



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 09:48 PM UTC

PDB ID : 4LCM / pdb_00004lcm
Title : Simvastatin Synthase (LOVD), from *Aspergillus Terreus*, LovD9 mutant (simh9014)
Authors : Gao, X.; Sawaya, M.R.; Yeates, T.O.; Tang, Y.
Deposited on : 2013-06-21
Resolution : 3.19 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

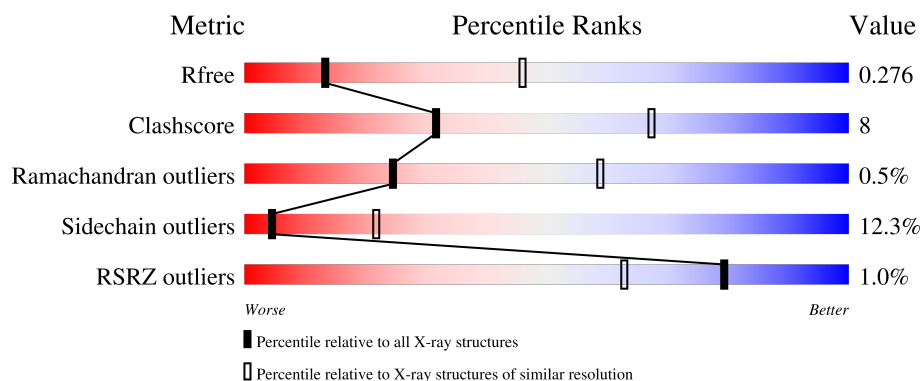
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	421	
1	B	421	
1	C	421	
1	D	421	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12356 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 Transesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	0	0
			3086	1942	555	565	24			
1	B	393	Total	C	N	O	S	0	0	0
			3086	1942	555	565	24			
1	C	393	Total	C	N	O	S	0	0	0
			3086	1942	555	565	24			
1	D	393	Total	C	N	O	S	0	0	0
			3086	1942	555	565	24			

There are 156 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	expression tag	UNP Q9Y7D1
A	0	VAL	-	expression tag	UNP Q9Y7D1
A	4	ASN	ILE	engineered mutation	UNP Q9Y7D1
A	9	VAL	ALA	engineered mutation	UNP Q9Y7D1
A	26	GLU	LYS	engineered mutation	UNP Q9Y7D1
A	28	SER	ARG	engineered mutation	UNP Q9Y7D1
A	35	LEU	ILE	engineered mutation	UNP Q9Y7D1
A	40	ALA	CYS	engineered mutation	UNP Q9Y7D1
A	43	ARG	ASN	engineered mutation	UNP Q9Y7D1
A	60	ASN	CYS	engineered mutation	UNP Q9Y7D1
A	96	ARG	ASP	engineered mutation	UNP Q9Y7D1
A	109	CYS	SER	engineered mutation	UNP Q9Y7D1
A	123	PRO	ALA	engineered mutation	UNP Q9Y7D1
A	157	VAL	MET	engineered mutation	UNP Q9Y7D1
A	164	GLY	SER	engineered mutation	UNP Q9Y7D1
A	172	ASN	SER	engineered mutation	UNP Q9Y7D1
A	174	PHE	LEU	engineered mutation	UNP Q9Y7D1
A	178	LEU	ALA	engineered mutation	UNP Q9Y7D1
A	191	GLY	ASN	engineered mutation	UNP Q9Y7D1
A	192	ILE	LEU	engineered mutation	UNP Q9Y7D1
A	241	MET	GLN	engineered mutation	UNP Q9Y7D1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	247	SER	ALA	engineered mutation	UNP Q9Y7D1
A	250	LYS	ARG	engineered mutation	UNP Q9Y7D1
A	256	THR	SER	engineered mutation	UNP Q9Y7D1
A	261	HIS	ALA	engineered mutation	UNP Q9Y7D1
A	275	SER	GLY	engineered mutation	UNP Q9Y7D1
A	297	GLY	GLN	engineered mutation	UNP Q9Y7D1
A	335	MET	LEU	engineered mutation	UNP Q9Y7D1
A	361	MET	LEU	engineered mutation	UNP Q9Y7D1
A	370	ILE	VAL	engineered mutation	UNP Q9Y7D1
A	383	VAL	ALA	engineered mutation	UNP Q9Y7D1
A	391	SER	ASN	engineered mutation	UNP Q9Y7D1
A	404	LYS	HIS	engineered mutation	UNP Q9Y7D1
A	414	HIS	-	expression tag	UNP Q9Y7D1
A	415	HIS	-	expression tag	UNP Q9Y7D1
A	416	HIS	-	expression tag	UNP Q9Y7D1
A	417	HIS	-	expression tag	UNP Q9Y7D1
A	418	HIS	-	expression tag	UNP Q9Y7D1
A	419	HIS	-	expression tag	UNP Q9Y7D1
B	-1	MET	-	expression tag	UNP Q9Y7D1
B	0	VAL	-	expression tag	UNP Q9Y7D1
B	4	ASN	ILE	engineered mutation	UNP Q9Y7D1
B	9	VAL	ALA	engineered mutation	UNP Q9Y7D1
B	26	GLU	LYS	engineered mutation	UNP Q9Y7D1
B	28	SER	ARG	engineered mutation	UNP Q9Y7D1
B	35	LEU	ILE	engineered mutation	UNP Q9Y7D1
B	40	ALA	CYS	engineered mutation	UNP Q9Y7D1
B	43	ARG	ASN	engineered mutation	UNP Q9Y7D1
B	60	ASN	CYS	engineered mutation	UNP Q9Y7D1
B	96	ARG	ASP	engineered mutation	UNP Q9Y7D1
B	109	CYS	SER	engineered mutation	UNP Q9Y7D1
B	123	PRO	ALA	engineered mutation	UNP Q9Y7D1
B	157	VAL	MET	engineered mutation	UNP Q9Y7D1
B	164	GLY	SER	engineered mutation	UNP Q9Y7D1
B	172	ASN	SER	engineered mutation	UNP Q9Y7D1
B	174	PHE	LEU	engineered mutation	UNP Q9Y7D1
B	178	LEU	ALA	engineered mutation	UNP Q9Y7D1
