



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:05 AM UTC

PDB ID : 3ZLH / pdb_00003zlh
Title : Structure of group A Streptococcal enolase
Authors : Cork, A.J.; Ericsson, D.J.; Law, R.H.P.; Casey, L.W.; Valkov, E.; Bertozzi, C.; Stamp, A.; Aquilina, J.A.; Whisstock, J.C.; Walker, M.J.; Kobe, B.
Deposited on : 2013-01-31
Resolution : 2.90 Å(reported)

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

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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

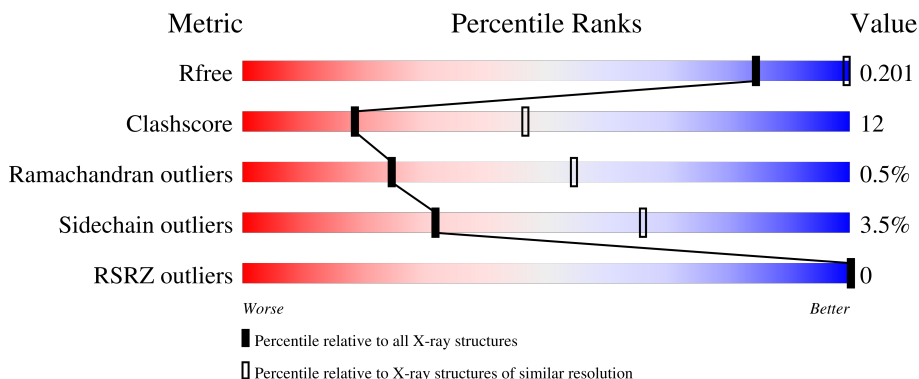
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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

Mol	Chain	Length	Quality of chain
1	A	455	 69% 22% • 5%
1	B	455	 71% 21% • 5%
1	C	455	 68% 22% • 7%
1	D	455	 69% 24% • 5%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 13172 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 ENOLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	0	0
			3302	2081	555	655	11			
1	B	432	Total	C	N	O	S	0	0	0
			3311	2086	557	657	11			
1	C	425	Total	C	N	O	S	0	0	0
			3254	2053	544	646	11			
1	D	432	Total	C	N	O	S	0	0	0
			3305	2082	555	657	11			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q5XD01
A	-18	GLY	-	expression tag	UNP Q5XD01
A	-17	SER	-	expression tag	UNP Q5XD01
A	-16	SER	-	expression tag	UNP Q5XD01
A	-15	HIS	-	expression tag	UNP Q5XD01
A	-14	HIS	-	expression tag	UNP Q5XD01
A	-13	HIS	-	expression tag	UNP Q5XD01
A	-12	HIS	-	expression tag	UNP Q5XD01
A	-11	HIS	-	expression tag	UNP Q5XD01
A	-10	HIS	-	expression tag	UNP Q5XD01
A	-9	SER	-	expression tag	UNP Q5XD01
A	-8	SER	-	expression tag	UNP Q5XD01
A	-7	GLY	-	expression tag	UNP Q5XD01
A	-6	LEU	-	expression tag	UNP Q5XD01
A	-5	VAL	-	expression tag	UNP Q5XD01
A	-4	PRO	-	expression tag	UNP Q5XD01
A	-3	ARG	-	expression tag	UNP Q5XD01
A	-2	GLY	-	expression tag	UNP Q5XD01
A	-1	SER	-	expression tag	UNP Q5XD01
A	0	HIS	-	expression tag	UNP Q5XD01
B	-19	MET	-	expression tag	UNP Q5XD01

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP Q5XD01
B	-17	SER	-	expression tag	UNP Q5XD01
B	-16	SER	-	expression tag	UNP Q5XD01
B	-15	HIS	-	expression tag	UNP Q5XD01
B	-14	HIS	-	expression tag	UNP Q5XD01
B	-13	HIS	-	expression tag	UNP Q5XD01
B	-12	HIS	-	expression tag	UNP Q5XD01
B	-11	HIS	-	expression tag	UNP Q5XD01
B	-10	HIS	-	expression tag	UNP Q5XD01
B	-9	SER	-	expression tag	UNP Q5XD01
B	-8	SER	-	expression tag	UNP Q5XD01
B	-7	GLY	-	expression tag	UNP Q5XD01
B	-6	LEU	-	expression tag	UNP Q5XD01
B	-5	VAL	-	expression tag	UNP Q5XD01
B	-4	PRO	-	expression tag	UNP Q5XD01
B	-3	ARG	-	expression tag	UNP Q5XD01
B	-2	GLY	-	expression tag	UNP Q5XD01
B	-1	SER	-	expression tag	UNP Q5XD01
B	0	HIS	-	expression tag	UNP Q5XD01
C	-19	MET	-	expression tag	UNP Q5XD01
C	-18	GLY	-	expression tag	UNP Q5XD01
C	-17	SER	-	expression tag	UNP Q5XD01
C	-16	SER	-	expression tag	UNP Q5XD01
C	-15	HIS	-	expression tag	UNP Q5XD01
C	-14	HIS	-	expression tag	UNP Q5XD01
C	-13	HIS	-	expression tag	UNP Q5XD01
C	-12	HIS	-	expression tag	UNP Q5XD01
C	-11	HIS	-	expression tag	UNP Q5XD01
C	-10	HIS	-	expression tag	UNP Q5XD01
C	-9	SER	-	expression tag	UNP Q5XD01
C	-8	SER	-	expression tag	UNP Q5XD01
C	-7	GLY	-	expression tag	UNP Q5XD01
C	-6	LEU	-	expression tag	UNP Q5XD01
C	-5	VAL	-	expression tag	UNP Q5XD01
C	-4	PRO	-	expression tag	UNP Q5XD01
C	-3	ARG	-	expression tag	UNP Q5XD01
C	-2	GLY	-	expression tag	UNP Q5XD01
C	-1	SER	-	expression tag	UNP Q5XD01
C	0	HIS	-	expression tag	UNP Q5XD01
D	-19	MET	-	expression tag	UNP Q5XD01
D	-18	GLY	-	expression tag	UNP Q5XD01
D	-17	SER	-	expression tag	UNP Q5XD01

