C:165:LEU:HB3	3:C:575:HOH:O	2.13	0.48
1:D:18:VAL:CG2	1:D:279:MET:HE1	2.41	0.48
1:A:372:GLU:HA	1:A:372:GLU:OE2	2.14	0.48
1:C:241:VAL:HG12	1:C:255:MET:SD	2.54	0.48
1:D:34:ILE:HD11	3:D:740:HOH:O	2.14	0.48
1:A:108:LYS:HE3	1:C:287:ARG:HB2	1.95	0.47
1:C:424:LEU:HD13	1:C:435:SER:HB2	1.97	0.46
1:C:117:MET:HE1	1:D:132:SER:OG	2.16	0.46
1:A:76:VAL:O	1:A:116:ALA:HA	2.16	0.46
1:D:76:VAL:O	1:D:116:ALA:HA	2.16	0.46
1:B:336:GLU:HG2	1:B:397:GLN:HA	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:424:LEU:HD13	1:D:435:SER:HB2	1.97	0.45
1:C:76:VAL:O	1:C:116:ALA:HA	2.16	0.45
1:B:416:LYS:HE2	3:B:786:HOH:O	2.17	0.45
1:B:424:LEU:HD13	1:B:435:SER:HB2	1.99	0.45
1:A:63:HIS:O	1:A:287:ARG:NH2	2.44	0.45
1:B:76:VAL:O	1:B:116:ALA:HA	2.17	0.45
1:A:84:GLU:OE1	1:B:82:LEU:HD23	2.17	0.45
1:A:316:GLN:HG3	1:B:113:TYR:CG	2.51	0.45
1:A:241:VAL:HG12	1:A:255:MET:SD	2.57	0.44
1:C:13:SER:CB	3:C:601:HOH:O	2.65	0.44
1:C:206:ARG:HG2	3:C:636:HOH:O	2.18	0.43
1:C:13:SER:HB3	3:C:601:HOH:O	2.18	0.43
1:A:108:LYS:HE3	1:C:287:ARG:CB	2.48	0.43
1:A:355:ALA:HB2	1:A:363:LEU:HD13	1.99	0.43
1:C:38:HIS:HD2	1:C:40:ASP:H	1.66	0.43
1:C:355:ALA:HB2	1:C:363:LEU:HD13	2.00	0.43
3:A:671:HOH:O	1:B:112:MET:HE1	2.19	0.43
1:B:38:HIS:CD2	1:B:40:ASP:H	2.35	0.43
1:D:198:HIS:HE1	1:D:350:GLU:OE1	2.02	0.43
1:D:336:GLU:HG2	1:D:397:GLN:HA	2.00	0.42
1:D:376:PRO:HG2	1:D:414:MET:HE3	2.02	0.42
1:A:241:VAL:HG21	2:A:501:ACO:C2A	2.50	0.42
1:C:81:PHE:CG	1:C:82:LEU:HG	2.55	0.42
1:D:359:HIS:CE1	3:D:736:HOH:O	2.73	0.41
1:B:357:TYR:CD1	1:C:277:ARG:HD3	2.55	0.41
1:A:124:ALA:HA	1:A:425:ASN:OD1	2.21	0.41
1:B:198:HIS:HE1	1:B:350:GLU:OE1	2.03	0.41
1:C:198:HIS:HE1	1:C:350:GLU:OE1	2.03	0.41
1:A:198:HIS:HE1	1:A:350:GLU:OE1	2.03	0.41
1:A:376:PRO:HG2	1:A:414:MET:HE3	2.02	0.41
1:B:131:ILE:CD1	1:B:296:CYS:HB2	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	428/457 (94%)	420 (98%)	8 (2%)	0	100	100
1	B	428/457 (94%)	418 (98%)	10 (2%)	0	100	100
1	C	428/457 (94%)	420 (98%)	8 (2%)	0	100	100
1	D	428/457 (94%)	419 (98%)	9 (2%)	0	100	100
All	All	1712/1828 (94%)	1677 (98%)	35 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	338/362 (93%)	333 (98%)	5 (2%)	57	68
1	B	338/362 (93%)	333 (98%)	5 (2%)	57	68
1	C	338/362 (93%)	332 (98%)	6 (2%)	51	63
1	D	338/362 (93%)	330 (98%)	8 (2%)	43	54
All	All	1352/1448 (93%)	1328 (98%)	24 (2%)	51	63