B	191	GLY	ASN	engineered mutation	UNP Q9Y7D1
B	192	ILE	LEU	engineered mutation	UNP Q9Y7D1
B	241	MET	GLN	engineered mutation	UNP Q9Y7D1
B	247	SER	ALA	engineered mutation	UNP Q9Y7D1
B	250	LYS	ARG	engineered mutation	UNP Q9Y7D1
B	256	THR	SER	engineered mutation	UNP Q9Y7D1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	261	HIS	ALA	engineered mutation	UNP Q9Y7D1
B	275	SER	GLY	engineered mutation	UNP Q9Y7D1
B	297	GLY	GLN	engineered mutation	UNP Q9Y7D1
B	335	MET	LEU	engineered mutation	UNP Q9Y7D1
B	361	MET	LEU	engineered mutation	UNP Q9Y7D1
B	370	ILE	VAL	engineered mutation	UNP Q9Y7D1
B	383	VAL	ALA	engineered mutation	UNP Q9Y7D1
B	391	SER	ASN	engineered mutation	UNP Q9Y7D1
B	404	LYS	HIS	engineered mutation	UNP Q9Y7D1
B	414	HIS	-	expression tag	UNP Q9Y7D1
B	415	HIS	-	expression tag	UNP Q9Y7D1
B	416	HIS	-	expression tag	UNP Q9Y7D1
B	417	HIS	-	expression tag	UNP Q9Y7D1
B	418	HIS	-	expression tag	UNP Q9Y7D1
B	419	HIS	-	expression tag	UNP Q9Y7D1
C	-1	MET	-	expression tag	UNP Q9Y7D1
C	0	VAL	-	expression tag	UNP Q9Y7D1
C	4	ASN	ILE	engineered mutation	UNP Q9Y7D1
C	9	VAL	ALA	engineered mutation	UNP Q9Y7D1
C	26	GLU	LYS	engineered mutation	UNP Q9Y7D1
C	28	SER	ARG	engineered mutation	UNP Q9Y7D1
C	35	LEU	ILE	engineered mutation	UNP Q9Y7D1
C	40	ALA	CYS	engineered mutation	UNP Q9Y7D1
C	43	ARG	ASN	engineered mutation	UNP Q9Y7D1
C	60	ASN	CYS	engineered mutation	UNP Q9Y7D1
C	96	ARG	ASP	engineered mutation	UNP Q9Y7D1
C	109	CYS	SER	engineered mutation	UNP Q9Y7D1
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C	157	VAL	MET	engineered mutation	UNP Q9Y7D1
C	164	GLY	SER	engineered mutation	UNP Q9Y7D1
C	172	ASN	SER	engineered mutation	UNP Q9Y7D1
C	174	PHE	LEU	engineered mutation	UNP Q9Y7D1
C	178	LEU	ALA	engineered mutation	UNP Q9Y7D1
C	191	GLY	ASN	engineered mutation	UNP Q9Y7D1
C	192	ILE	LEU	engineered mutation	UNP Q9Y7D1
C	241	MET	GLN	engineered mutation	UNP Q9Y7D1
C	247	SER	ALA	engineered mutation	UNP Q9Y7D1
C	250	LYS	ARG	engineered mutation	UNP Q9Y7D1
C	256	THR	SER	engineered mutation	UNP Q9Y7D1
C	261	HIS	ALA	engineered mutation	UNP Q9Y7D1
C	275	SER	GLY	engineered mutation	UNP Q9Y7D1
C	297	GLY	GLN	engineered mutation	UNP Q9Y7D1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	335	MET	LEU	engineered mutation	UNP Q9Y7D1
C	361	MET	LEU	engineered mutation	UNP Q9Y7D1
C	370	ILE	VAL	engineered mutation	UNP Q9Y7D1
C	383	VAL	ALA	engineered mutation	UNP Q9Y7D1
C	391	SER	ASN	engineered mutation	UNP Q9Y7D1
C	404	LYS	HIS	engineered mutation	UNP Q9Y7D1
C	414	HIS	-	expression tag	UNP Q9Y7D1
C	415	HIS	-	expression tag	UNP Q9Y7D1
C	416	HIS	-	expression tag	UNP Q9Y7D1
C	417	HIS	-	expression tag	UNP Q9Y7D1
C	418	HIS	-	expression tag	UNP Q9Y7D1
C	419	HIS	-	expression tag	UNP Q9Y7D1
D	-1	MET	-	expression tag	UNP Q9Y7D1
D	0	VAL	-	expression tag	UNP Q9Y7D1
D	4	ASN	ILE	engineered mutation	UNP Q9Y7D1
D	9	VAL	ALA	engineered mutation	UNP Q9Y7D1
D	26	GLU	LYS	engineered mutation	UNP Q9Y7D1
D	28	SER	ARG	engineered mutation	UNP Q9Y7D1
D	35	LEU	ILE	engineered mutation	UNP Q9Y7D1
D	40	ALA	CYS	engineered mutation	UNP Q9Y7D1
D	43	ARG	ASN	engineered mutation	UNP Q9Y7D1
D	60	ASN	CYS	engineered mutation	UNP Q9Y7D1
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D	261	HIS	ALA	engineered mutation	UNP Q9Y7D1
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D	297	GLY	GLN	engineered mutation	UNP Q9Y7D1
D	335	MET	LEU	engineered mutation	UNP Q9Y7D1
D	361	MET	LEU	engineered mutation	UNP Q9Y7D1
D	370	ILE	VAL	engineered mutation	UNP Q9Y7D1