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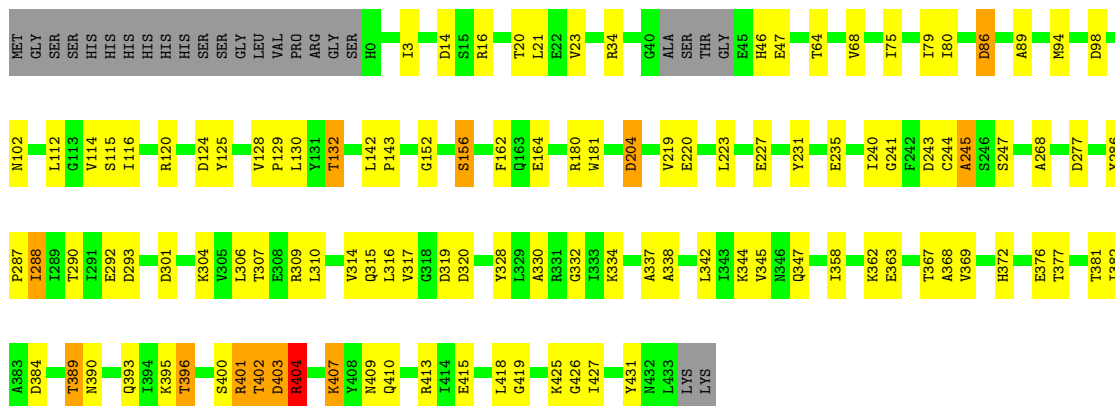
Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP Q5XD01
D	-15	HIS	-	expression tag	UNP Q5XD01
D	-14	HIS	-	expression tag	UNP Q5XD01
D	-13	HIS	-	expression tag	UNP Q5XD01
D	-12	HIS	-	expression tag	UNP Q5XD01
D	-11	HIS	-	expression tag	UNP Q5XD01
D	-10	HIS	-	expression tag	UNP Q5XD01
D	-9	SER	-	expression tag	UNP Q5XD01
D	-8	SER	-	expression tag	UNP Q5XD01
D	-7	GLY	-	expression tag	UNP Q5XD01
D	-6	LEU	-	expression tag	UNP Q5XD01
D	-5	VAL	-	expression tag	UNP Q5XD01
D	-4	PRO	-	expression tag	UNP Q5XD01
D	-3	ARG	-	expression tag	UNP Q5XD01
D	-2	GLY	-	expression tag	UNP Q5XD01
D	-1	SER	-	expression tag	UNP Q5XD01
D	0	HIS	-	expression tag	UNP Q5XD01

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.

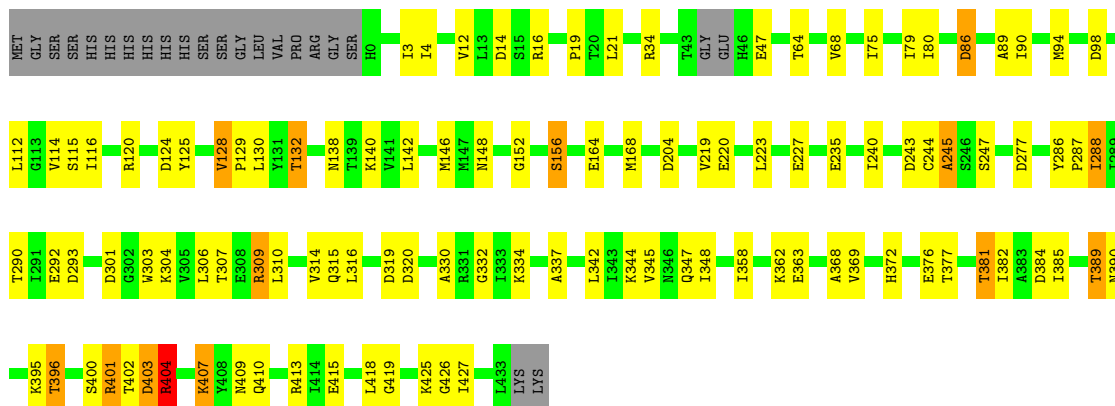
• Molecule 1: ENOLASE

Chain A: 



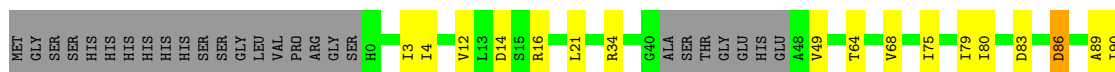
• Molecule 1: ENOLASE

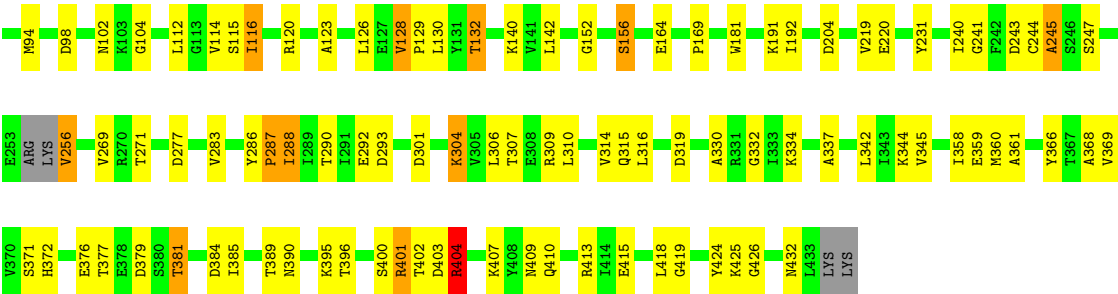
Chain B: 