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	SER
1	A	45	LYS
1	A	118	ARG
1	A	292	LYS
1	A	372	GLU
1	B	13	SER
1	B	45	LYS
1	B	118	ARG
1	B	326	ASN
1	B	363	LEU

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Mol	Chain	Res	Type
1	C	13	SER
1	C	45	LYS
1	C	118	ARG
1	C	159	ARG
1	C	372	GLU
1	C	426	MET
1	D	13	SER
1	D	45	LYS
1	D	118	ARG
1	D	277	ARG
1	D	319	GLN
1	D	326	ASN
1	D	328	LYS
1	D	363	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
1	A	176	GLN
1	A	198	HIS
1	A	316	GLN
1	A	319	GLN
1	A	425	ASN
1	B	38	HIS
1	B	46	ASN
1	B	58	GLN
1	B	63	HIS
1	B	80	ASN
1	B	153	GLN
1	B	198	HIS
1	B	260	GLN
1	B	316	GLN
1	B	366	ASN
1	B	408	GLN
1	B	425	ASN
1	C	38	HIS
1	C	80	ASN
1	C	198	HIS
1	C	260	GLN
1	C	326	ASN
1	C	408	GLN

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Mol	Chain	Res	Type
1	C	425	ASN
1	D	80	ASN
1	D	153	GLN
1	D	198	HIS
1	D	260	GLN
1	D	359	HIS
1	D	366	ASN
1	D	425	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ACO	D	501	-	51,53,53	1.31	5 (9%)	73,79,79	2.10	24 (32%)
2	ACO	A	501	-	51,53,53	1.30	4 (7%)	73,79,79	1.85	20 (27%)
2	ACO	B	501	-	51,53,53	1.31	5 (9%)	73,79,79	1.69	14 (19%)

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
2	ACO	D	501	-	-	22/51/67/67	0/3/3/3
2	ACO	A	501	-	-	31/51/67/67	0/3/3/3
2	ACO	B	501	-	-	10/51/67/67	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	ACO	C5A-C4A	4.84	1.47	1.39
2	B	501	ACO	C5A-C4A	4.76	1.47	1.39
2	A	501	ACO	C5A-C4A	4.17	1.46	1.39
2	D	501	ACO	P2A-O3A	3.77	1.63	1.59
2	A	501	ACO	P1A-O3A	-3.28	1.56	1.59
2	B	501	ACO	C2B-C3B	-2.92	1.46	1.53
2	A	501	ACO	P2A-O3A	-2.89	1.56	1.59
2	B	501	ACO	P1A-O3A	-2.77	1.56	1.59
2	D	501	ACO	C5A-C6A	2.62	1.48	1.41
2	B	501	ACO	C4A-N9A	-2.61	1.32	1.37
2	A	501	ACO	C5A-C6A	2.60	1.48	1.41
2	D	501	ACO	P1A-O3A	2.55	1.62	1.59
2	D	501	ACO	C8A-N7A	2.49	1.36	1.31
2	B	501	ACO	C5A-C6A	2.37	1.47	1.41

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	ACO	C5A-C4A-N3A	-5.31	119.40	126.72
2	D	501	ACO	N3A-C2A-N1A	-4.70	121.47	128.58
2	B	501	ACO	C7P-C6P-C5P	-4.64	104.66	112.39
2	D	501	ACO	C6P-C5P-N4P	4.57	124.67	116.34
2	D	501	ACO	C7P-N8P-C9P	4.53	130.68	122.55
2	A	501	ACO	C5A-C4A-N3A	-4.52	120.49	126.72
2	D	501	ACO	N3A-C4A-N9A	4.48	134.79	127.17
2	D	501	ACO	C2A-N3A-C4A	4.17	122.02	111.83
2	B	501	ACO	C4A-N9A-C8A	4.16	110.11	105.74