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D	391	SER	ASN	engineered mutation	UNP Q9Y7D1
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D	416	HIS	-	expression tag	UNP Q9Y7D1
D	417	HIS	-	expression tag	UNP Q9Y7D1
D	418	HIS	-	expression tag	UNP Q9Y7D1
D	419	HIS	-	expression tag	UNP Q9Y7D1

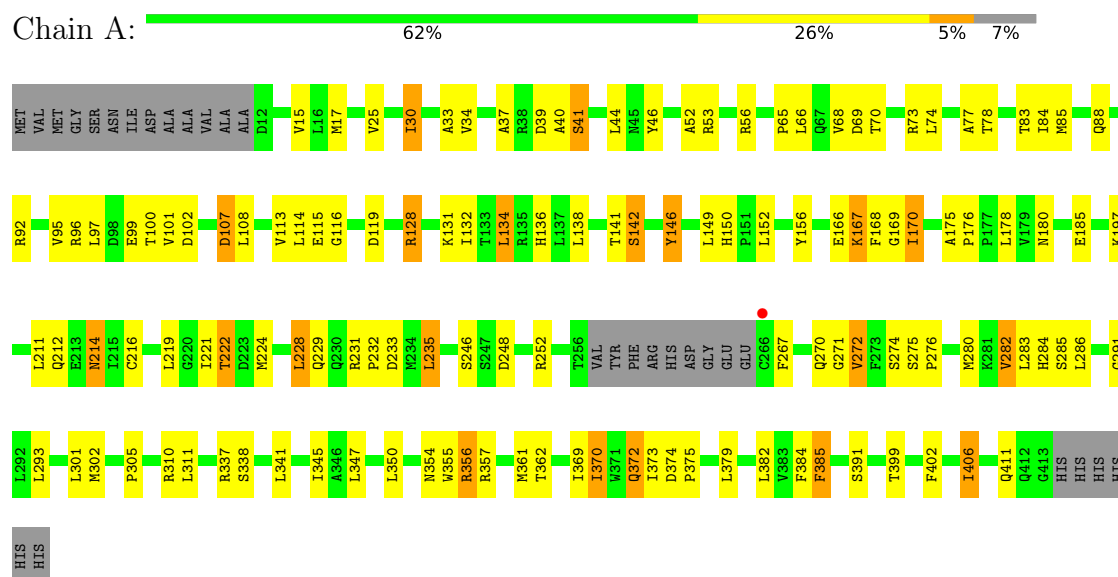
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total O 3 3	0	0
2	B	2	Total O 2 2	0	0
2	C	4	Total O 4 4	0	0
2	D	3	Total O 3 3	0	0