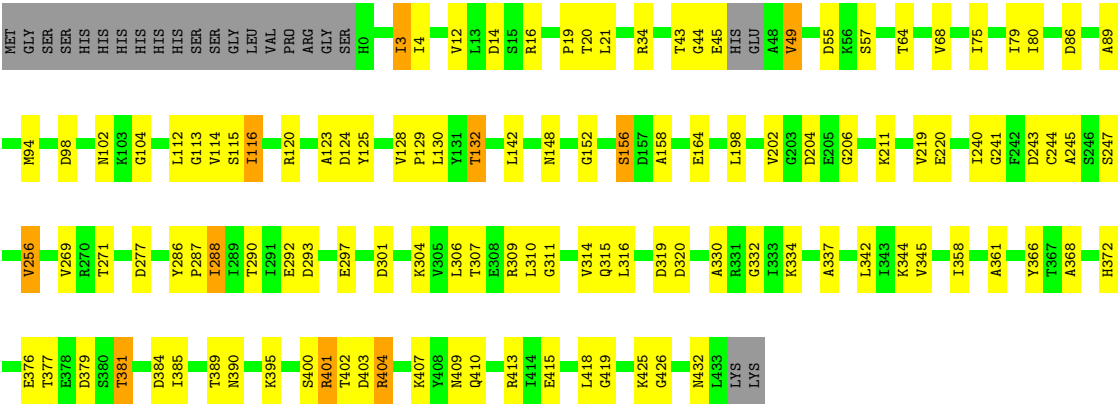
• Molecule 1: ENOLASE

Chain C: 





● Molecule 1: ENOLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	185.79Å 185.79Å 56.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.75 – 2.90 19.75 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.3 (19.75-2.90) 99.7 (19.75-2.90)	Depositor EDS
R_{merge}	0.43	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.88Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.153 , 0.205 0.150 , 0.201	Depositor DCC
R_{free} test set	2166 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	45.5	Xtriage
Anisotropy	0.259	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 0.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.18$	Xtriage
Estimated twinning fraction	0.407 for h,-k,-l	Xtriage
Reported twinning fraction	0.545 for H, K, L 0.455 for K, H, -L	Depositor
Outliers	0 of 43076 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13172	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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](#)

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	1.13	5/3355 (0.1%)	1.19	13/4534 (0.3%)
1	B	1.14	3/3364 (0.1%)	1.20	13/4547 (0.3%)
1	C	1.11	2/3305 (0.1%)	1.20	12/4467 (0.3%)
1	D	1.13	3/3357 (0.1%)	1.19	12/4537 (0.3%)
All	All	1.13	13/13381 (0.1%)	1.19	50/18085 (0.3%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	319	ASP	CA-C	-6.71	1.46	1.53
1	D	319	ASP	CA-C	-6.12	1.47	1.53
1	A	319	ASP	CA-C	-5.69	1.47	1.53
1	A	403	ASP	N-CA	5.62	1.53	1.46
1	A	407	LYS	CA-C	-5.45	1.45	1.52
1	D	377	THR	CA-C	5.44	1.59	1.52
1	B	319	ASP	CA-C	-5.33	1.48	1.53
1	A	396	THR	N-CA	5.25	1.52	1.46
1	A	143	PRO	N-CA	-5.22	1.41	1.47
1	D	20	THR	C-O	-5.19	1.20	1.24
1	B	348	ILE	N-CA	5.15	1.51	1.45
1	B	396	THR	N-CA	5.11	1.52	1.46
1	C	304	LYS	N-CA	5.00	1.52	1.46

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	396	THR	CB-CA-C	-8.46	97.92	111.28
1	B	396	THR	CB-CA-C	-8.45	97.93	111.28
1	D	404	ARG	N-CA-C	-6.63	103.86	112.68
1	A	404	ARG	N-CA-C	-6.30	104.30	112.68
1	A	21	LEU	N-CA-C	6.23	118.65	109.24
1	D	4	ILE	CB-CA-C	-6.22	103.80	111.08
1	C	381	THR	CB-CA-C	-5.97	99.58	110.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	404	ARG	N-CA-C	-5.97	104.74	112.68
1	B	21	LEU	N-CA-C	5.95	118.44	109.23
1	D	123	ALA	N-CA-C	-5.92	104.95	111.82
1	D	381	THR	CB-CA-C	-5.89	99.73	110.63
1	D	20	THR	N-CA-C	-5.89	100.11	108.23
1	B	381	THR	CB-CA-C	-5.88	101.02	110.79
1	A	320	ASP	N-CA-C	-5.82	106.18	113.28
1	B	245	ALA	N-CA-C	-5.76	103.30	111.24
1	A	393	GLN	N-CA-C	5.73	118.11	109.23
1	C	21	LEU	N-CA-C	5.71	118.07	109.23
1	B	128	VAL	CA-C-N	-5.70	114.39	120.03
1	B	128	VAL	C-N-CA	-5.70	114.39	120.03
1	D	3	ILE	N-CA-C	5.68	116.94	109.21
1	B	309	ARG	N-CA-C	5.67	118.31	111.40
1	B	4	ILE	CB-CA-C	-5.66	104.45	111.08
1	A	369	VAL	CB-CA-C	-5.58	105.89	111.80
1	C	123	ALA	N-CA-C	-5.54	105.40	111.82
1	A	381	THR	CB-CA-C	-5.50	101.67	110.79
1	A	245	ALA	N-CA-C	-5.38	103.81	111.24
1	A	204	ASP	N-CA-C	5.38	118.64	111.75
1	C	4	ILE	CB-CA-C	-5.38	104.79	111.08
1	D	21	LEU	N-CA-C	5.35	117.52	109.23
1	B	389	THR	CB-CA-C	-5.34	100.86	110.37
1	A	20	THR	N-CA-C	-5.32	100.88	108.23
1	A	162	PHE	N-CA-C	-5.30	102.01	109.96
1	C	83	ASP	N-CA-C	-5.27	101.88	110.20
1	D	320	ASP	N-CA-C	-5.27	106.70	113.23
1	D	202	VAL	N-CA-C	5.26	116.15	108.53
1	D	198	LEU	CA-C-N	5.25	128.44	120.77
1	D	198	LEU	C-N-CA	5.25	128.44	120.77
1	C	369	VAL	CB-CA-C	-5.24	106.24	111.80
1	C	245	ALA	N-CA-C	-5.20	103.84	111.17
1	B	320	ASP	N-CA-C	-5.18	106.81	113.23
1	C	371	SER	N-CA-C	5.14	117.68	109.81
1	A	268	ALA	N-CA-C	5.07	116.75	108.63
1	B	404	ARG	N-CA-C	-5.07	105.94	112.68
1	C	283	VAL	N-CA-C	5.07	115.79	110.62
1	B	369	VAL	CB-CA-C	-5.06	106.43	111.80
1	B	407	LYS	CA-CB-CG	-5.05	104.00	114.10
1	C	128	VAL	CA-C-N	-5.05	115.03	120.03
1	C	128	VAL	C-N-CA	-5.05	115.03	120.03
1	D	49	VAL	N-CA-CB	5.05	117.17	110.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	389	THR	CB-CA-C	-5.01	101.45	110.37

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	3302	0	3254	78	1
1	B	3311	0	3265	74	1
1	C	3254	0	3208	78	0
1	D	3305	0	3261	85	0
All	All	13172	0	12988	313	1

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 12.