2	B	501	ACO	N3A-C4A-N9A	4.06	134.07	127.17
2	A	501	ACO	C4A-C5A-N7A	-4.03	105.97	110.58
2	D	501	ACO	CAP-C9P-N8P	4.02	124.11	116.48
2	B	501	ACO	C5A-C4A-N3A	-3.89	121.36	126.72
2	D	501	ACO	C3P-N4P-C5P	3.76	129.81	122.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ACO	C2P-S1P-C	3.67	119.08	101.42
2	A	501	ACO	N3A-C2A-N1A	-3.66	123.03	128.58
2	D	501	ACO	O6A-CCP-CBP	3.58	116.30	110.55
2	D	501	ACO	C2P-C3P-N4P	-3.57	104.97	112.41
2	B	501	ACO	N3A-C2A-N1A	-3.52	123.25	128.58
2	A	501	ACO	N3A-C4A-N9A	3.43	133.00	127.17
2	A	501	ACO	C6A-C5A-N7A	3.38	138.60	132.09
2	A	501	ACO	C2A-N3A-C4A	3.35	120.01	111.83
2	A	501	ACO	C4A-N9A-C8A	3.19	109.09	105.74
2	A	501	ACO	C5A-N7A-C8A	2.98	108.13	103.45
2	B	501	ACO	C2A-N3A-C4A	2.97	119.10	111.83
2	D	501	ACO	O5P-C5P-N4P	-2.97	117.19	123.03
2	A	501	ACO	O5P-C5P-C6P	-2.95	116.67	122.02
2	A	501	ACO	C2A-N1A-C6A	2.88	123.46	118.73
2	D	501	ACO	C4A-N9A-C8A	2.87	108.75	105.74
2	D	501	ACO	C4A-C5A-N7A	-2.75	107.44	110.58
2	A	501	ACO	P3B-O3B-C3B	-2.74	116.12	123.43
2	A	501	ACO	C2B-C1B-N9A	-2.72	106.55	113.30
2	B	501	ACO	O5A-P2A-O4A	2.68	124.92	112.44
2	B	501	ACO	C6A-C5A-N7A	2.62	137.15	132.09
2	D	501	ACO	C6A-C5A-N7A	2.62	137.15	132.09
2	B	501	ACO	OAP-CAP-CBP	-2.59	104.18	110.18
2	D	501	ACO	O9P-C9P-N8P	-2.59	117.51	122.98
2	A	501	ACO	N9A-C8A-N7A	-2.57	110.28	113.94
2	B	501	ACO	O9A-P3B-O7A	2.51	120.62	110.83
2	A	501	ACO	C3B-C2B-C1B	2.49	105.38	99.89
2	D	501	ACO	CDP-CBP-CCP	2.47	112.30	108.22
2	D	501	ACO	CDP-CBP-CAP	2.46	112.95	108.77
2	A	501	ACO	C2P-C3P-N4P	2.44	117.50	112.41
2	D	501	ACO	O3A-P1A-O1A	-2.43	103.39	110.70
2	D	501	ACO	O2A-P1A-O1A	2.41	123.63	112.44
2	B	501	ACO	O5P-C5P-C6P	-2.40	117.67	122.02
2	A	501	ACO	C3P-N4P-C5P	2.38	127.25	122.82
2	B	501	ACO	C4A-N9A-C1B	-2.37	121.09	126.63
2	B	501	ACO	C2A-N1A-C6A	2.35	122.59	118.73
2	A	501	ACO	O3A-P1A-O1A	2.34	117.73	110.70
2	D	501	ACO	C2A-N1A-C6A	2.29	122.50	118.73
2	D	501	ACO	C5A-N7A-C8A	2.28	107.04	103.45
2	A	501	ACO	O6A-CCP-CBP	2.20	114.09	110.55
2	D	501	ACO	O5P-C5P-C6P	-2.16	118.11	122.02
2	D	501	ACO	N9A-C8A-N7A	-2.09	110.97	113.94
2	B	501	ACO	C2P-C3P-N4P	2.08	116.76	112.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	ACO	CEP-CBP-CDP	-2.06	105.09	109.20
2	A	501	ACO	C4A-N9A-C1B	-2.01	121.92	126.63

There are no chirality outliers.

