



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 11:10 am BST

PDB ID : 1R52
Title : Crystal structure of the bifunctional chorismate synthase from *Saccharomyces cerevisiae*
Authors : Quevillon-Cheruel, S.; Leulliot, N.; Meyer, P.; Graille, M.; Bremang, M.; Blondeau, K.; Sorel, I.; Poupon, A.; Janin, J.; van Tilbeurgh, H.
Deposited on : 2003-10-09
Resolution : 2.89 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

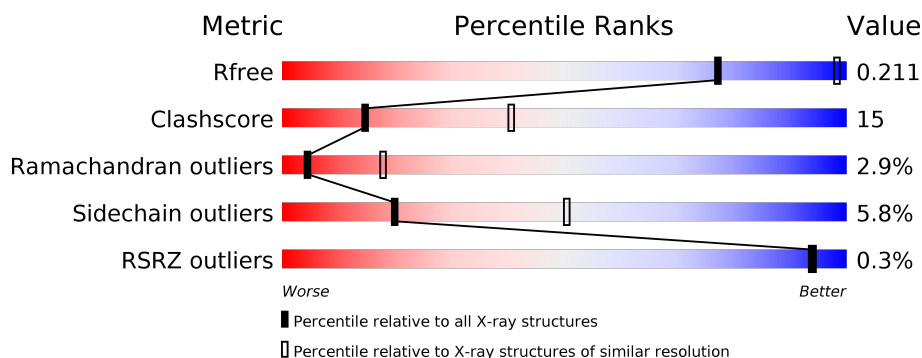
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (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 on 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	382	<div> <div>46%</div> <div>23%</div> <div>5%</div> <div>26%</div> </div>
1	B	382	<div> <div>48%</div> <div>21%</div> <div>• •</div> <div>26%</div> </div>
1	C	382	<div> <div>48%</div> <div>23%</div> <div>• •</div> <div>26%</div> </div>
1	D	382	<div> <div>48%</div> <div>21%</div> <div>• •</div> <div>26%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8636 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 Chorismate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	S	45	0	0
			2138	1347	371	405	15			
1	B	282	Total	C	N	O	S	44	0	0
			2130	1340	373	402	15			
1	C	283	Total	C	N	O	S	46	0	0
			2139	1345	374	405	15			
1	D	282	Total	C	N	O	S	41	0	0
			2131	1341	373	402	15			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	377	HIS	-	EXPRESSION TAG	UNP P28777
A	378	HIS	-	EXPRESSION TAG	UNP P28777
A	379	HIS	-	EXPRESSION TAG	UNP P28777
A	380	HIS	-	EXPRESSION TAG	UNP P28777
A	381	HIS	-	EXPRESSION TAG	UNP P28777
A	382	HIS	-	EXPRESSION TAG	UNP P28777
B	377	HIS	-	EXPRESSION TAG	UNP P28777
B	378	HIS	-	EXPRESSION TAG	UNP P28777
B	379	HIS	-	EXPRESSION TAG	UNP P28777
B	380	HIS	-	EXPRESSION TAG	UNP P28777
B	381	HIS	-	EXPRESSION TAG	UNP P28777
B	382	HIS	-	EXPRESSION TAG	UNP P28777
C	377	HIS	-	EXPRESSION TAG	UNP P28777
C	378	HIS	-	EXPRESSION TAG	UNP P28777
C	379	HIS	-	EXPRESSION TAG	UNP P28777
C	380	HIS	-	EXPRESSION TAG	UNP P28777
C	381	HIS	-	EXPRESSION TAG	UNP P28777
C	382	HIS	-	EXPRESSION TAG	UNP P28777
D	377	HIS	-	EXPRESSION TAG	UNP P28777
D	378	HIS	-	EXPRESSION TAG	UNP P28777
D	379	HIS	-	EXPRESSION TAG	UNP P28777

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Chain	Residue	Modelled	Actual	Comment	Reference
D	380	HIS	-	EXPRESSION TAG	UNP P28777
D	381	HIS	-	EXPRESSION TAG	UNP P28777
D	382	HIS	-	EXPRESSION TAG	UNP P28777

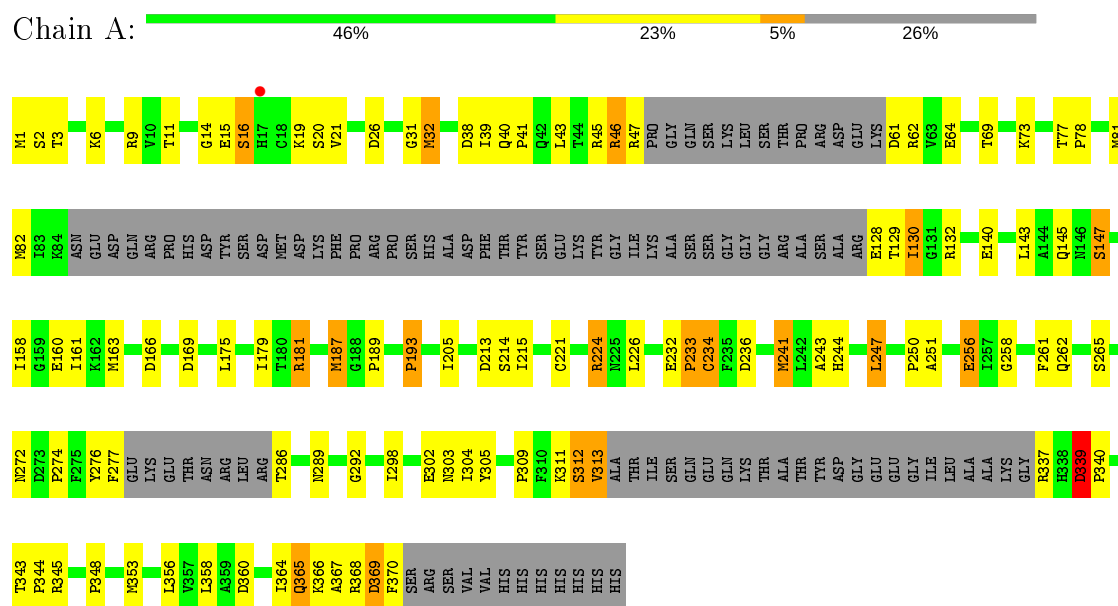
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	25	Total	O	0	0
			25	25		
2	B	26	Total	O	0	0
			26	26		
2	C	27	Total	O	0	0
			27	27		
2	D	20	Total	O	0	0
			20	20		