All (313) 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:204:ASP:O	1:B:401:ARG:NH1	1.71	1.22
1:B:307:THR:HG21	1:B:337:ALA:O	1.77	0.84
1:D:307:THR:HG21	1:D:337:ALA:O	1.82	0.80
1:C:307:THR:HG21	1:C:337:ALA:O	1.82	0.79
1:D:44:GLY:O	1:D:297:GLU:OE1	1.99	0.79
1:B:138:ASN:OD1	1:B:140:LYS:NZ	2.14	0.77
1:A:128:VAL:CG1	1:A:129:PRO:HD2	2.16	0.76
1:B:75:ILE:HG21	1:B:114:VAL:HG21	1.69	0.75
1:A:307:THR:HG21	1:A:337:ALA:O	1.87	0.75
1:A:204:ASP:O	1:A:401:ARG:NH1	2.21	0.73
1:C:75:ILE:HG21	1:C:114:VAL:HG21	1.71	0.72
1:D:75:ILE:HG21	1:D:114:VAL:HG21	1.71	0.72
1:B:204:ASP:C	1:B:401:ARG:NH1	2.47	0.72
1:A:75:ILE:HG21	1:A:114:VAL:HG21	1.73	0.71
1:B:376:GLU:CG	1:B:407:LYS:HZ2	2.04	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:376:GLU:HG2	1:A:407:LYS:HZ2	1.59	0.67
1:A:376:GLU:CG	1:A:407:LYS:HZ2	2.06	0.67
1:A:128:VAL:HG13	1:A:129:PRO:HD2	1.75	0.67
1:B:376:GLU:HG2	1:B:407:LYS:HZ2	1.60	0.66
1:B:382:ILE:HD11	1:B:396:THR:HG22	1.77	0.66
1:C:14:ASP:OD1	1:C:14:ASP:C	2.38	0.66
1:D:307:THR:HG21	1:D:337:ALA:C	2.21	0.66
1:A:382:ILE:HD11	1:A:396:THR:HG22	1.77	0.65
1:D:128:VAL:CG1	1:D:129:PRO:CD	2.76	0.63
1:A:128:VAL:CG1	1:A:129:PRO:CD	2.77	0.63
1:D:128:VAL:HG12	1:D:132:THR:OG1	2.00	0.62
1:C:307:THR:HG21	1:C:337:ALA:C	2.24	0.62
1:D:372:HIS:CD2	1:D:404:ARG:NH1	2.67	0.61
1:D:128:VAL:CG1	1:D:129:PRO:HD2	2.29	0.61
1:C:376:GLU:HG3	1:C:407:LYS:HZ1	1.66	0.61
1:A:14:ASP:OD1	1:A:14:ASP:C	2.43	0.61
1:D:286:TYR:HB2	1:D:288:ILE:HD12	1.83	0.61
1:C:286:TYR:HB2	1:C:288:ILE:HD12	1.83	0.60
1:C:372:HIS:CD2	1:C:404:ARG:HH11	2.20	0.60
1:C:152:GLY:HA2	1:C:156:SER:HB2	1.84	0.60
1:D:376:GLU:HG3	1:D:407:LYS:HZ1	1.66	0.60
1:A:128:VAL:HG12	1:A:129:PRO:HD2	1.83	0.59
1:A:243:ASP:HA	1:A:292:GLU:HB3	1.84	0.59
1:C:372:HIS:CD2	1:C:404:ARG:NH1	2.70	0.59
1:A:286:TYR:HB2	1:A:288:ILE:HD12	1.83	0.59
1:B:14:ASP:C	1:B:14:ASP:OD1	2.43	0.59
1:D:14:ASP:C	1:D:14:ASP:OD1	2.44	0.59
1:B:307:THR:HG21	1:B:337:ALA:C	2.28	0.58
1:B:372:HIS:CD2	1:B:396:THR:HA	2.37	0.58
1:A:372:HIS:CD2	1:A:396:THR:HA	2.37	0.58
1:B:286:TYR:HB2	1:B:288:ILE:HD12	1.84	0.58
1:A:235:GLU:HG2	1:A:427:ILE:CD1	2.34	0.58
1:B:152:GLY:HA2	1:B:156:SER:HB2	1.85	0.58
1:D:243:ASP:HA	1:D:292:GLU:HB3	1.84	0.58
1:D:372:HIS:CD2	1:D:404:ARG:HH11	2.21	0.58
1:C:181:TRP:CD1	1:C:231:TYR:HH	2.22	0.58
1:C:376:GLU:HG3	1:C:407:LYS:NZ	2.19	0.58
1:D:256:VAL:CG1	1:D:269:VAL:HG13	2.33	0.58
1:B:243:ASP:HA	1:B:292:GLU:HB3	1.85	0.58
1:D:152:GLY:HA2	1:D:156:SER:HB2	1.86	0.58
1:A:152:GLY:HA2	1:A:156:SER:HB2	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:86:ASP:HB3	1:B:89:ALA:HB3	1.86	0.57
1:D:256:VAL:HG11	1:D:269:VAL:CG1	2.35	0.57
1:B:332:GLY:HA2	1:B:337:ALA:HB3	1.86	0.57
1:C:400:SER:O	1:C:401:ARG:HG3	2.05	0.57
1:A:307:THR:HG21	1:A:337:ALA:C	2.30	0.57
1:A:332:GLY:HA2	1:A:337:ALA:HB3	1.87	0.56
1:C:243:ASP:HA	1:C:292:GLU:HB3	1.86	0.56
1:B:376:GLU:HG2	1:B:407:LYS:NZ	2.19	0.56
1:A:86:ASP:HB3	1:A:89:ALA:HB3	1.86	0.56
1:D:86:ASP:HB3	1:D:89:ALA:HB3	1.87	0.56
1:D:128:VAL:HG12	1:D:129:PRO:CD	2.36	0.56
1:D:376:GLU:HG3	1:D:407:LYS:NZ	2.20	0.56
1:A:376:GLU:HG2	1:A:407:LYS:NZ	2.20	0.56
1:B:372:HIS:CE1	1:B:404:ARG:HH11	2.24	0.55
1:B:277:ASP:OD1	1:B:309:ARG:NH1	2.32	0.55
1:C:116:ILE:HG23	1:C:381:THR:HG21	1.88	0.55
1:B:235:GLU:HG2	1:B:427:ILE:HD12	1.87	0.55
1:A:382:ILE:HD11	1:A:396:THR:CG2	2.37	0.55
1:B:382:ILE:HD11	1:B:396:THR:CG2	2.36	0.55
1:C:332:GLY:HA2	1:C:337:ALA:HB3	1.87	0.55
1:D:128:VAL:HG12	1:D:129:PRO:HD2	1.87	0.55
1:D:128:VAL:HG13	1:D:129:PRO:CD	2.37	0.55
1:D:400:SER:O	1:D:401:ARG:HG3	2.07	0.55
1:C:49:VAL:CG2	1:C:104:GLY:HA2	2.37	0.54
1:D:49:VAL:CG2	1:D:104:GLY:HA2	2.37	0.54
1:B:138:ASN:CG	1:B:140:LYS:HZ1	2.09	0.54
1:D:277:ASP:OD1	1:D:309:ARG:NH1	2.33	0.54
1:A:181:TRP:CD1	1:A:231:TYR:HH	2.26	0.54
1:D:400:SER:C	1:D:401:ARG:HG3	2.33	0.54
1:C:277:ASP:OD1	1:C:309:ARG:NH1	2.35	0.53
1:C:306:LEU:CD1	1:C:310:LEU:HD12	2.37	0.53
1:D:332:GLY:HA2	1:D:337:ALA:HB3	1.90	0.53
1:A:277:ASP:OD1	1:A:309:ARG:NH1	2.35	0.53
1:C:256:VAL:CG1	1:C:269:VAL:HG13	2.39	0.53
1:A:128:VAL:HG13	1:A:129:PRO:CD	2.39	0.53
1:B:358:ILE:HG23	1:B:368:ALA:CB	2.39	0.53
1:D:16:ARG:NH1	1:D:204:ASP:OD2	2.40	0.53
1:C:86:ASP:HB3	1:C:89:ALA:HB3	1.90	0.53
1:D:116:ILE:HG23	1:D:381:THR:HG21	1.90	0.53
1:D:409:ASN:O	1:D:410:GLN:C	2.51	0.53
1:B:128:VAL:CG1	1:B:129:PRO:CD	2.88	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:409:ASN:O	1:C:410:GLN:C	2.52	0.52
1:C:400:SER:C	1:C:401:ARG:HG3	2.33	0.52
1:A:372:HIS:CE1	1:A:404:ARG:HH11	2.27	0.52
1:A:16:ARG:NH1	1:A:204:ASP:OD2	2.41	0.52
1:C:16:ARG:NH1	1:C:204:ASP:OD2	2.41	0.52