All (63) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	ACO	C5B-O5B-P1A-O1A
2	A	501	ACO	C5B-O5B-P1A-O3A
2	A	501	ACO	CCP-O6A-P2A-O3A
2	A	501	ACO	CCP-O6A-P2A-O4A
2	A	501	ACO	CCP-O6A-P2A-O5A
2	A	501	ACO	CAP-CBP-CCP-O6A
2	A	501	ACO	OAP-CAP-CBP-CCP
2	A	501	ACO	C9P-CAP-CBP-CCP
2	A	501	ACO	OAP-CAP-CBP-CDP
2	A	501	ACO	C9P-CAP-CBP-CDP
2	A	501	ACO	OAP-CAP-CBP-CEP
2	A	501	ACO	C9P-CAP-CBP-CEP
2	A	501	ACO	C5P-C6P-C7P-N8P
2	A	501	ACO	C6P-C5P-N4P-C3P
2	A	501	ACO	O5P-C5P-N4P-C3P
2	A	501	ACO	S1P-C2P-C3P-N4P
2	A	501	ACO	C3P-C2P-S1P-C
2	A	501	ACO	O-C-S1P-C2P
2	A	501	ACO	CH3-C-S1P-C2P
2	B	501	ACO	C3P-C2P-S1P-C
2	D	501	ACO	O4B-C4B-C5B-O5B
2	D	501	ACO	C5B-O5B-P1A-O1A
2	D	501	ACO	C5B-O5B-P1A-O2A
2	D	501	ACO	CCP-O6A-P2A-O3A
2	D	501	ACO	CCP-O6A-P2A-O4A
2	D	501	ACO	CCP-O6A-P2A-O5A
2	D	501	ACO	CDP-CBP-CCP-O6A
2	D	501	ACO	CAP-CBP-CCP-O6A
2	D	501	ACO	CAP-C9P-N8P-C7P
2	D	501	ACO	C3P-C2P-S1P-C
2	D	501	ACO	O-C-S1P-C2P
2	D	501	ACO	CH3-C-S1P-C2P
2	D	501	ACO	C6P-C5P-N4P-C3P