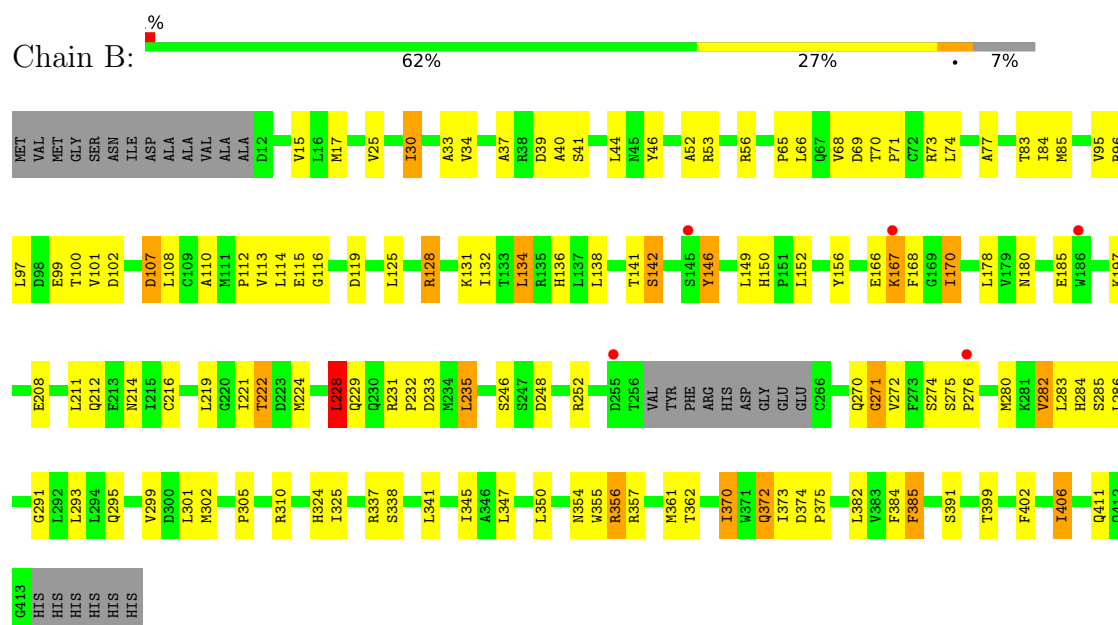
3 Residue-property plots [i](#)

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.

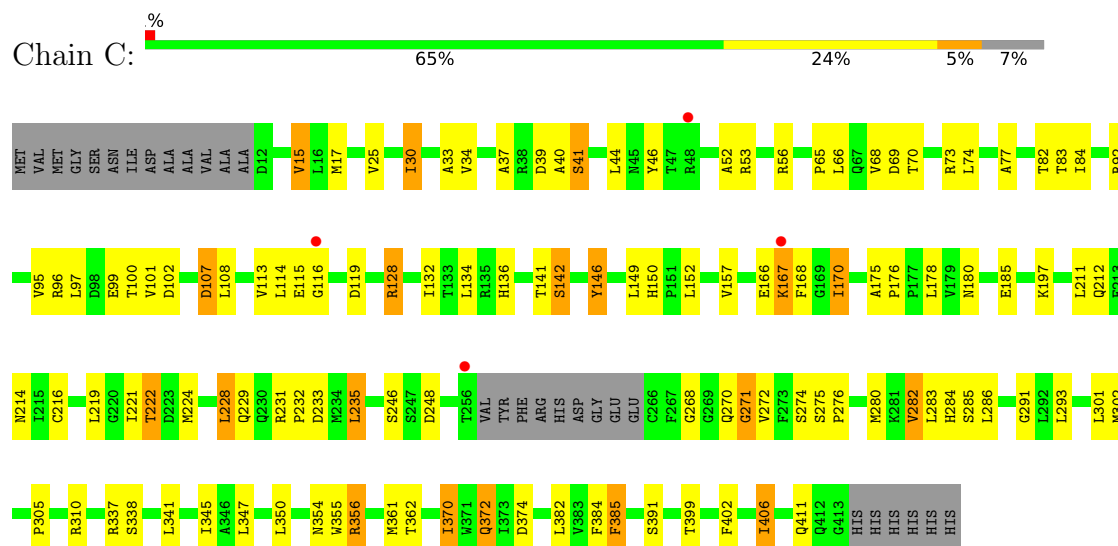
• Molecule 1: Transesterase



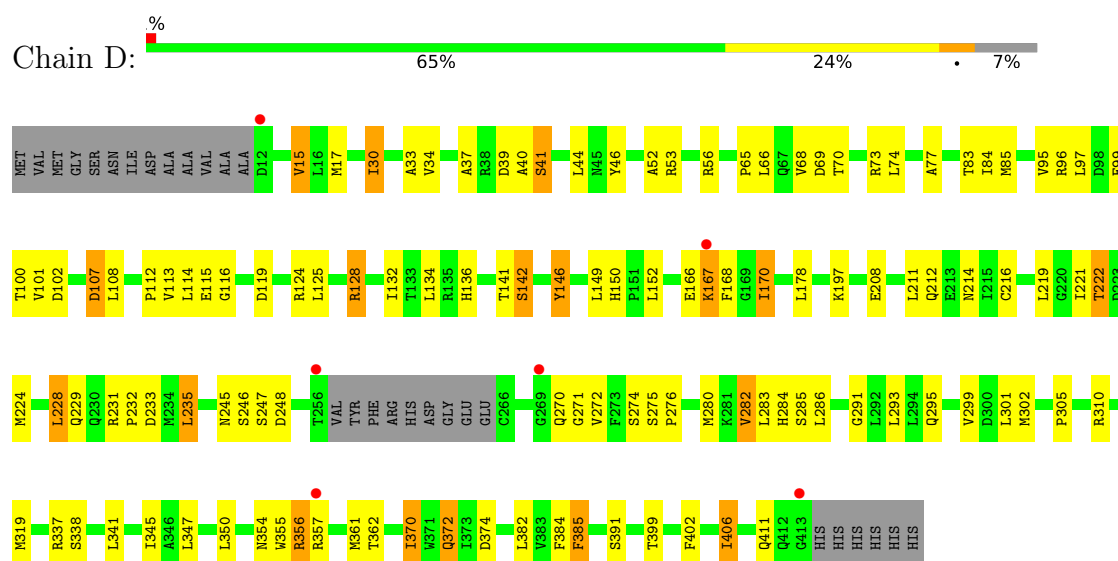
• Molecule 1: Transesterase



- Molecule 1: Transesterase



- Molecule 1: Transesterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.31Å 83.94Å 99.20Å 102.56° 89.02° 107.64°	Depositor
Resolution (Å)	19.66 – 3.19 19.66 – 3.19	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.66-3.19) 92.3 (19.66-3.19)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 3.22Å)	Xtriage
Refinement program	BUSTER-TNT, BUSTER 2.8.0	Depositor
R, R_{free}	0.208 , 0.252 0.226 , 0.276	Depositor DCC
R_{free} test set	1302 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	43.1	Xtriage
Anisotropy	0.556	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 58.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.029 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	12356	wwPDB-VP
Average B, all atoms (Å ²)	50.0	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 23.64 % 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 4.4796e-03. 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