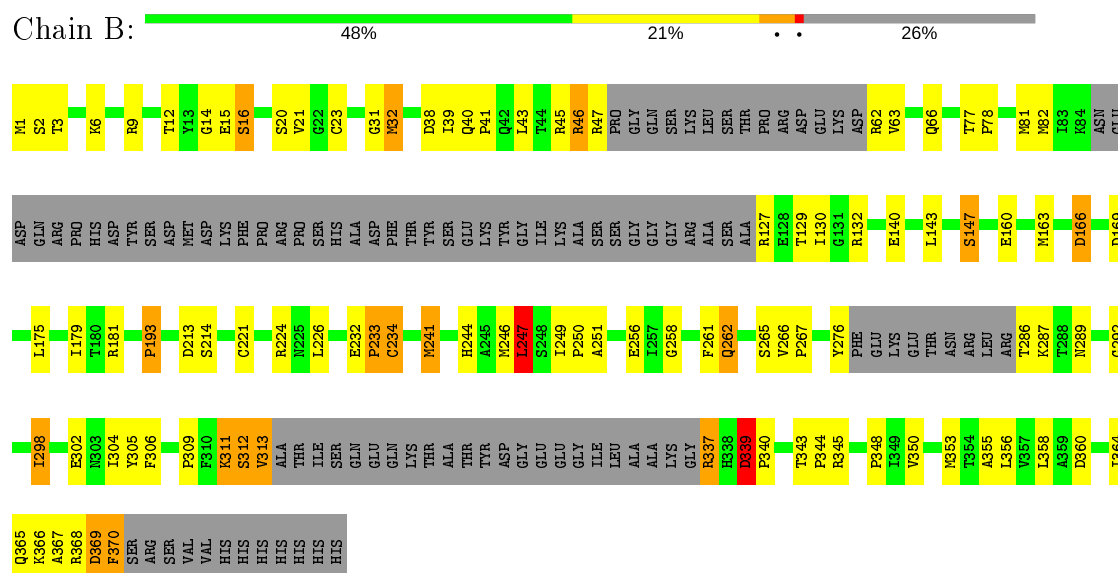
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