2	D	501	ACO	O9P-C9P-N8P-C7P
2	D	501	ACO	C3B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
2	D	501	ACO	O5P-C5P-N4P-C3P
2	A	501	ACO	C2B-C3B-O3B-P3B
2	B	501	ACO	C2B-C3B-O3B-P3B
2	A	501	ACO	C4B-C3B-O3B-P3B
2	B	501	ACO	C4B-C3B-O3B-P3B
2	B	501	ACO	O4B-C4B-C5B-O5B
2	B	501	ACO	C3B-C4B-C5B-O5B
2	A	501	ACO	P2A-O3A-P1A-O1A
2	A	501	ACO	CDP-CBP-CCP-O6A
2	A	501	ACO	CEP-CBP-CCP-O6A
2	D	501	ACO	CEP-CBP-CCP-O6A
2	B	501	ACO	O-C-S1P-C2P
2	A	501	ACO	O9P-C9P-N8P-C7P
2	A	501	ACO	C5B-O5B-P1A-O2A
2	B	501	ACO	C5B-O5B-P1A-O1A
2	D	501	ACO	C5B-O5B-P1A-O3A
2	A	501	ACO	CAP-C9P-N8P-C7P
2	A	501	ACO	O9P-C9P-CAP-CBP
2	D	501	ACO	O9P-C9P-CAP-CBP
2	D	501	ACO	C3B-O3B-P3B-O9A
2	A	501	ACO	N8P-C9P-CAP-CBP
2	D	501	ACO	N8P-C9P-CAP-CBP
2	B	501	ACO	CH3-C-S1P-C2P
2	A	501	ACO	P2A-O3A-P1A-O2A
2	B	501	ACO	C3B-O3B-P3B-O9A
2	D	501	ACO	CBP-CCP-O6A-P2A
2	A	501	ACO	C6P-C7P-N8P-C9P
2	B	501	ACO	P1A-O3A-P2A-O5A

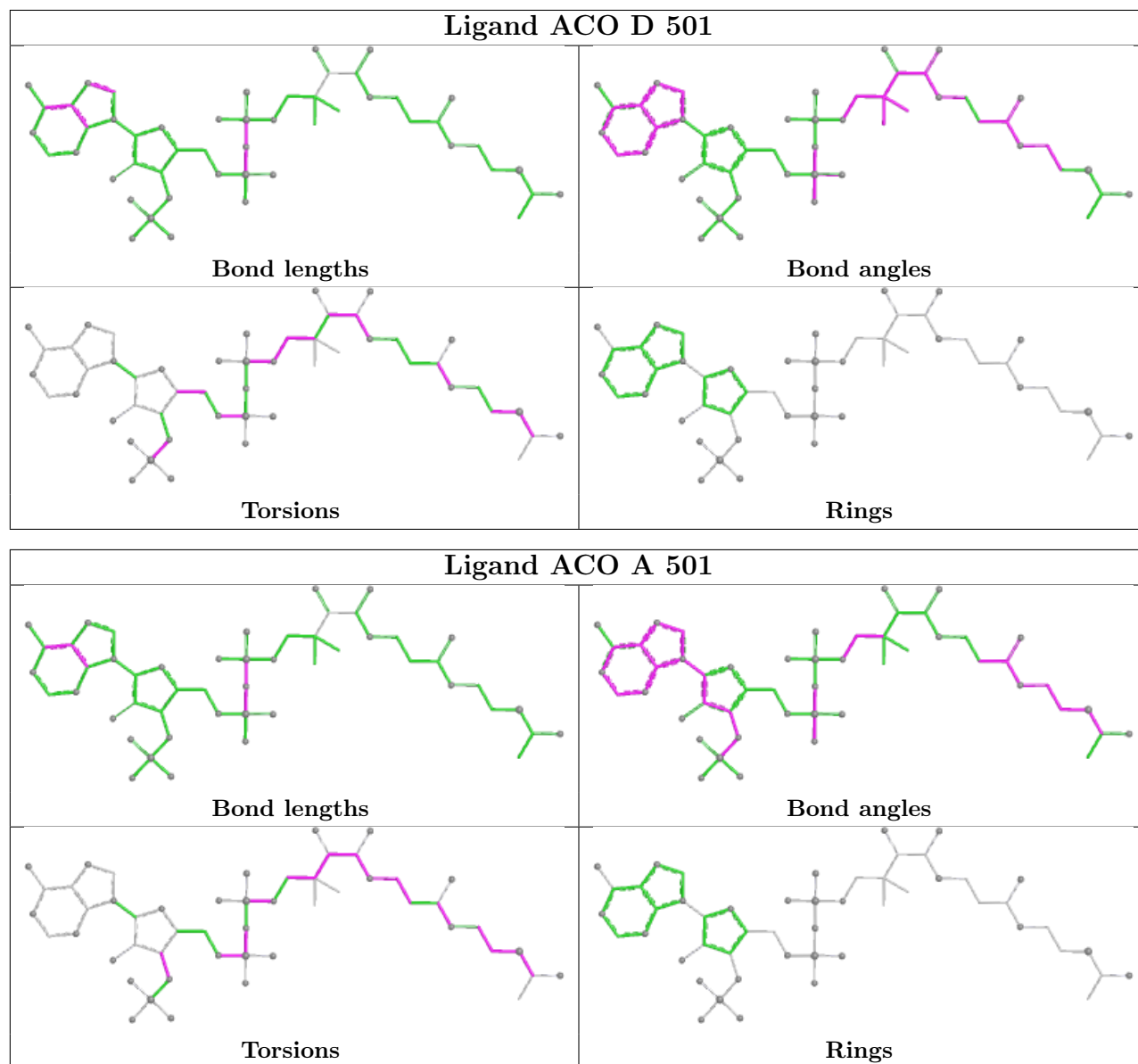
There are no ring outliers.