5.1 Standard geometry

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	0.85	0/3152	1.47	34/4264 (0.8%)
1	B	0.83	0/3152	1.46	37/4264 (0.9%)
1	C	0.83	0/3152	1.46	30/4264 (0.7%)
1	D	0.84	1/3152 (0.0%)	1.47	31/4264 (0.7%)
All	All	0.84	1/12608 (0.0%)	1.47	132/17056 (0.8%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	319	MET	SD-CE	-5.17	1.66	1.79

All (132) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	232	PRO	CA-C-N	7.63	130.36	120.44
1	C	232	PRO	C-N-CA	7.63	130.36	120.44
1	D	232	PRO	CA-C-N	7.63	130.36	120.44
1	D	232	PRO	C-N-CA	7.63	130.36	120.44
1	B	232	PRO	CA-C-N	7.57	130.28	120.44
1	B	232	PRO	C-N-CA	7.57	130.28	120.44
1	A	232	PRO	CA-C-N	7.52	130.22	120.44
1	A	232	PRO	C-N-CA	7.52	130.22	120.44
1	C	272	VAL	N-CA-C	7.35	119.21	109.21
1	D	272	VAL	N-CA-C	7.29	119.13	109.21
1	A	272	VAL	N-CA-C	7.13	118.91	109.21
1	B	272	VAL	N-CA-C	7.06	118.81	109.21
1	A	69	ASP	CA-CB-CG	6.77	119.37	112.60
1	B	69	ASP	CA-CB-CG	6.66	119.26	112.60
1	D	69	ASP	CA-CB-CG	6.65	119.25	112.60
1	A	40	ALA	CA-C-N	6.46	129.23	120.38
1	A	40	ALA	C-N-CA	6.46	129.23	120.38
1	A	141	THR	CA-C-N	6.30	128.72	120.28
1	A	141	THR	C-N-CA	6.30	128.72	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	107	ASP	CA-CB-CG	6.20	118.80	112.60
1	C	69	ASP	CA-CB-CG	6.19	118.79	112.60
1	B	385	PHE	CA-CB-CG	6.13	119.94	113.80
1	C	40	ALA	CA-C-N	6.12	128.77	120.38
1	C	40	ALA	C-N-CA	6.12	128.77	120.38
1	B	141	THR	CA-C-N	6.11	128.47	120.28
1	B	141	THR	C-N-CA	6.11	128.47	120.28
1	C	141	THR	CA-C-N	6.10	128.45	120.28
1	C	141	THR	C-N-CA	6.10	128.45	120.28
1	D	141	THR	CA-C-N	6.06	128.40	120.28
1	D	141	THR	C-N-CA	6.06	128.40	120.28
1	D	40	ALA	CA-C-N	6.03	128.65	120.38
1	D	40	ALA	C-N-CA	6.03	128.65	120.38
1	C	385	PHE	CA-CB-CG	6.03	119.83	113.80
1	B	40	ALA	CA-C-N	6.01	128.61	120.38
1	B	40	ALA	C-N-CA	6.01	128.61	120.38
1	B	107	ASP	CA-CB-CG	5.94	118.54	112.60
1	D	146	TYR	N-CA-C	-5.93	99.73	109.40
1	A	107	ASP	CA-CB-CG	5.88	118.48	112.60
1	D	385	PHE	CA-CB-CG	5.88	119.68	113.80
1	A	214	ASN	N-CA-C	5.85	120.70	113.50
1	C	146	TYR	N-CA-C	-5.81	99.92	109.40
1	C	102	ASP	CA-C-N	5.80	128.31	120.65
1	C	102	ASP	C-N-CA	5.80	128.31	120.65
1	C	41	SER	N-CA-C	5.73	118.27	111.33
1	C	107	ASP	CA-CB-CG	5.68	118.28	112.60
1	C	214	ASN	N-CA-C	5.68	120.49	113.50
1	A	385	PHE	CA-CB-CG	5.67	119.47	113.80
1	B	102	ASP	CA-C-N	5.64	128.10	120.65
1	B	102	ASP	C-N-CA	5.64	128.10	120.65
1	B	41	SER	N-CA-C	5.63	118.14	111.33
1	A	146	TYR	N-CA-C	-5.62	100.24	109.40
1	A	102	ASP	CA-C-N	5.58	128.02	120.65
1	A	102	ASP	C-N-CA	5.58	128.02	120.65
1	B	146	TYR	N-CA-C	-5.57	100.32	109.40
1	D	102	ASP	CA-C-N	5.55	127.98	120.65
1	D	102	ASP	C-N-CA	5.55	127.98	120.65
1	A	115	GLU	N-CA-C	-5.55	104.63	111.40
1	D	115	GLU	N-CA-C	-5.54	104.64	111.40
1	D	214	ASN	N-CA-C	5.51	120.28	113.50
1	D	41	SER	N-CA-C	5.51	117.99	111.33
1	D	291	GLY	CA-C-N	5.49	127.58	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	291	GLY	C-N-CA	5.49	127.58	120.44
1	B	115	GLU	N-CA-C	-5.48	104.71	111.40
1	A	131	LYS	N-CA-C	5.47	118.61	110.52
1	C	115	GLU	N-CA-C	-5.44	104.76	111.40
1	A	246	SER	CA-C-N	5.44	131.77	121.81
1	A	246	SER	C-N-CA	5.44	131.77	121.81
1	A	41	SER	N-CA-C	5.42	117.89	111.33
1	D	246	SER	CA-C-N	5.41	131.71	121.81
1	D	246	SER	C-N-CA	5.41	131.71	121.81
1	B	246	SER	CA-C-N	5.41	131.71	121.81
1	B	246	SER	C-N-CA	5.41	131.71	121.81
1	A	101	VAL	CA-C-N	5.41	131.87	121.54
1	A	101	VAL	C-N-CA	5.41	131.87	121.54
1	C	246	SER	CA-C-N	5.40	131.69	121.81
1	C	246	SER	C-N-CA	5.40	131.69	121.81
1	B	101	VAL	CA-C-N	5.40	131.85	121.54
1	B	101	VAL	C-N-CA	5.40	131.85	121.54
1	C	101	VAL	CA-C-N	5.37	131.80	121.54
1	C	101	VAL	C-N-CA	5.37	131.80	121.54
1	B	214	ASN	N-CA-C	5.35	120.08	113.50
1	B	291	GLY	CA-C-N	5.33	127.37	120.44
1	B	291	GLY	C-N-CA	5.33	127.37	120.44
1	D	101	VAL	CA-C-N	5.31	131.69	121.54
1	D	101	VAL	C-N-CA	5.31	131.69	121.54
1	D	284	HIS	CA-C-N	5.29	127.32	120.44
1	D	284	HIS	C-N-CA	5.29	127.32	120.44
1	A	284	HIS	CA-C-N	5.26	127.28	120.44
1	A	284	HIS	C-N-CA	5.26	127.28	120.44
1	B	85	MET	CA-C-N	5.25	127.32	120.28
1	B	85	MET	C-N-CA	5.25	127.32	120.28
1	A	291	GLY	CA-C-N	5.25	127.26	120.44
1	A	291	GLY	C-N-CA	5.25	127.26	120.44
1	B	357	ARG	CA-C-N	5.24	128.29	120.90
1	B	357	ARG	C-N-CA	5.24	128.29	120.90
1	D	282	VAL	CA-C-N	5.22	127.28	120.28
1	D	282	VAL	C-N-CA	5.22	127.28	120.28
1	B	284	HIS	CA-C-N	5.21	127.21	120.44
1	B	284	HIS	C-N-CA	5.21	127.21	120.44
1	D	208	GLU	CB-CG-CD	5.19	121.42	112.60
1	C	284	HIS	CA-C-N	5.19	127.18	120.44
1	C	284	HIS	C-N-CA	5.19	127.18	120.44
1	C	282	VAL	CA-C-N	5.18	127.23	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	282	VAL	C-N-CA	5.18	127.23	120.28
1	B	131	LYS	N-CA-C	5.18	118.45	110.42
1	D	85	MET	CA-C-N	5.16	127.19	120.28
1	D	85	MET	C-N-CA	5.16	127.19	120.28
1	B	208	GLU	CB-CG-CD	5.16	121.36	112.60
1	C	25	VAL	CA-C-N	5.15	127.14	120.44
1	C	25	VAL	C-N-CA	5.15	127.14	120.44
1	B	25	VAL	CA-C-N	5.13	127.11	120.44
1	B	25	VAL	C-N-CA	5.13	127.11	120.44
1	B	271	GLY	N-CA-C	5.12	125.31	113.18
1	A	25	VAL	CA-C-N	5.11	127.08	120.44
1	A	25	VAL	C-N-CA	5.11	127.08	120.44
1	C	271	GLY	N-CA-C	5.10	125.27	113.18
1	A	357	ARG	CA-C-N	5.10	128.09	120.90
1	A	357	ARG	C-N-CA	5.10	128.09	120.90
1	A	85	MET	CA-C-N	5.09	127.11	120.28
1	A	85	MET	C-N-CA	5.09	127.11	120.28
1	B	282	VAL	CA-C-N	5.09	127.11	120.28
1	B	282	VAL	C-N-CA	5.09	127.11	120.28
1	D	357	ARG	CA-C-N	5.09	128.07	120.90
1	D	357	ARG	C-N-CA	5.09	128.07	120.90
1	C	157	VAL	N-CA-C	5.08	115.85	110.72
1	A	369	ILE	N-CA-CB	5.07	117.21	110.26
1	A	282	VAL	CA-C-N	5.07	127.08	120.28
1	A	282	VAL	C-N-CA	5.07	127.08	120.28
1	C	291	GLY	CA-C-N	5.07	127.03	120.44
1	C	291	GLY	C-N-CA	5.07	127.03	120.44
1	B	228	LEU	CA-C-N	5.01	127.69	120.38
1	B	228	LEU	C-N-CA	5.01	127.69	120.38