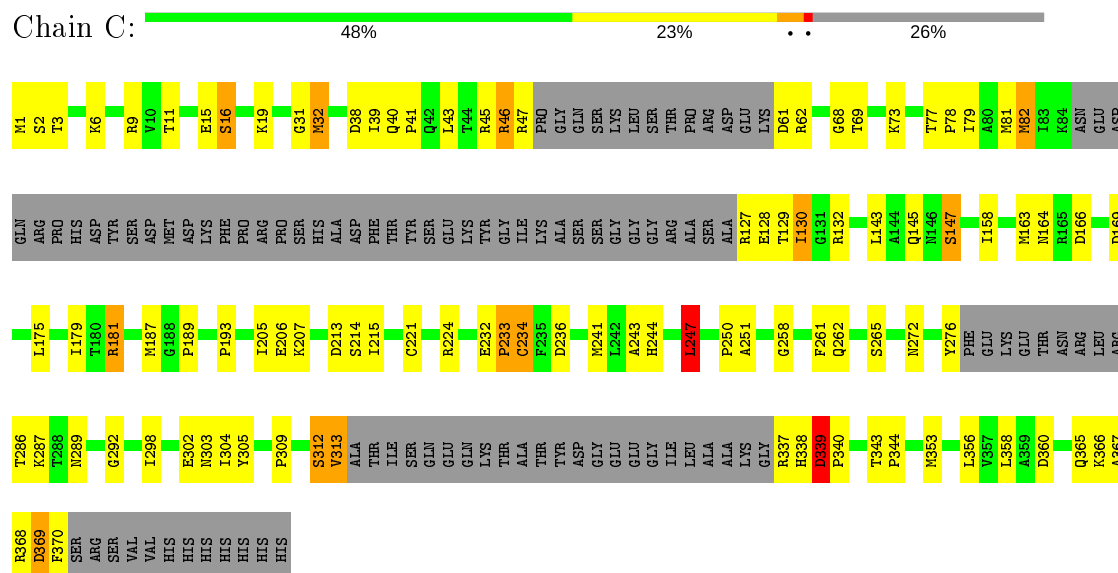
• Molecule 1: Chorismate synthase



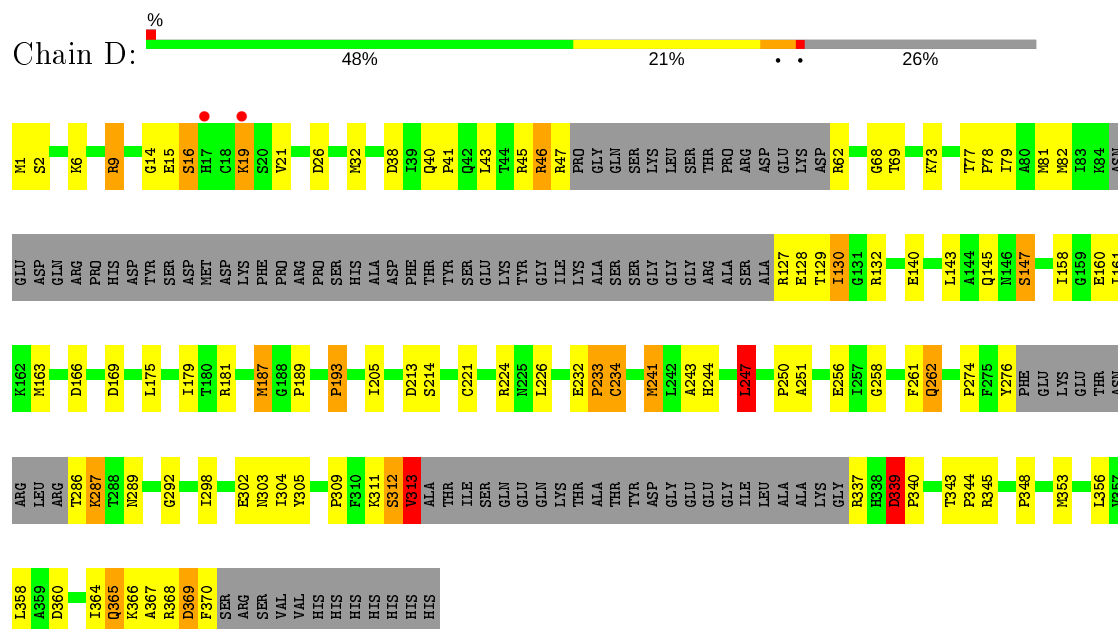
• Molecule 1: Chorismate synthase



• Molecule 1: Chorismate synthase



• Molecule 1: Chorismate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.82Å 75.60Å 91.33Å 114.44° 108.43° 89.98°	Depositor
Resolution (Å)	29.75 – 2.89 29.73 – 2.89	Depositor EDS
% Data completeness (in resolution range)	89.6 (29.75-2.89) 89.6 (29.73-2.89)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.52 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.188 , 0.224 0.176 , 0.211	Depositor DCC
R_{free} test set	1344 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	47.8	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 21.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.467 for h,-k,-h-l 0.467 for -h,k,-k-l 0.467 for -h,-k,h+k+l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8636	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.50% 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

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	1.15	10/2172 (0.5%)	1.12	16/2927 (0.5%)
1	B	1.13	7/2163 (0.3%)	1.14	14/2914 (0.5%)
1	C	1.18	6/2173 (0.3%)	1.41	20/2928 (0.7%)
1	D	1.12	7/2165 (0.3%)	1.15	17/2917 (0.6%)
All	All	1.15	30/8673 (0.3%)	1.21	67/11686 (0.6%)

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	313	VAL	CB-CG1	14.57	1.83	1.52
1	C	61	ASP	CB-CG	-12.56	1.25	1.51
1	C	286	THR	CB-OG1	11.37	1.66	1.43
1	A	276	TYR	CB-CG	-8.92	1.38	1.51
1	A	61	ASP	CB-CG	-8.44	1.34	1.51
1	C	337	ARG	CB-CG	-7.97	1.31	1.52
1	D	286	THR	CB-OG1	7.79	1.58	1.43
1	A	47	ARG	CB-CG	-7.69	1.31	1.52
1	A	19	LYS	CB-CG	-6.85	1.34	1.52
1	B	276	TYR	CB-CG	-6.83	1.41	1.51
1	D	19	LYS	C-O	-6.35	1.11	1.23
1	A	256	GLU	CD-OE1	-6.33	1.18	1.25
1	B	286	THR	CB-OG1	-6.31	1.30	1.43
1	C	19	LYS	C-O	-6.18	1.11	1.23
1	B	47	ARG	CB-CG	-5.84	1.36	1.52
1	A	277	PHE	CB-CG	-5.79	1.41	1.51
1	B	241	MET	CG-SD	5.63	1.95	1.81
1	C	62	ARG	CB-CG	-5.62	1.37	1.52
1	D	241	MET	CG-SD	5.51	1.95	1.81
1	A	187	MET	SD-CE	5.45	2.08	1.77
1	D	187	MET	SD-CE	5.41	2.08	1.77
1	D	286	THR	CB-CG2	-5.41	1.34	1.52
1	B	312	SER	CB-OG	5.39	1.49	1.42
1	D	256	GLU	CD-OE1	-5.36	1.19	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	312	SER	CB-OG	5.27	1.49	1.42
1	A	241	MET	CG-SD	5.26	1.94	1.81
1	B	256	GLU	CD-OE1	-5.17	1.20	1.25
1	A	19	LYS	C-O	-5.17	1.13	1.23
1	B	355	ALA	CA-CB	-5.14	1.41	1.52
1	A	312	SER	CB-OG	5.03	1.48	1.42