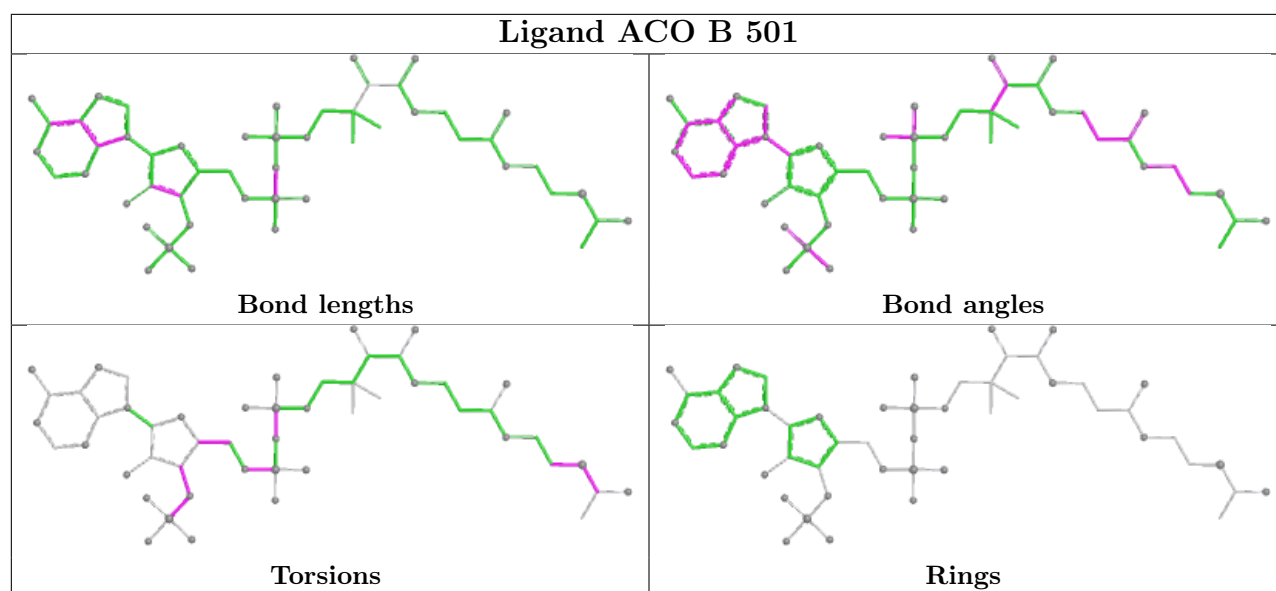
3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	ACO	3	0
2	A	501	ACO	3	0
2	B	501	ACO	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	430/457 (94%)	0.50	19 (4%) 39 38	15, 31, 59, 99	0
1	B	430/457 (94%)	0.00	4 (0%) 81 82	12, 22, 42, 77	0
1	C	430/457 (94%)	0.38	21 (4%) 35 34	13, 28, 59, 90	0
1	D	430/457 (94%)	0.21	6 (1%) 73 74	14, 26, 51, 89	0
All	All	1720/1828 (94%)	0.27	50 (2%) 53 54	12, 26, 54, 99	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	12	ALA	4.0
1	B	12	ALA	3.7
1	D	43	LYS	3.7
1	A	237	ASP	3.7
1	A	12	ALA	3.7
1	C	37	LYS	3.7
1	A	37	LYS	3.5
1	C	12	ALA	3.3
1	A	39	PRO	3.1
1	C	238	PRO	3.1
1	A	238	PRO	3.1
1	D	158	ALA	3.1
1	A	103	GLY	3.0