1:C:128:VAL:CG1	1:C:129:PRO:CD	2.87	0.52
1:C:418:LEU:O	1:C:419:GLY:C	2.53	0.52
1:D:372:HIS:HD2	1:D:404:ARG:NH1	2.08	0.52
1:C:358:ILE:HG23	1:C:368:ALA:CB	2.40	0.51
1:D:128:VAL:CG1	1:D:132:THR:OG1	2.58	0.51
1:D:307:THR:CG2	1:D:337:ALA:O	2.57	0.51
1:A:290:THR:HA	1:A:315:GLN:O	2.10	0.51
1:C:256:VAL:HG11	1:C:269:VAL:CG1	2.40	0.51
1:A:306:LEU:CD1	1:A:310:LEU:HD12	2.41	0.51
1:D:376:GLU:CG	1:D:407:LYS:HZ1	2.24	0.51
1:B:16:ARG:NH1	1:B:204:ASP:OD2	2.44	0.51
1:C:130:LEU:HB3	1:C:384:ASP:OD2	2.11	0.51
1:C:425:LYS:O	1:C:426:GLY:C	2.54	0.51
1:D:358:ILE:HG23	1:D:368:ALA:CB	2.41	0.51
1:A:389:THR:O	1:A:390:ASN:C	2.54	0.51
1:A:128:VAL:HG12	1:A:129:PRO:CD	2.39	0.50
1:C:376:GLU:CG	1:C:407:LYS:NZ	2.74	0.50
1:C:376:GLU:CG	1:C:407:LYS:HZ1	2.23	0.50
1:A:358:ILE:HG23	1:A:368:ALA:CB	2.40	0.50
1:A:409:ASN:O	1:A:410:GLN:C	2.53	0.50
1:C:94:MET:HE1	1:C:115:SER:HB2	1.94	0.50
1:C:244:CYS:HB2	1:C:293:ASP:O	2.11	0.50
1:A:219:VAL:O	1:A:220:GLU:C	2.55	0.50
1:A:94:MET:HE1	1:A:115:SER:HB2	1.92	0.49
1:B:389:THR:O	1:B:390:ASN:C	2.52	0.49
1:C:372:HIS:ND1	1:C:396:THR:HA	2.27	0.49
1:D:306:LEU:CD1	1:D:310:LEU:HD12	2.42	0.49
1:D:376:GLU:CG	1:D:407:LYS:NZ	2.75	0.49
1:A:425:LYS:O	1:A:426:GLY:C	2.55	0.49
1:C:389:THR:O	1:C:390:ASN:C	2.53	0.49
1:D:94:MET:HE1	1:D:115:SER:HB2	1.94	0.49
1:A:130:LEU:HB3	1:A:384:ASP:OD2	2.13	0.49
1:A:307:THR:HA	1:A:316:LEU:HD13	1.94	0.49
1:C:372:HIS:HD2	1:C:404:ARG:NH1	2.10	0.49
1:B:286:TYR:HB3	1:B:287:PRO:HD2	1.95	0.49
1:B:409:ASN:O	1:B:410:GLN:C	2.52	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:410:GLN:O	1:B:413:ARG:HB2	2.12	0.49
1:C:290:THR:HA	1:C:315:GLN:O	2.12	0.49
1:A:400:SER:O	1:A:401:ARG:HG3	2.12	0.49
1:D:290:THR:HA	1:D:315:GLN:O	2.13	0.49
1:A:307:THR:HG21	1:A:338:ALA:HA	1.94	0.49
1:B:244:CYS:HB2	1:B:293:ASP:O	2.13	0.49
1:D:120:ARG:HH11	1:D:381:THR:HG22	1.78	0.49
1:D:244:CYS:HB2	1:D:293:ASP:O	2.13	0.49
1:D:128:VAL:HG13	1:D:129:PRO:HD2	1.94	0.48
1:D:389:THR:O	1:D:390:ASN:C	2.56	0.48
1:A:235:GLU:HG2	1:A:427:ILE:HD12	1.94	0.48
1:C:306:LEU:HD11	1:C:310:LEU:HD12	1.93	0.48
1:B:130:LEU:HB3	1:B:384:ASP:OD2	2.14	0.48
1:C:120:ARG:HH11	1:C:381:THR:HG22	1.78	0.48
1:A:418:LEU:O	1:A:419:GLY:C	2.56	0.48
1:B:94:MET:HE1	1:B:115:SER:HB2	1.95	0.48
1:C:330:ALA:O	1:C:334:LYS:HG3	2.14	0.48
1:A:244:CYS:HB2	1:A:293:ASP:O	2.14	0.48
1:B:204:ASP:C	1:B:401:ARG:HH11	2.13	0.48
1:B:306:LEU:CD1	1:B:310:LEU:HD12	2.44	0.48
1:B:244:CYS:O	1:B:245:ALA:C	2.56	0.48
1:D:128:VAL:HG12	1:D:129:PRO:N	2.28	0.48
1:D:286:TYR:HB3	1:D:287:PRO:HD2	1.96	0.47
1:D:307:THR:HA	1:D:316:LEU:HD13	1.95	0.47
1:D:410:GLN:O	1:D:413:ARG:HB2	2.13	0.47
1:A:128:VAL:HG12	1:A:132:THR:OG1	2.13	0.47
1:B:47:GLU:CD	1:B:347:GLN:HG3	2.39	0.47
1:B:342:LEU:HD23	1:B:344:LYS:HE3	1.96	0.47
1:D:142:LEU:HD11	1:D:415:GLU:HA	1.96	0.47
1:B:372:HIS:CE1	1:B:404:ARG:NH1	2.82	0.47
1:B:330:ALA:O	1:B:334:LYS:HG3	2.13	0.47
1:D:130:LEU:HB3	1:D:384:ASP:OD2	2.14	0.47
1:D:425:LYS:O	1:D:426:GLY:C	2.57	0.47
1:A:244:CYS:O	1:A:245:ALA:C	2.57	0.47
1:A:376:GLU:HG3	1:A:377:THR:H	1.80	0.47
1:C:286:TYR:HB3	1:C:287:PRO:HD2	1.96	0.47
1:C:307:THR:CG2	1:C:337:ALA:O	2.58	0.47
1:B:307:THR:HA	1:B:316:LEU:HD13	1.96	0.47
1:C:410:GLN:O	1:C:413:ARG:HB2	2.14	0.47
1:B:142:LEU:HD11	1:B:415:GLU:HA	1.97	0.47
1:C:142:LEU:HD11	1:C:415:GLU:HA	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:219:VAL:O	1:D:220:GLU:C	2.56	0.47
1:D:342:LEU:HD23	1:D:344:LYS:HE3	1.96	0.47
1:B:164:GLU:OE1	1:B:395:LYS:HE2	2.15	0.47
1:B:290:THR:HA	1:B:315:GLN:O	2.15	0.47
1:C:244:CYS:O	1:C:245:ALA:C	2.57	0.47
1:D:301:ASP:HA	1:D:304:LYS:HD2	1.96	0.47
1:A:372:HIS:CE1	1:A:404:ARG:NH1	2.84	0.46
1:D:330:ALA:O	1:D:334:LYS:HG3	2.15	0.46
1:A:410:GLN:O	1:A:413:ARG:HB2	2.15	0.46
1:C:128:VAL:HG12	1:C:129:PRO:CD	2.46	0.46
1:D:206:GLY:N	1:D:401:ARG:HD3	2.30	0.46
1:D:400:SER:O	1:D:401:ARG:CB	2.64	0.46
1:A:240:ILE:CG2	1:A:241:GLY:N	2.79	0.46
1:C:129:PRO:HD2	1:C:132:THR:OG1	2.16	0.46
1:C:307:THR:HA	1:C:316:LEU:HD13	1.98	0.46
1:C:379:ASP:OD2	1:C:381:THR:HG23	2.16	0.46
1:D:120:ARG:HH11	1:D:381:THR:CG2	2.29	0.46
1:A:34:ARG:O	1:A:120:ARG:HD3	2.15	0.46
1:A:286:TYR:HB3	1:A:287:PRO:HD2	1.97	0.46
1:D:44:GLY:O	1:D:45:GLU:HB2	2.15	0.46
1:C:116:ILE:CG2	1:C:381:THR:HG21	2.45	0.45
1:C:342:LEU:HD23	1:C:344:LYS:HE3	1.96	0.45
1:C:400:SER:O	1:C:401:ARG:CB	2.64	0.45
1:D:164:GLU:OE1	1:D:395:LYS:HE2	2.16	0.45
1:A:330:ALA:O	1:A:334:LYS:HG3	2.15	0.45
1:A:400:SER:C	1:A:401:ARG:HG3	2.41	0.45
1:B:34:ARG:O	1:B:120:ARG:HD3	2.16	0.45
1:B:219:VAL:O	1:B:220:GLU:C	2.59	0.45
1:D:34:ARG:O	1:D:120:ARG:HD3	2.16	0.45
1:A:223:LEU:O	1:A:227:GLU:HG2	2.15	0.45