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	276	TYR	CB-CG-CD1	-27.23	104.66	121.00
1	C	276	TYR	CB-CG-CD2	23.81	135.28	121.00
1	C	313	VAL	CA-CB-CG2	-18.45	83.22	110.90
1	C	313	VAL	CA-CB-CG1	-18.19	83.62	110.90
1	C	313	VAL	CG1-CB-CG2	-12.15	91.47	110.90
1	B	38	ASP	CB-CG-OD2	10.61	127.84	118.30
1	D	276	TYR	CB-CG-CD2	-10.15	114.91	121.00
1	B	9	ARG	NE-CZ-NH1	9.02	124.81	120.30
1	B	213	ASP	CB-CG-OD2	8.99	126.39	118.30
1	D	38	ASP	CB-CG-OD2	8.80	126.22	118.30
1	C	9	ARG	NE-CZ-NH1	8.70	124.65	120.30
1	C	38	ASP	CB-CG-OD2	8.65	126.08	118.30
1	A	213	ASP	CB-CG-OD2	8.63	126.06	118.30
1	D	213	ASP	CB-CG-OD2	8.52	125.97	118.30
1	B	337	ARG	CA-CB-CG	8.46	132.02	113.40
1	C	213	ASP	CB-CG-OD2	8.41	125.86	118.30
1	D	339	ASP	CB-CG-OD2	8.02	125.52	118.30
1	D	337	ARG	CA-CB-CG	-8.00	95.80	113.40
1	D	47	ARG	NE-CZ-NH2	7.60	124.10	120.30
1	A	9	ARG	NE-CZ-NH1	7.55	124.07	120.30
1	C	47	ARG	NE-CZ-NH2	7.55	124.07	120.30
1	B	47	ARG	NE-CZ-NH2	7.43	124.01	120.30
1	C	337	ARG	NE-CZ-NH2	7.42	124.01	120.30
1	A	62	ARG	NE-CZ-NH2	7.42	124.01	120.30
1	D	127	ARG	NE-CZ-NH2	7.40	124.00	120.30
1	C	127	ARG	NE-CZ-NH2	7.38	123.99	120.30
1	B	62	ARG	NE-CZ-NH2	7.38	123.99	120.30
1	D	337	ARG	NE-CZ-NH2	7.35	123.97	120.30
1	C	62	ARG	NE-CZ-NH2	7.32	123.96	120.30
1	B	337	ARG	NE-CZ-NH2	7.32	123.96	120.30
1	D	62	ARG	NE-CZ-NH2	7.31	123.95	120.30
1	A	277	PHE	CB-CG-CD1	-7.30	115.69	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	337	ARG	NE-CZ-NH2	7.29	123.95	120.30
1	B	127	ARG	NE-CZ-NH2	7.24	123.92	120.30
1	A	47	ARG	NE-CZ-NH2	7.22	123.91	120.30
1	A	38	ASP	CB-CG-OD2	7.15	124.74	118.30
1	D	169	ASP	CB-CG-OD2	6.99	124.59	118.30
1	A	277	PHE	CB-CG-CD2	6.89	125.62	120.80
1	C	236	ASP	CB-CG-OD2	6.88	124.49	118.30
1	C	169	ASP	CB-CG-OD2	6.77	124.39	118.30
1	A	26	ASP	CB-CG-OD1	6.72	124.35	118.30
1	B	247	LEU	CA-CB-CG	-6.58	100.18	115.30
1	D	9	ARG	NE-CZ-NH2	6.50	123.55	120.30
1	D	247	LEU	CA-CB-CG	-6.49	100.38	115.30
1	D	276	TYR	CB-CG-CD1	6.41	124.85	121.00
1	D	26	ASP	CB-CG-OD1	6.38	124.04	118.30
1	C	247	LEU	CA-CB-CG	-6.33	100.75	115.30
1	B	166	ASP	CB-CG-OD2	6.17	123.85	118.30
1	A	181	ARG	NE-CZ-NH2	-6.07	117.27	120.30
1	B	224	ARG	NE-CZ-NH1	-6.04	117.28	120.30
1	A	236	ASP	CB-CG-OD2	5.86	123.57	118.30
1	C	62	ARG	CA-CB-CG	5.84	126.25	113.40
1	A	339	ASP	CB-CG-OD2	5.83	123.55	118.30
1	A	169	ASP	CB-CG-OD2	5.80	123.52	118.30
1	C	339	ASP	CB-CG-OD2	5.80	123.52	118.30
1	C	181	ARG	NE-CZ-NH2	-5.78	117.41	120.30
1	A	247	LEU	CA-CB-CG	-5.71	102.17	115.30
1	D	286	THR	CA-CB-CG2	5.62	120.27	112.40
1	A	369	ASP	CB-CG-OD2	5.62	123.35	118.30
1	B	224	ARG	CG-CD-NE	-5.59	100.07	111.80
1	C	360	ASP	CB-CG-OD2	5.52	123.27	118.30
1	B	169	ASP	CB-CG-OD2	5.39	123.16	118.30
1	C	286	THR	CA-CB-OG1	-5.21	98.05	109.00
1	A	224	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	D	313	VAL	CB-CA-C	-5.16	101.60	111.40
1	D	38	ASP	CB-CG-OD1	-5.08	113.73	118.30
1	B	370	PHE	N-CA-C	5.04	124.60	111.00

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	2138	0	2151	80	0
1	B	2130	0	2151	75	0
1	C	2139	0	2161	76	0
1	D	2131	0	2157	81	0
2	A	25	0	0	2	0
2	B	26	0	0	0	0
2	C	27	0	0	0	0
2	D	20	0	0	2	0
All	All	8636	0	8620	256	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 15.