1	C	159	ARG	3.0
1	A	36	LYS	3.0
1	A	247	GLU	2.9
1	C	36	LYS	2.9
1	A	106	GLY	2.9
1	C	177	LEU	2.8
1	A	43	LYS	2.8
1	B	238	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	243	PHE	2.7
1	A	42	GLY	2.7
1	A	239	SER	2.7
1	C	227	LYS	2.6
1	A	227	LYS	2.6
1	C	241	VAL	2.6
1	C	39	PRO	2.5
1	D	238	PRO	2.5
1	A	44	LYS	2.5
1	C	197	ASN	2.5
1	A	200	THR	2.4
1	C	368	ASP	2.3
1	C	45	LYS	2.3
1	C	235	GLY	2.3
1	A	236	GLU	2.3
1	C	247	GLU	2.3
1	D	237	ASP	2.3
1	C	255	MET	2.3
1	A	241	VAL	2.3
1	C	366	ASN	2.2
1	B	106	GLY	2.2
1	D	37	LYS	2.2
1	A	417	ILE	2.2
1	C	43	LYS	2.2
1	A	41	PHE	2.1
1	C	13	SER	2.1
1	C	180	PHE	2.1
1	B	45	LYS	2.0
1	C	365	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

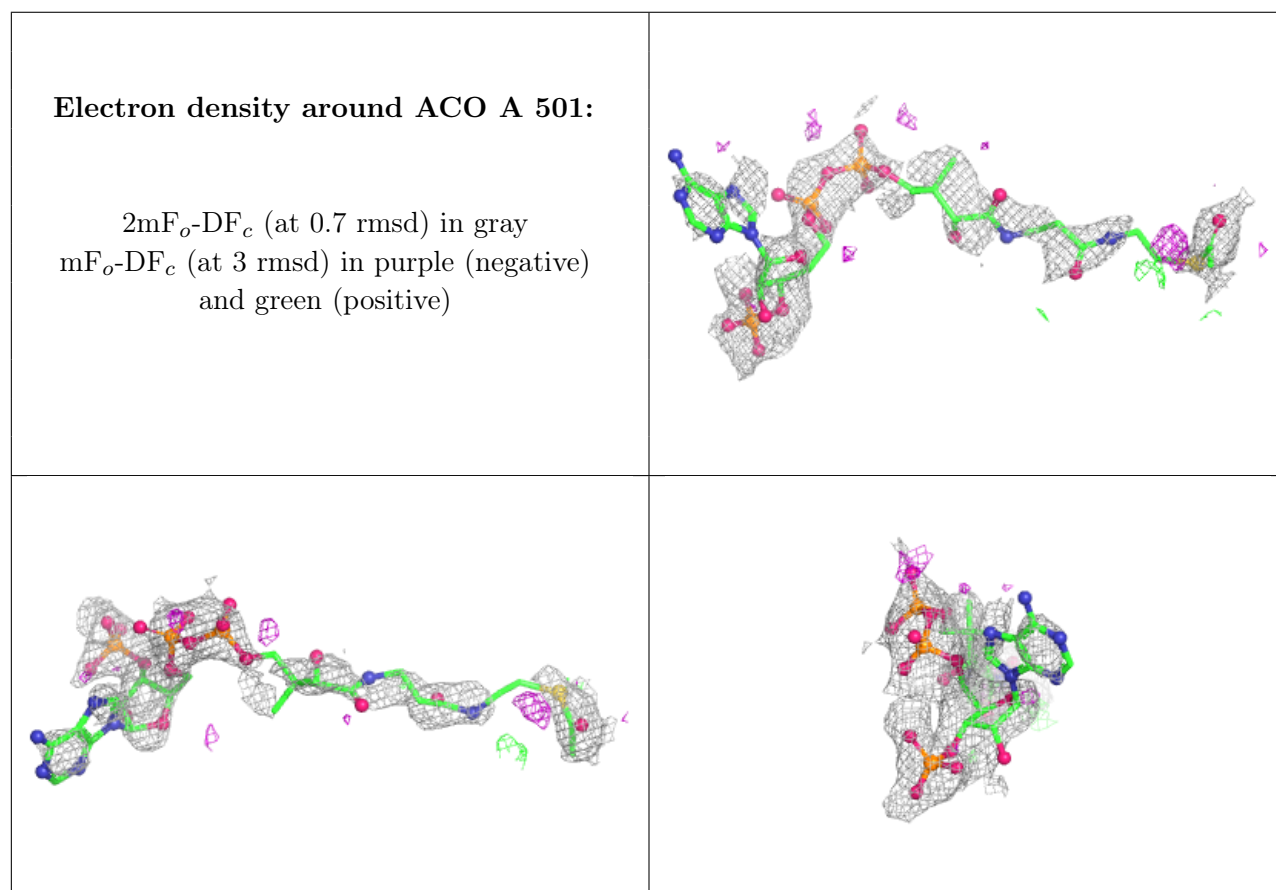