1:C:34:ARG:O	1:C:120:ARG:HD3	2.17	0.45
1:D:256:VAL:CG1	1:D:269:VAL:CG1	2.94	0.45
1:B:418:LEU:O	1:B:419:GLY:C	2.59	0.45
1:B:128:VAL:HG12	1:B:129:PRO:CD	2.47	0.45
1:B:303:TRP:O	1:B:307:THR:HG23	2.17	0.45
1:A:342:LEU:HD23	1:A:344:LYS:HE3	1.98	0.45
1:D:244:CYS:O	1:D:245:ALA:C	2.60	0.44
1:A:128:VAL:HG12	1:A:129:PRO:N	2.31	0.44
1:B:64:THR:O	1:B:68:VAL:HG23	2.17	0.44
1:A:164:GLU:OE1	1:A:395:LYS:HE2	2.17	0.44
1:B:223:LEU:O	1:B:227:GLU:HG2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:219:VAL:O	1:C:220:GLU:C	2.60	0.44
1:D:116:ILE:CG2	1:D:381:THR:HG21	2.47	0.44
1:A:180:ARG:NH1	1:A:180:ARG:HG2	2.32	0.44
1:B:112:LEU:HD11	1:B:345:VAL:CG1	2.47	0.44
1:B:376:GLU:CG	1:B:407:LYS:NZ	2.77	0.44
1:C:240:ILE:CG2	1:C:241:GLY:N	2.80	0.44
1:D:418:LEU:O	1:D:419:GLY:C	2.61	0.44
1:D:64:THR:O	1:D:68:VAL:HG23	2.18	0.44
1:D:379:ASP:OD2	1:D:381:THR:HG23	2.18	0.44
1:A:47:GLU:CD	1:A:347:GLN:HG3	2.43	0.44
1:D:256:VAL:HG22	1:D:271:THR:HA	1.99	0.44
1:A:124:ASP:O	1:A:125:TYR:C	2.60	0.43
1:D:307:THR:O	1:D:311:GLY:HA3	2.18	0.43
1:A:102:ASN:OD1	1:A:102:ASN:C	2.61	0.43
1:C:376:GLU:HG3	1:C:377:THR:H	1.84	0.43
1:A:400:SER:O	1:A:401:ARG:CB	2.67	0.43
1:B:128:VAL:HG13	1:B:129:PRO:CD	2.49	0.43
1:C:432:ASN:OD1	1:C:432:ASN:N	2.50	0.43
1:B:128:VAL:HG12	1:B:129:PRO:HD2	2.01	0.43
1:D:55:ASP:OD1	1:D:57:SER:OG	2.36	0.43
1:C:128:VAL:HG12	1:C:129:PRO:HD2	2.01	0.43
1:A:376:GLU:CG	1:A:407:LYS:NZ	2.78	0.43
1:B:146:MET:HE2	1:B:168:MET:SD	2.59	0.43
1:C:64:THR:O	1:C:68:VAL:HG23	2.19	0.43
1:A:363:GLU:HA	1:B:362:LYS:O	2.19	0.43
1:B:128:VAL:CG1	1:B:129:PRO:HD2	2.49	0.43
1:A:307:THR:HA	1:A:316:LEU:CD1	2.49	0.43
1:B:79:ILE:O	1:B:80:ILE:C	2.62	0.43
1:C:164:GLU:OE1	1:C:395:LYS:HE2	2.18	0.43
1:B:301:ASP:HA	1:B:304:LYS:HD2	2.01	0.42
1:C:256:VAL:HG22	1:C:271:THR:HA	2.00	0.42
1:C:301:ASP:HA	1:C:304:LYS:HD2	2.00	0.42
1:C:361:ALA:O	1:C:366:TYR:HB2	2.20	0.42
1:B:142:LEU:CD1	1:B:415:GLU:HG3	2.49	0.42
1:B:306:LEU:HD11	1:B:310:LEU:HD12	2.01	0.42
1:D:79:ILE:O	1:D:80:ILE:C	2.62	0.42
1:A:142:LEU:HD11	1:A:415:GLU:HA	2.01	0.42
1:C:120:ARG:HH11	1:C:381:THR:CG2	2.32	0.42
1:C:381:THR:O	1:C:385:ILE:HG13	2.20	0.42
1:A:301:ASP:HA	1:A:304:LYS:HD2	2.01	0.42
1:B:124:ASP:O	1:B:125:TYR:C	2.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:THR:O	1:A:68:VAL:HG23	2.19	0.42
1:B:400:SER:C	1:B:401:ARG:HG3	2.45	0.42
1:B:425:LYS:O	1:B:426:GLY:C	2.60	0.42
1:B:400:SER:O	1:B:401:ARG:CB	2.67	0.42
1:C:102:ASN:OD1	1:C:102:ASN:C	2.62	0.42
1:B:381:THR:O	1:B:385:ILE:HG13	2.20	0.42
1:D:102:ASN:OD1	1:D:102:ASN:C	2.63	0.42
1:B:79:ILE:HD12	1:B:90:ILE:HG23	2.02	0.42
1:D:12:VAL:O	1:D:19:PRO:HA	2.20	0.42
1:C:128:VAL:CG1	1:C:129:PRO:HD2	2.49	0.42
1:D:129:PRO:HD2	1:D:132:THR:OG1	2.20	0.42
1:C:128:VAL:HG13	1:C:129:PRO:CD	2.49	0.41
1:C:292:GLU:O	1:C:293:ASP:C	2.63	0.41
1:D:240:ILE:CG2	1:D:241:GLY:N	2.83	0.41
1:B:219:VAL:HG13	1:B:240:ILE:HD13	2.01	0.41
1:D:306:LEU:HD11	1:D:310:LEU:HD12	2.01	0.41
1:B:129:PRO:HD2	1:B:132:THR:OG1	2.20	0.41
1:A:306:LEU:HD11	1:A:310:LEU:HD12	2.01	0.41
1:D:361:ALA:O	1:D:366:TYR:HB2	2.20	0.41
1:D:124:ASP:O	1:D:125:TYR:C	2.63	0.41
1:A:306:LEU:HD12	1:A:310:LEU:HD12	2.03	0.41
1:A:362:LYS:O	1:B:363:GLU:HA	2.19	0.41
1:A:367:THR:HA	1:A:431:TYR:OH	2.21	0.41
1:B:112:LEU:HD11	1:B:345:VAL:HG11	2.01	0.41
1:D:158:ALA:HB2	1:D:211:LYS:HA	2.03	0.41
1:D:432:ASN:OD1	1:D:432:ASN:N	2.49	0.41
1:B:12:VAL:O	1:B:19:PRO:HA	2.20	0.41
1:C:79:ILE:HD12	1:C:90:ILE:HG23	2.03	0.41
1:D:219:VAL:HG13	1:D:240:ILE:HD13	2.02	0.41
1:D:292:GLU:O	1:D:293:ASP:C	2.61	0.41
1:C:112:LEU:HD11	1:C:345:VAL:CG1	2.50	0.41
1:A:112:LEU:HD11	1:A:345:VAL:CG1	2.50	0.41
1:C:191:LYS:O	1:C:192:ILE:C	2.62	0.41
1:C:307:THR:HA	1:C:316:LEU:CD1	2.51	0.41
1:C:359:GLU:O	1:C:360:MET:C	2.64	0.41
1:D:112:LEU:HD11	1:D:345:VAL:CG1	2.50	0.41
1:D:381:THR:O	1:D:385:ILE:HG13	2.20	0.41
1:A:79:ILE:O	1:A:80:ILE:C	2.62	0.41
1:C:126:LEU:O	1:C:128:VAL:HG23	2.21	0.41
1:C:256:VAL:CG1	1:C:269:VAL:CG1	2.99	0.41
1:A:46:HIS:HB3	1:A:328:TYR:CE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:VAL:CG1	1:A:132:THR:OG1	2.68	0.40
1:A:180:ARG:HG2	1:A:180:ARG:HH11	1.85	0.40
1:B:376:GLU:HG3	1:B:377:THR:H	1.85	0.40
1:D:112:LEU:O	1:D:113:GLY:C	2.64	0.40
1:A:23:VAL:HG11	1:A:114:VAL:HA	2.03	0.40
1:C:79:ILE:O	1:C:80:ILE:C	2.64	0.40
1:D:148:ASN:ND2	1:D:164:GLU:HG2	2.36	0.40
1:B:148:ASN:ND2	1:B:164:GLU:HG2	2.37	0.40
1:D:400:SER:C	1:D:401:ARG:CG	2.94	0.40
1:A:142:LEU:CD1	1:A:415:GLU:HG3	2.51	0.40
1:C:169:PRO:O	1:C:424:TYR:OH	2.36	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:THR:OG1	1:B:403:ASP:OD1[3_555]	2.16	0.04