All (256) 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:C:187:MET:CE	1:C:187:MET:SD	2.05	1.44
1:A:187:MET:SD	1:A:187:MET:CE	2.08	1.41
1:D:187:MET:CE	1:D:187:MET:SD	2.08	1.40
1:B:368:ARG:HD3	2:D:385:HOH:O	1.68	0.92
1:A:367:ALA:HB2	1:C:367:ALA:HB2	1.61	0.81
1:D:9:ARG:HD3	2:D:388:HOH:O	1.80	0.81
1:B:367:ALA:HB2	1:D:367:ALA:HB2	1.62	0.80
1:B:369:ASP:O	1:D:366:LYS:NZ	2.16	0.77
1:B:175:LEU:HD23	1:B:179:ILE:HB	1.68	0.75
1:B:366:LYS:NZ	1:D:369:ASP:O	2.18	0.74
2:A:402:HOH:O	1:C:368:ARG:HD3	1.92	0.70
1:A:233:PRO:HD3	1:A:298:ILE:HG22	1.75	0.69
1:A:262:GLN:HG2	2:A:403:HOH:O	1.92	0.69
1:B:233:PRO:HD3	1:B:298:ILE:HG22	1.74	0.69
1:B:233:PRO:O	1:B:234:CYS:HB2	1.92	0.68
1:D:233:PRO:HD3	1:D:298:ILE:HG22	1.75	0.68
1:A:143:LEU:O	1:A:147:SER:OG	2.12	0.68
1:C:312:SER:O	1:C:313:VAL:CB	2.42	0.67
1:A:64:GLU:CD	1:D:19:LYS:HZ2	1.98	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:233:PRO:O	1:C:234:CYS:HB2	1.94	0.67
1:D:175:LEU:HD23	1:D:179:ILE:HB	1.76	0.66
1:D:353:MET:HE2	1:D:356:LEU:HD12	1.76	0.66
1:C:175:LEU:HD23	1:C:179:ILE:HB	1.78	0.66
1:A:175:LEU:HD23	1:A:179:ILE:HB	1.79	0.65
1:A:365:GLN:HE21	1:A:366:LYS:N	1.95	0.65
1:B:250:PRO:O	1:B:251:ALA:HB3	1.96	0.65
1:C:233:PRO:HD3	1:C:298:ILE:HG22	1.77	0.65
1:A:311:LYS:HE2	1:A:313:VAL:O	1.98	0.64
1:D:365:GLN:HE21	1:D:366:LYS:N	1.96	0.64
1:D:43:LEU:O	1:D:46:ARG:HG2	1.97	0.64
1:C:43:LEU:O	1:C:46:ARG:HG2	1.98	0.64
1:A:166:ASP:HA	1:A:262:GLN:NE2	2.14	0.62
1:D:45:ARG:NH1	1:D:181:ARG:HD3	2.15	0.61
1:A:233:PRO:O	1:A:234:CYS:HB2	2.00	0.61
1:C:312:SER:O	1:C:313:VAL:HB	2.00	0.61
1:C:309:PRO:HD3	1:D:261:PHE:CZ	2.35	0.61
1:D:353:MET:HA	1:D:353:MET:HE2	1.83	0.61
1:D:214:SER:HA	1:D:340:PRO:HB3	1.82	0.61
1:D:166:ASP:HA	1:D:262:GLN:NE2	2.16	0.60
1:C:128:GLU:HG3	1:C:130:ILE:HG12	1.82	0.60
1:D:233:PRO:O	1:D:234:CYS:HB2	2.01	0.60
1:D:166:ASP:HA	1:D:262:GLN:HE22	1.66	0.60
1:A:289:ASN:ND2	1:A:302:GLU:HB3	2.16	0.59
1:A:214:SER:HA	1:A:340:PRO:HB3	1.84	0.59
1:B:360:ASP:O	1:B:364:ILE:HG13	2.02	0.59
1:A:272:ASN:HD21	1:B:313:VAL:HG21	1.67	0.58
1:A:45:ARG:NH1	1:A:181:ARG:HD3	2.18	0.58
1:A:43:LEU:O	1:A:46:ARG:HG2	2.03	0.58
1:C:272:ASN:HD21	1:D:313:VAL:HG21	1.68	0.58
1:A:2:SER:OG	1:D:78:PRO:HG3	2.04	0.57
1:B:214:SER:HA	1:B:340:PRO:HB3	1.85	0.57
1:A:244:HIS:HB2	1:B:244:HIS:HB2	1.85	0.57
1:C:353:MET:HE2	1:C:356:LEU:HD12	1.85	0.57
1:B:311:LYS:HE2	1:B:313:VAL:O	2.04	0.57
1:A:366:LYS:NZ	1:C:369:ASP:O	2.26	0.57
1:A:309:PRO:HD3	1:B:261:PHE:CZ	2.39	0.57
1:A:261:PHE:CZ	1:B:309:PRO:HD3	2.40	0.56
1:B:353:MET:HE2	1:B:356:LEU:HD12	1.88	0.56
1:C:214:SER:HA	1:C:340:PRO:HB3	1.87	0.56
1:A:370:PHE:HA	1:C:366:LYS:NZ	2.20	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:THR:O	1:A:132:ARG:HB2	2.05	0.56
1:A:1:MET:HG2	1:A:3:THR:H	1.71	0.56
1:A:292:GLY:O	1:A:304:ILE:HA	2.05	0.56
1:B:43:LEU:O	1:B:46:ARG:HG2	2.05	0.56
1:B:339:ASP:HB3	1:B:340:PRO:CA	2.36	0.55
1:A:132:ARG:HD3	1:A:353:MET:HE3	1.89	0.55
1:D:163:MET:CE	1:D:193:PRO:HD2	2.36	0.55
1:B:40:GLN:N	1:B:41:PRO:HD2	2.22	0.55