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	3086	0	3069	55	0
1	B	3086	0	3069	53	0
1	C	3086	0	3069	48	0
1	D	3086	0	3069	48	0
2	A	3	0	0	0	0
2	B	2	0	0	0	0
2	C	4	0	0	0	0
2	D	3	0	0	0	0
All	All	12356	0	12276	200	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 8.

All (200) 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:96:ARG:HB3	1:A:99:GLU:HB2	1.56	0.87
1:C:96:ARG:HB3	1:C:99:GLU:HB2	1.56	0.87
1:B:96:ARG:HB3	1:B:99:GLU:HB2	1.56	0.86
1:D:96:ARG:HB3	1:D:99:GLU:HB2	1.57	0.85
1:C:15:VAL:HG11	1:D:15:VAL:HG11	1.59	0.82
1:D:146:TYR:HB2	1:D:149:LEU:HD12	1.72	0.72
1:C:146:TYR:HB2	1:C:149:LEU:HD12	1.72	0.71
1:A:146:TYR:HB2	1:A:149:LEU:HD12	1.73	0.70
1:B:146:TYR:HB2	1:B:149:LEU:HD12	1.73	0.69
1:D:52:ALA:HA	1:D:65:PRO:HA	1.78	0.65
1:C:52:ALA:HA	1:C:65:PRO:HA	1.78	0.64
1:B:52:ALA:HA	1:B:65:PRO:HA	1.78	0.64
1:A:52:ALA:HA	1:A:65:PRO:HA	1.78	0.64
1:D:37:ALA:HB3	1:D:46:TYR:HB3	1.84	0.60
1:B:37:ALA:HB3	1:B:46:TYR:HB3	1.84	0.59
1:C:37:ALA:HB3	1:C:46:TYR:HB3	1.85	0.59
1:D:68:VAL:HA	1:D:276:PRO:HG2	1.83	0.59
1:C:68:VAL:HA	1:C:276:PRO:HG2	1.83	0.59
1:B:68:VAL:HA	1:B:276:PRO:HG2	1.83	0.59
1:A:37:ALA:HB3	1:A:46:TYR:HB3	1.85	0.58
1:A:68:VAL:HA	1:A:276:PRO:HG2	1.83	0.58
1:D:167:LYS:HE3	1:D:168:PHE:CZ	2.39	0.58
1:A:167:LYS:HE3	1:A:168:PHE:CZ	2.39	0.58
1:C:167:LYS:HE3	1:C:168:PHE:CZ	2.39	0.57
1:A:74:LEU:HB3	1:A:77:ALA:HB3	1.87	0.56
1:B:167:LYS:HE3	1:B:168:PHE:CZ	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:SER:HB2	1:A:347:LEU:HD11	1.88	0.56
1:D:74:LEU:HB3	1:D:77:ALA:HB3	1.87	0.56
1:B:74:LEU:HB3	1:B:77:ALA:HB3	1.88	0.56
1:D:219:LEU:HD11	1:D:293:LEU:HD11	1.88	0.55
1:C:74:LEU:HB3	1:C:77:ALA:HB3	1.88	0.55
1:C:356:ARG:HH22	1:C:362:THR:HG21	1.72	0.55
1:B:168:PHE:HB2	1:B:170:ILE:HG12	1.89	0.55
1:D:356:ARG:HH22	1:D:362:THR:HG21	1.72	0.55
1:C:219:LEU:HD11	1:C:293:LEU:HD11	1.88	0.54
1:D:168:PHE:HB2	1:D:170:ILE:HG12	1.88	0.54
1:B:356:ARG:HH22	1:B:362:THR:HG21	1.73	0.54
1:D:338:SER:HB2	1:D:347:LEU:HD11	1.89	0.54
1:B:338:SER:HB2	1:B:347:LEU:HD11	1.89	0.53
1:C:338:SER:HB2	1:C:347:LEU:HD11	1.89	0.53
1:B:219:LEU:HD11	1:B:293:LEU:HD11	1.88	0.53
1:D:372:GLN:HE22	1:D:374:ASP:HB2	1.74	0.53
1:A:372:GLN:HE22	1:A:374:ASP:HB2	1.74	0.53
1:B:372:GLN:HE22	1:B:374:ASP:HB2	1.74	0.53
1:D:355:TRP:CD1	1:D:356:ARG:H	2.27	0.52
1:A:128:ARG:HG3	1:A:128:ARG:HH11	1.75	0.52
1:A:168:PHE:HB2	1:A:170:ILE:HG12	1.91	0.52
1:A:219:LEU:HD11	1:A:293:LEU:HD11	1.89	0.52
1:A:356:ARG:HH22	1:A:362:THR:HG21	1.73	0.52
1:A:355:TRP:CD1	1:A:356:ARG:H	2.27	0.52
1:B:128:ARG:HH11	1:B:128:ARG:HG3	1.75	0.52
1:B:355:TRP:CD1	1:B:356:ARG:H	2.27	0.52
1:C:128:ARG:NH1	1:C:132:ILE:HD11	2.25	0.52
1:C:355:TRP:CD1	1:C:356:ARG:H	2.27	0.52
1:C:372:GLN:HE22	1:C:374:ASP:HB2	1.75	0.51
1:B:128:ARG:NH1	1:B:132:ILE:HD11	2.25	0.51
1:C:128:ARG:HH11	1:C:128:ARG:HG3	1.75	0.51
1:B:33:ALA:HB2	1:B:385:PHE:CE1	2.45	0.51
1:C:17:MET:HG3	1:C:402:PHE:HA	1.92	0.51
1:C:33:ALA:HB2	1:C:385:PHE:CE1	2.45	0.51
1:D:128:ARG:NH1	1:D:132:ILE:HD11	2.26	0.51
1:B:305:PRO:HB3	1:B:338:SER:HB3	1.93	0.51
1:A:128:ARG:NH1	1:A:132:ILE:HD11	2.25	0.51
1:B:17:MET:HG3	1:B:402:PHE:HA	1.93	0.51
1:D:33:ALA:HB2	1:D:385:PHE:CE1	2.45	0.51
1:A:17:MET:HG3	1:A:402:PHE:HA	1.93	0.51
1:A:33:ALA:HB2	1:A:385:PHE:CE1	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:MET:HA	1:A:283:LEU:HD12	1.93	0.50
1:B:280:MET:HA	1:B:283:LEU:HD12	1.94	0.50
1:D:228:LEU:HD13	1:D:235:LEU:HG	1.93	0.50
1:C:228:LEU:HD13	1:C:235:LEU:HG	1.94	0.50
1:A:305:PRO:HB3	1:A:338:SER:HB3	1.94	0.50
1:C:280:MET:HA	1:C:283:LEU:HD12	1.93	0.50
1:D:17:MET:HG3	1:D:402:PHE:HA	1.93	0.49
1:C:168:PHE:HB2	1:C:170:ILE:HG12	1.93	0.49
1:D:350:LEU:HD22	1:D:356:ARG:HB2	1.94	0.49
1:D:128:ARG:HG3	1:D:128:ARG:HH11	1.76	0.49
1:D:280:MET:HA	1:D:283:LEU:HD12	1.94	0.49
1:B:228:LEU:HD13	1:B:235:LEU:HG	1.94	0.49
1:B:74:LEU:HD21	1:B:274:SER:HB3	1.95	0.48
1:C:212:GLN:HA	1:C:216:CYS:HB2	1.95	0.48
1:C:350:LEU:HD22	1:C:356:ARG:HB2	1.94	0.48
1:B:212:GLN:HA	1:B:216:CYS:HB2	1.95	0.48
1:C:305:PRO:HB3	1:C:338:SER:HB3	1.94	0.48
1:B:350:LEU:HD22	1:B:356:ARG:HB2	1.94	0.48
1:D:212:GLN:HA	1:D:216:CYS:HB2	1.96	0.48
1:A:350:LEU:HD22	1:A:356:ARG:HB2	1.95	0.48
1:D:305:PRO:HB3	1:D:338:SER:HB3	1.94	0.48
1:A:283:LEU:HA	1:A:286:LEU:HD12	1.95	0.48
1:B:150:HIS:HE1	1:B:152:LEU:HD12	1.78	0.48
1:B:283:LEU:HA	1:B:286:LEU:HD12	1.95	0.48
1:D:283:LEU:HA	1:D:286:LEU:HD12	1.95	0.48
1:A:74:LEU:HD21	1:A:274:SER:HB3	1.96	0.47
1:A:228:LEU:HD13	1:A:235:LEU:HG	1.95	0.47
1:C:283:LEU:HA	1:C:286:LEU:HD12	1.94	0.47
1:D:116:GLY:HA2	1:D:152:LEU:HD22	1.97	0.47
1:A:169:GLY:HA3	1:B:110:ALA:O	2.14	0.47
1:C:70:THR:O	1:C:275:SER:HA	2.15	0.47
1:D:150:HIS:HE1	1:D:152:LEU:HD12	1.78	0.47
1:A:212:GLN:HA	1:A:216:CYS:HB2	1.97	0.47