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	426/455 (94%)	392 (92%)	32 (8%)	2 (0%)	24	54
1	B	428/455 (94%)	390 (91%)	36 (8%)	2 (0%)	24	54
1	C	419/455 (92%)	378 (90%)	37 (9%)	4 (1%)	12	39
1	D	428/455 (94%)	394 (92%)	33 (8%)	1 (0%)	43	72
All	All	1701/1820 (94%)	1554 (91%)	138 (8%)	9 (0%)	24	54

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	86	ASP
1	A	401	ARG
1	B	401	ARG
1	C	86	ASP
1	C	401	ARG
1	D	401	ARG
1	A	86	ASP
1	C	12	VAL
1	C	287	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	342/362 (94%)	330 (96%)	12 (4%)	32	66
1	B	343/362 (95%)	332 (97%)	11 (3%)	34	68
1	C	337/362 (93%)	324 (96%)	13 (4%)	28	63
1	D	342/362 (94%)	330 (96%)	12 (4%)	32	66
All	All	1364/1448 (94%)	1316 (96%)	48 (4%)	32	66

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ILE
1	A	98	ASP
1	A	116	ILE
1	A	132	THR
1	A	156	SER
1	A	247	SER
1	A	288	ILE
1	A	314	VAL
1	A	317	VAL
1	A	402	THR
1	A	403	ASP
1	A	404	ARG
1	B	3	ILE