1:D:292:GLY:O	1:D:304:ILE:HA	2.07	0.55
1:B:221:CYS:O	1:B:305:TYR:HA	2.08	0.54
1:A:166:ASP:HA	1:A:262:GLN:HE21	1.73	0.54
1:A:40:GLN:N	1:A:41:PRO:HD2	2.23	0.54
1:A:366:LYS:NZ	1:C:370:PHE:HA	2.22	0.54
1:D:311:LYS:HE2	1:D:313:VAL:O	2.07	0.53
1:D:163:MET:HE2	1:D:193:PRO:HD2	1.90	0.53
1:B:166:ASP:HA	1:B:262:GLN:NE2	2.23	0.53
1:B:313:VAL:HG12	1:B:337:ARG:O	2.09	0.53
1:D:274:PRO:HA	1:D:287:LYS:HB3	1.91	0.53
1:C:292:GLY:O	1:C:304:ILE:HA	2.09	0.53
1:D:128:GLU:HG3	1:D:130:ILE:HG12	1.90	0.53
1:C:6:LYS:HE2	1:D:1:MET:CE	2.39	0.53
1:C:68:GLY:HA3	1:C:79:ILE:HG12	1.90	0.52
1:C:258:GLY:HA2	1:C:305:TYR:CZ	2.44	0.52
1:D:143:LEU:O	1:D:147:SER:OG	2.27	0.52
1:B:163:MET:CE	1:B:193:PRO:HD2	2.40	0.52
1:A:226:LEU:HD11	1:A:358:LEU:HD22	1.92	0.52
1:C:368:ARG:HG3	1:C:368:ARG:HH11	1.74	0.52
1:D:353:MET:CE	1:D:356:LEU:HD12	2.40	0.52
1:A:250:PRO:O	1:A:251:ALA:HB3	2.10	0.51
1:B:132:ARG:HD3	1:B:353:MET:HE3	1.92	0.51
1:B:339:ASP:HB3	1:B:340:PRO:C	2.30	0.51
1:C:258:GLY:HA2	1:C:305:TYR:CE1	2.44	0.51
1:D:250:PRO:O	1:D:251:ALA:HB3	2.11	0.51
1:D:258:GLY:HA2	1:D:305:TYR:CZ	2.45	0.51
1:D:345:ARG:O	1:D:348:PRO:HD2	2.11	0.51
1:A:221:CYS:O	1:A:305:TYR:HA	2.09	0.51
1:C:1:MET:CE	1:D:6:LYS:HE2	2.41	0.51
1:C:163:MET:CE	1:C:193:PRO:HD2	2.40	0.51
1:D:40:GLN:N	1:D:41:PRO:HD2	2.25	0.51
1:B:233:PRO:O	1:B:234:CYS:CB	2.59	0.51
1:C:339:ASP:HB3	1:C:340:PRO:CA	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:14:GLY:CA	1:D:21:VAL:HG12	2.40	0.51
1:A:160:GLU:HG2	1:B:265:SER:HB2	1.92	0.51
1:C:272:ASN:HD21	1:D:313:VAL:CG2	2.24	0.50
1:C:143:LEU:O	1:C:147:SER:OG	2.26	0.50
1:B:163:MET:HE1	1:B:193:PRO:HD2	1.94	0.50
1:A:6:LYS:HE2	1:B:1:MET:CE	2.42	0.50
1:C:244:HIS:HB2	1:D:244:HIS:HB2	1.94	0.50
1:A:128:GLU:HG3	1:A:130:ILE:HG12	1.94	0.49
1:A:15:GLU:O	1:A:16:SER:CB	2.60	0.49
1:A:64:GLU:CD	1:D:19:LYS:NZ	2.65	0.49
1:A:6:LYS:HE2	1:B:1:MET:HE3	1.94	0.49
1:A:339:ASP:HB3	1:A:340:PRO:CA	2.42	0.49
1:B:1:MET:HG2	1:B:3:THR:H	1.77	0.49
1:B:343:THR:HB	1:B:344:PRO:HD3	1.95	0.49
1:C:39:ILE:HG21	1:C:81:MET:HE1	1.94	0.49
1:C:15:GLU:O	1:C:16:SER:CB	2.60	0.49
1:C:45:ARG:NH1	1:C:181:ARG:HD3	2.27	0.49
1:D:233:PRO:HD3	1:D:298:ILE:CG2	2.43	0.49
1:A:343:THR:HB	1:A:344:PRO:HD3	1.95	0.49
1:B:175:LEU:CD2	1:B:179:ILE:HB	2.41	0.49
1:B:370:PHE:CZ	1:D:366:LYS:HD3	2.47	0.49
1:A:78:PRO:HG3	1:D:2:SER:OG	2.12	0.49
1:C:31:GLY:O	1:C:32:MET:C	2.50	0.48
1:C:163:MET:HE2	1:C:193:PRO:HD2	1.96	0.48
1:C:6:LYS:HE2	1:D:1:MET:HE2	1.95	0.48
1:D:15:GLU:O	1:D:16:SER:CB	2.61	0.48
1:B:15:GLU:O	1:B:16:SER:CB	2.61	0.48
1:A:1:MET:CE	1:B:6:LYS:HE2	2.43	0.48
1:C:272:ASN:ND2	1:D:313:VAL:HG21	2.28	0.48
1:D:368:ARG:HG3	1:D:368:ARG:HH11	1.77	0.48
1:A:256:GLU:HG2	1:B:261:PHE:CZ	2.48	0.48
1:B:292:GLY:O	1:B:304:ILE:HA	2.14	0.48
1:B:339:ASP:HB3	1:B:340:PRO:HA	1.96	0.48
1:B:345:ARG:O	1:B:348:PRO:HD2	2.14	0.48
1:D:221:CYS:O	1:D:305:TYR:HA	2.14	0.48
1:D:339:ASP:HB3	1:D:340:PRO:CA	2.44	0.48
1:D:77:THR:HB	1:D:78:PRO:HD2	1.96	0.48
1:C:353:MET:HE2	1:C:353:MET:HA	1.96	0.47