1:C:74:LEU:HD21	1:C:274:SER:HB3	1.96	0.47
1:D:70:THR:O	1:D:275:SER:HA	2.15	0.47
1:A:150:HIS:HE1	1:A:152:LEU:HD12	1.80	0.46
1:A:70:THR:O	1:A:275:SER:HA	2.14	0.46
1:C:150:HIS:HE1	1:C:152:LEU:HD12	1.80	0.46
1:A:116:GLY:HA2	1:A:152:LEU:HD22	1.96	0.46
1:B:116:GLY:HA2	1:B:152:LEU:HD22	1.97	0.46
1:B:222:THR:O	1:B:231:ARG:NH1	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:THR:O	1:A:231:ARG:NH1	2.49	0.46
1:A:108:LEU:O	1:A:128:ARG:NH2	2.49	0.46
1:A:34:VAL:HG23	1:A:66:LEU:HD23	1.98	0.46
1:D:74:LEU:HD21	1:D:274:SER:HB3	1.97	0.46
1:B:30:ILE:HD13	1:B:385:PHE:HE1	1.81	0.45
1:D:34:VAL:HG23	1:D:66:LEU:HD23	1.97	0.45
1:D:84:ILE:HD11	1:D:286:LEU:HD11	1.97	0.45
1:B:34:VAL:HG23	1:B:66:LEU:HD23	1.98	0.45
1:A:355:TRP:CG	1:A:356:ARG:H	2.35	0.45
1:A:382:LEU:HD13	1:A:384:PHE:CD1	2.51	0.45
1:C:34:VAL:HG23	1:C:66:LEU:HD23	1.99	0.45
1:C:116:GLY:HA2	1:C:152:LEU:HD22	1.98	0.45
1:C:222:THR:O	1:C:231:ARG:NH1	2.49	0.45
1:C:302:MET:HB3	1:C:361:MET:HE1	1.99	0.45
1:D:302:MET:HB3	1:D:361:MET:HE1	1.98	0.45
1:B:302:MET:HB3	1:B:361:MET:HE1	1.99	0.45
1:B:108:LEU:O	1:B:128:ARG:NH2	2.50	0.45
1:B:70:THR:O	1:B:275:SER:HA	2.16	0.45
1:B:382:LEU:HD13	1:B:384:PHE:CD1	2.52	0.45
1:B:355:TRP:CG	1:B:356:ARG:H	2.35	0.45
1:C:382:LEU:HD13	1:C:384:PHE:CD1	2.52	0.45
1:B:84:ILE:HD11	1:B:286:LEU:HD11	1.97	0.44
1:C:108:LEU:O	1:C:128:ARG:NH2	2.50	0.44
1:C:355:TRP:CG	1:C:356:ARG:H	2.35	0.44
1:D:355:TRP:CG	1:D:356:ARG:H	2.35	0.44
1:A:30:ILE:HD13	1:A:385:PHE:HE1	1.83	0.44
1:A:136:HIS:HB3	1:A:142:SER:HA	1.99	0.44
1:D:222:THR:O	1:D:231:ARG:NH1	2.50	0.44
1:B:370:ILE:HG21	1:B:399:THR:HG21	1.98	0.44
1:D:382:LEU:HD13	1:D:384:PHE:CD1	2.52	0.44
1:A:370:ILE:HG21	1:A:399:THR:HG21	1.98	0.44
1:C:30:ILE:HD13	1:C:385:PHE:HE1	1.82	0.44
1:D:108:LEU:O	1:D:128:ARG:NH2	2.51	0.44
1:C:82:THR:OG1	1:C:268:GLY:O	2.33	0.44
1:C:136:HIS:HB3	1:C:142:SER:HA	1.99	0.44
1:D:211:LEU:HD21	1:D:224:MET:HG2	1.99	0.44
1:A:44:LEU:HD11	1:A:406:ILE:HD12	2.00	0.43
1:B:211:LEU:HD21	1:B:224:MET:HG2	2.00	0.43
1:C:44:LEU:HD11	1:C:406:ILE:HD12	2.00	0.43
1:D:136:HIS:HB3	1:D:142:SER:HA	1.99	0.43
1:A:282:VAL:O	1:A:285:SER:HB3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:MET:HB3	1:A:361:MET:HE1	1.99	0.43
1:B:136:HIS:HB3	1:B:142:SER:HA	2.00	0.43
1:A:180:ASN:HD22	1:A:185:GLU:HG2	1.84	0.43
1:A:211:LEU:HD21	1:A:224:MET:HG2	2.00	0.43
1:D:370:ILE:HG21	1:D:399:THR:HG21	2.00	0.43
1:A:167:LYS:HD2	1:B:156:TYR:OH	2.19	0.43
1:D:30:ILE:HD13	1:D:385:PHE:HE1	1.82	0.43
1:D:44:LEU:HD11	1:D:406:ILE:HD12	2.01	0.43
1:C:211:LEU:HD21	1:C:224:MET:HG2	2.00	0.43
1:D:282:VAL:O	1:D:285:SER:HB3	2.19	0.43
1:D:216:CYS:HA	1:D:221:ILE:HB	2.01	0.43
1:C:84:ILE:HD11	1:C:286:LEU:HD11	2.00	0.43
1:B:282:VAL:O	1:B:285:SER:HB3	2.19	0.42
1:B:356:ARG:NH2	1:B:362:THR:HG21	2.34	0.42
1:C:370:ILE:HG21	1:C:399:THR:HG21	2.00	0.42
1:A:39:ASP:HB3	1:A:44:LEU:HB3	2.01	0.42
1:D:356:ARG:NH2	1:D:362:THR:HG21	2.34	0.42
1:C:39:ASP:OD2	1:C:41:SER:OG	2.36	0.42
1:A:156:TYR:OH	1:B:167:LYS:HD2	2.20	0.42
1:C:180:ASN:HD22	1:C:185:GLU:HG2	1.84	0.42
1:A:84:ILE:HD11	1:A:286:LEU:HD11	2.00	0.42
1:B:39:ASP:HB3	1:B:44:LEU:HB3	2.02	0.42
1:D:39:ASP:HB3	1:D:44:LEU:HB3	2.02	0.42
1:A:39:ASP:OD2	1:A:41:SER:OG	2.34	0.42
1:C:39:ASP:HB3	1:C:44:LEU:HB3	2.02	0.42
1:D:39:ASP:OD2	1:D:41:SER:OG	2.35	0.42
1:A:216:CYS:HA	1:A:221:ILE:HB	2.01	0.41
1:C:356:ARG:NH2	1:C:362:THR:HG21	2.34	0.41
1:B:44:LEU:HD11	1:B:406:ILE:HD12	2.02	0.41
1:C:216:CYS:HA	1:C:221:ILE:HB	2.02	0.41
1:A:175:ALA:N	1:A:176:PRO:HD3	2.36	0.41
1:A:78:THR:HG23	1:A:272:VAL:HB	2.03	0.41
1:A:185:GLU:HA	1:A:311:LEU:HD22	2.03	0.41
1:A:356:ARG:NH2	1:A:362:THR:HG21	2.35	0.41
1:C:175:ALA:N	1:C:176:PRO:HD3	2.36	0.41
1:C:282:VAL:O	1:C:285:SER:HB3	2.20	0.41
1:D:112:PRO:HB2	1:D:125:LEU:HB3	2.02	0.41
1:B:324:HIS:CD2	1:B:325:ILE:HG13	2.56	0.41
1:A:88:GLN:OE1	1:A:214:ASN:HB3	2.22	0.40
1:A:373:ILE:O	1:A:375:PRO:HD3	2.20	0.40
1:B:216:CYS:HA	1:B:221:ILE:HB	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:LEU:HD22	1:A:138:LEU:HD11	2.03	0.40
1:B:70:THR:HA	1:B:71:PRO:HD3	2.00	0.40
1:B:134:LEU:HD22	1:B:138:LEU:HD11	2.04	0.40
1:B:180:ASN:HD22	1:B:185:GLU:HG2	1.85	0.40
1:B:295:GLN:O	1:B:299:VAL:HG23	2.22	0.40
1:B:373:ILE:O	1:B:375:PRO:HD3	2.22	0.40
1:A:44:LEU:HD22	1:A:379:LEU:CD2	2.51	0.40
1:B:112:PRO:HB2	1:B:125:LEU:HB3	2.03	0.40
1:D:245:ASN:C	1:D:247:SER:H	2.30	0.40
1:D:295:GLN:O	1:D:299:VAL:HG23	2.22	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	389/421 (92%)	358 (92%)	29 (8%)	2 (0%)	24	59
1	B	389/421 (92%)	358 (92%)	29 (8%)	2 (0%)	24	59
1	C	389/421 (92%)	358 (92%)	29 (8%)	2 (0%)	24	59
1	D	389/421 (92%)	359 (92%)	28 (7%)	2 (0%)	24	59
All	All	1556/1684 (92%)	1433 (92%)	115 (7%)	8 (0%)	24	59