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

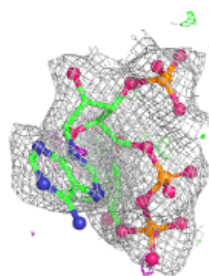
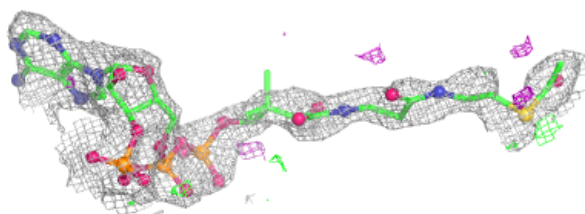
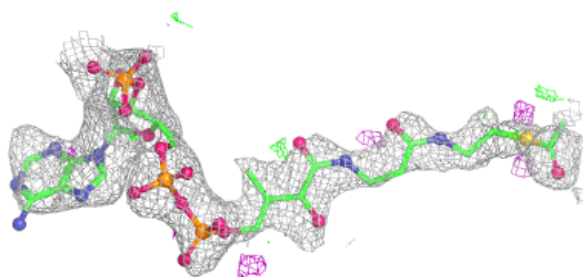
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ACO	A	501	51/51	0.72	0.21	42,97,126,131	0
2	ACO	D	501	51/51	0.79	0.16	45,64,78,83	0
2	ACO	B	501	51/51	0.94	0.07	16,21,28,44	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

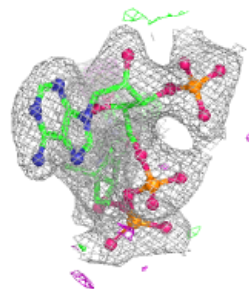
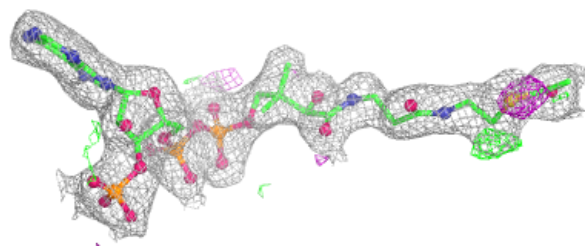
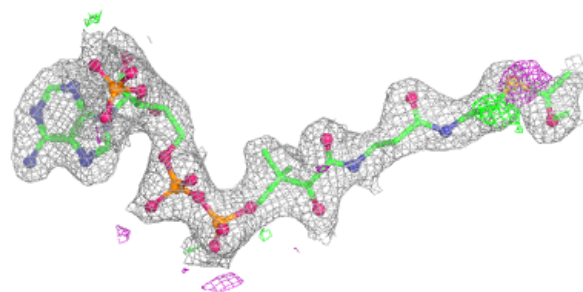


Electron density around ACO D 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around ACO B 501:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.