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Mol	Chain	Res	Type
1	B	98	ASP
1	B	116	ILE
1	B	132	THR
1	B	156	SER
1	B	247	SER
1	B	288	ILE
1	B	314	VAL
1	B	402	THR
1	B	403	ASP
1	B	404	ARG
1	C	3	ILE
1	C	98	ASP
1	C	116	ILE
1	C	132	THR
1	C	140	LYS
1	C	156	SER
1	C	247	SER
1	C	256	VAL
1	C	288	ILE
1	C	314	VAL
1	C	402	THR
1	C	403	ASP
1	C	404	ARG
1	D	3	ILE
1	D	43	THR
1	D	98	ASP
1	D	116	ILE
1	D	132	THR
1	D	156	SER
1	D	247	SER
1	D	256	VAL
1	D	288	ILE
1	D	314	VAL
1	D	402	THR
1	D	403	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	GLN
1	A	372	HIS
1	B	87	GLN

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Mol	Chain	Res	Type
1	B	372	HIS
1	C	87	GLN
1	C	275	GLN
1	D	87	GLN
1	D	423	GLN

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](#)

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	430/455 (94%)	-2.03	0 100 100	27, 46, 72, 104	0
1	B	432/455 (94%)	-2.03	0 100 100	25, 44, 63, 78	0
1	C	425/455 (93%)	-2.04	0 100 100	27, 44, 65, 88	0
1	D	432/455 (94%)	-2.03	0 100 100	26, 43, 63, 94	0
All	All	1719/1820 (94%)	-2.03	0 100 100	25, 44, 66, 104	0

There are no RSRZ outliers to report.

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.