1:C:243:ALA:CB	1:D:247:LEU:HD13	2.45	0.47
1:D:226:LEU:HD11	1:D:358:LEU:HD22	1.95	0.47
1:A:353:MET:HE2	1:A:356:LEU:HD12	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:ASP:O	1:A:364:ILE:HG13	2.15	0.47
1:D:339:ASP:HB3	1:D:340:PRO:HA	1.96	0.47
1:B:129:THR:O	1:B:132:ARG:HB2	2.15	0.47
1:B:258:GLY:HA2	1:B:305:TYR:CZ	2.50	0.47
1:D:129:THR:O	1:D:132:ARG:HB2	2.15	0.47
1:A:274:PRO:O	1:A:286:THR:HA	2.15	0.47
1:D:68:GLY:HA3	1:D:79:ILE:HG12	1.97	0.47
1:A:370:PHE:CZ	1:C:366:LYS:HD3	2.50	0.47
1:B:366:LYS:HD3	1:D:370:PHE:CZ	2.49	0.47
1:B:140:GLU:OE2	1:B:181:ARG:NH2	2.48	0.46
1:C:250:PRO:O	1:C:251:ALA:HB3	2.15	0.46
1:C:343:THR:HB	1:C:344:PRO:HD3	1.96	0.46
1:A:64:GLU:OE1	1:D:19:LYS:NZ	2.48	0.46
1:D:343:THR:HB	1:D:344:PRO:HD3	1.96	0.46
1:D:289:ASN:ND2	1:D:302:GLU:HB3	2.30	0.46
1:B:366:LYS:NZ	1:D:370:PHE:HA	2.31	0.46
1:C:40:GLN:N	1:C:41:PRO:HD2	2.30	0.46
1:A:132:ARG:HD3	1:A:353:MET:CE	2.46	0.46
1:D:15:GLU:O	1:D:16:SER:HB3	2.16	0.46
1:A:365:GLN:C	1:A:365:GLN:HE21	2.20	0.45
1:B:77:THR:HB	1:B:78:PRO:HD2	1.98	0.45
1:A:130:ILE:HG12	1:A:130:ILE:H	1.62	0.45
1:C:339:ASP:HB3	1:C:340:PRO:C	2.36	0.45
1:A:1:MET:N	1:B:232:GLU:HB3	2.32	0.45
1:A:158:ILE:HG21	1:A:205:ILE:HG13	1.98	0.45
1:A:258:GLY:HA2	1:A:305:TYR:CZ	2.52	0.45
1:B:14:GLY:CA	1:B:21:VAL:HG12	2.47	0.45
1:C:1:MET:HG2	1:C:3:THR:H	1.82	0.45
1:A:265:SER:HB2	1:B:160:GLU:HG2	1.99	0.45
1:D:368:ARG:NH1	1:D:368:ARG:HG3	2.32	0.45
1:C:247:LEU:HD13	1:D:243:ALA:HB2	1.99	0.45
1:B:370:PHE:HA	1:D:366:LYS:NZ	2.32	0.45
1:C:289:ASN:ND2	1:C:302:GLU:HB3	2.31	0.45
1:D:360:ASP:O	1:D:364:ILE:HG13	2.16	0.44
1:A:39:ILE:HG21	1:A:81:MET:HE1	1.98	0.44
1:B:226:LEU:HD11	1:B:358:LEU:HD22	1.99	0.44
1:D:81:MET:HE2	1:D:81:MET:HB3	1.80	0.44
1:C:1:MET:CE	1:D:6:LYS:CE	2.96	0.44
1:C:261:PHE:CZ	1:D:309:PRO:HD3	2.52	0.44
1:B:249:ILE:CG2	1:B:250:PRO:HD2	2.47	0.44
1:A:345:ARG:O	1:A:348:PRO:HD2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:MET:HE2	1:A:193:PRO:HD2	1.99	0.44
1:D:289:ASN:ND2	1:D:303:ASN:O	2.50	0.44
1:D:161:ILE:N	1:D:161:ILE:HD13	2.33	0.44
1:D:158:ILE:HG21	1:D:205:ILE:HG13	1.98	0.44
1:A:339:ASP:HB3	1:A:340:PRO:C	2.38	0.43
1:C:289:ASN:ND2	1:C:303:ASN:O	2.51	0.43
1:A:272:ASN:HD21	1:B:313:VAL:CG2	2.29	0.43
1:A:31:GLY:O	1:A:32:MET:C	2.57	0.43
1:C:2:SER:O	1:C:11:THR:HA	2.19	0.43
1:B:81:MET:HE2	1:B:81:MET:HB3	1.86	0.43
1:C:232:GLU:HB3	1:D:1:MET:N	2.33	0.43
1:C:339:ASP:HB3	1:C:340:PRO:HA	2.00	0.43
1:D:14:GLY:HA3	1:D:21:VAL:HG12	2.00	0.43
1:C:233:PRO:O	1:C:234:CYS:CB	2.65	0.43
1:D:258:GLY:HA2	1:D:305:TYR:CE1	2.54	0.43
1:A:1:MET:HE3	1:B:6:LYS:HE2	2.00	0.43
1:C:243:ALA:HB2	1:D:247:LEU:HD13	2.01	0.43
1:C:313:VAL:HB	1:C:338:HIS:O	2.18	0.42
1:A:261:PHE:CE1	1:B:309:PRO:HD3	2.53	0.42
1:B:250:PRO:O	1:B:251:ALA:CB	2.65	0.42
1:B:39:ILE:HG12	1:B:81:MET:HE1	2.00	0.42
1:A:298:ILE:HG21	1:A:298:ILE:HD13	1.67	0.42
1:A:69:THR:HA	1:A:73:LYS:O	2.19	0.42
1:A:77:THR:HB	1:A:78:PRO:HD2	2.01	0.42
1:B:289:ASN:ND2	1:B:302:GLU:HB3	2.34	0.42
1:A:2:SER:O	1:A:11:THR:HA	2.20	0.42