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	271	GLY
1	B	271	GLY
1	C	271	GLY
1	D	271	GLY
1	A	356	ARG

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Mol	Chain	Res	Type
1	B	356	ARG
1	C	356	ARG
1	D	356	ARG

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	330/352 (94%)	288 (87%)	42 (13%)	4	20
1	B	330/352 (94%)	290 (88%)	40 (12%)	5	22
1	C	330/352 (94%)	290 (88%)	40 (12%)	5	22
1	D	330/352 (94%)	290 (88%)	40 (12%)	5	22
All	All	1320/1408 (94%)	1158 (88%)	162 (12%)	4	22

All (162) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	VAL
1	A	30	ILE
1	A	53	ARG
1	A	56	ARG
1	A	73	ARG
1	A	83	THR
1	A	92	ARG
1	A	95	VAL
1	A	97	LEU
1	A	100	THR
1	A	107	ASP
1	A	113	VAL
1	A	114	LEU
1	A	119	ASP
1	A	128	ARG
1	A	134	LEU
1	A	142	SER
1	A	166	GLU

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Mol	Chain	Res	Type
1	A	167	LYS
1	A	170	ILE
1	A	178	LEU
1	A	197	LYS
1	A	222	THR
1	A	228	LEU
1	A	229	GLN
1	A	233	ASP
1	A	235	LEU
1	A	248	ASP
1	A	252	ARG
1	A	267	PHE
1	A	270	GLN
1	A	301	LEU
1	A	310	ARG
1	A	337	ARG
1	A	341	LEU
1	A	345	ILE
1	A	354	ASN
1	A	370	ILE
1	A	372	GLN
1	A	391	SER
1	A	406	ILE
1	A	411	GLN
1	B	15	VAL
1	B	30	ILE
1	B	53	ARG
1	B	56	ARG
1	B	73	ARG
1	B	83	THR
1	B	95	VAL
1	B	97	LEU
1	B	100	THR
1	B	107	ASP
1	B	113	VAL
1	B	114	LEU
1	B	119	ASP
1	B	128	ARG
1	B	134	LEU
1	B	142	SER
1	B	166	GLU
1	B	167	LYS

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Mol	Chain	Res	Type
1	B	170	ILE
1	B	178	LEU
1	B	197	LYS
1	B	222	THR
1	B	228	LEU
1	B	229	GLN
1	B	233	ASP
1	B	235	LEU
1	B	248	ASP
1	B	252	ARG
1	B	270	GLN
1	B	301	LEU
1	B	310	ARG
1	B	337	ARG
1	B	341	LEU
1	B	345	ILE
1	B	354	ASN
1	B	370	ILE
1	B	372	GLN
1	B	391	SER
1	B	406	ILE
1	B	411	GLN
1	C	15	VAL
1	C	30	ILE
1	C	53	ARG
1	C	56	ARG
1	C	73	ARG
1	C	83	THR
1	C	92	ARG
1	C	95	VAL
1	C	97	LEU
1	C	100	THR
1	C	107	ASP
1	C	113	VAL
1	C	114	LEU
1	C	119	ASP
1	C	128	ARG
1	C	134	LEU
1	C	142	SER
1	C	166	GLU
1	C	167	LYS
1	C	170	ILE

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Mol	Chain	Res	Type
1	C	178	LEU
1	C	197	LYS
1	C	222	THR
1	C	228	LEU
1	C	229	GLN
1	C	233	ASP
1	C	235	LEU
1	C	248	ASP
1	C	270	GLN
1	C	301	LEU
1	C	310	ARG
1	C	337	ARG
1	C	341	LEU
1	C	345	ILE
1	C	354	ASN
1	C	370	ILE
1	C	372	GLN
1	C	391	SER
1	C	406	ILE
1	C	411	GLN
1	D	15	VAL
1	D	30	ILE
1	D	53	ARG
1	D	56	ARG
1	D	73	ARG
1	D	83	THR
1	D	95	VAL
1	D	97	LEU
1	D	100	THR
1	D	107	ASP
1	D	113	VAL
1	D	114	LEU
1	D	119	ASP
1	D	124	ARG
1	D	128	ARG
1	D	134	LEU
1	D	142	SER
1	D	166	GLU
1	D	167	LYS
1	D	170	ILE
1	D	178	LEU
1	D	197	LYS

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Mol	Chain	Res	Type
1	D	222	THR
1	D	228	LEU
1	D	229	GLN
1	D	233	ASP
1	D	235	LEU
1	D	248	ASP
1	D	270	GLN
1	D	301	LEU
1	D	310	ARG
1	D	337	ARG
1	D	341	LEU
1	D	345	ILE
1	D	354	ASN
1	D	370	ILE
1	D	372	GLN
1	D	391	SER
1	D	406	ILE
1	D	411	GLN

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

Mol	Chain	Res	Type
1	A	60	ASN
1	A	67	GLN
1	A	150	HIS
1	A	171	GLN
1	A	172	ASN
1	A	230	GLN
1	A	245	ASN
1	A	295	GLN
1	A	318	HIS
1	A	354	ASN
1	A	368	ASN
1	A	372	GLN
1	B	60	ASN
1	B	67	GLN
1	B	150	HIS
1	B	171	GLN
1	B	172	ASN
1	B	230	GLN
1	B	295	GLN
1	B	318	HIS

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Mol	Chain	Res	Type
1	B	354	ASN
1	B	372	GLN
1	C	60	ASN
1	C	67	GLN
1	C	150	HIS
1	C	171	GLN
1	C	172	ASN
1	C	212	GLN
1	C	230	GLN
1	C	245	ASN
1	C	295	GLN
1	C	318	HIS
1	C	354	ASN
1	C	372	GLN
1	D	60	ASN
1	D	67	GLN
1	D	150	HIS
1	D	171	GLN
1	D	172	ASN
1	D	212	GLN
1	D	230	GLN
1	D	245	ASN
1	D	295	GLN
1	D	317	GLN
1	D	318	HIS
1	D	354	ASN
1	D	372	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	393/421 (93%)	0.23	1 (0%) 90 81	14, 42, 75, 105	0
1	B	393/421 (93%)	0.38	5 (1%) 75 55	18, 51, 84, 115	0
1	C	393/421 (93%)	0.32	4 (1%) 79 63	22, 49, 87, 114	0
1	D	393/421 (93%)	0.31	6 (1%) 72 52	19, 46, 76, 116	0
All	All	1572/1684 (93%)	0.31	16 (1%) 79 63	14, 48, 81, 116	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	413	GLY	3.0
1	D	167	LYS	2.9
1	C	48	ARG	2.6
1	B	276	PRO	2.5
1	B	255	ASP	2.5
1	B	186	TRP	2.4
1	C	256	THR	2.3
1	D	256	THR	2.3
1	C	167	LYS	2.3
1	B	167	LYS	2.3
1	B	145	SER	2.2
1	D	357	ARG	2.1
1	D	12	ASP	2.1
1	C	116	GLY	2.1
1	A	266	CYS	2.0
1	D	269	GLY	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 [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.