1:A:289:ASN:ND2	1:A:303:ASN:O	2.53	0.42
1:C:265:SER:HB2	1:D:160:GLU:HG2	2.02	0.42
1:A:232:GLU:HB3	1:B:1:MET:N	2.34	0.42
1:B:66:GLN:HE22	1:C:82:MET:CE	2.32	0.42
1:C:206:GLU:O	1:C:207:LYS:C	2.57	0.42
1:C:158:ILE:HG21	1:C:205:ILE:HG13	2.01	0.42
1:A:368:ARG:HH11	1:A:368:ARG:HG3	1.84	0.42
1:B:246:MET:SD	1:B:350:VAL:HG13	2.60	0.42
1:C:163:MET:HG2	1:C:164:ASN:N	2.34	0.42
1:A:366:LYS:HD3	1:C:370:PHE:CE2	2.55	0.42
1:A:368:ARG:HG3	1:A:368:ARG:NH1	2.34	0.42
1:B:266:VAL:HA	1:B:267:PRO:HD3	1.90	0.42
1:B:43:LEU:CD1	1:B:63:VAL:HG21	2.49	0.42
1:A:140:GLU:OE2	1:A:181:ARG:NH2	2.52	0.42
1:C:166:ASP:HA	1:C:262:GLN:NE2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:77:THR:HB	1:C:78:PRO:HD2	2.01	0.42
1:D:175:LEU:CD2	1:D:179:ILE:HB	2.49	0.42
1:C:129:THR:O	1:C:132:ARG:HB2	2.20	0.41
1:C:214:SER:OG	1:C:312:SER:HB2	2.20	0.41
1:A:14:GLY:HA3	1:A:21:VAL:HB	2.03	0.41
1:B:339:ASP:CB	1:B:340:PRO:HA	2.50	0.41
1:C:1:MET:HE3	1:D:6:LYS:HE2	2.02	0.41
1:A:161:ILE:HD13	1:A:161:ILE:N	2.35	0.41
1:B:2:SER:OG	1:C:78:PRO:HG3	2.20	0.41
1:B:31:GLY:O	1:B:32:MET:C	2.58	0.41
1:C:368:ARG:HG3	1:C:368:ARG:NH1	2.34	0.41
1:C:6:LYS:CE	1:D:1:MET:CE	2.99	0.41
1:A:339:ASP:HB3	1:A:340:PRO:HA	2.03	0.41
1:A:15:GLU:O	1:A:16:SER:HB3	2.21	0.41
1:B:221:CYS:HB3	1:B:306:PHE:CZ	2.55	0.41
1:A:243:ALA:HB2	1:B:247:LEU:HD13	2.02	0.41
1:C:221:CYS:O	1:C:305:TYR:HA	2.19	0.41
1:C:272:ASN:HD21	1:D:313:VAL:CB	2.33	0.41
1:D:140:GLU:OE2	1:D:181:ARG:NH2	2.54	0.41
1:D:69:THR:HA	1:D:73:LYS:O	2.20	0.41
1:B:12:THR:HG22	1:B:23:CYS:HB3	2.03	0.41
1:C:1:MET:N	1:D:232:GLU:HB3	2.36	0.41
1:B:143:LEU:O	1:B:147:SER:OG	2.31	0.40
1:C:353:MET:CE	1:C:356:LEU:HD12	2.51	0.40
1:C:358:LEU:HD23	1:C:358:LEU:HA	1.95	0.40
1:B:298:ILE:HG21	1:B:298:ILE:HD13	1.68	0.40
1:A:364:ILE:HG21	1:A:364:ILE:HD13	1.91	0.40
1:A:81:MET:HE2	1:A:81:MET:HB3	1.91	0.40
1:C:69:THR:HA	1:C:73:LYS:O	2.21	0.40
1:B:45:ARG:NH1	1:B:181:ARG:HD3	2.37	0.40
1:B:353:MET:HE2	1:B:353:MET:HA	2.02	0.40
1:B:78:PRO:HG3	1:C:2:SER:OG	2.21	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	273/382 (72%)	249 (91%)	16 (6%)	8 (3%)	4	18
1	B	272/382 (71%)	250 (92%)	14 (5%)	8 (3%)	4	18
1	C	273/382 (72%)	247 (90%)	18 (7%)	8 (3%)	4	18
1	D	272/382 (71%)	248 (91%)	16 (6%)	8 (3%)	4	18
All	All	1090/1528 (71%)	994 (91%)	64 (6%)	32 (3%)	4	18

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	SER
1	A	147	SER
1	A	234	CYS
1	A	339	ASP
1	B	16	SER
1	B	147	SER
1	B	234	CYS
1	B	339	ASP
1	C	16	SER
1	C	147	SER
1	C	234	CYS
1	C	339	ASP
1	D	16	SER
1	D	147	SER
1	D	234	CYS
1	D	339	ASP
1	A	32	MET
1	B	32	MET
1	C	32	MET
1	D	32	MET
1	A	233	PRO
1	B	287	LYS
1	C	233	PRO
1	A	20	SER
1	B	20	SER
1	D	233	PRO
1	D	287	LYS
1	B	233	PRO

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Mol	Chain	Res	Type
1	C	287	LYS
1	D	189	PRO
1	A	189	PRO
1	C	189	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	234/318 (74%)	220 (94%)	14 (6%)	19	49
1	B	233/318 (73%)	219 (94%)	14 (6%)	19	49
1	C	235/318 (74%)	223 (95%)	12 (5%)	24	56
1	D	234/318 (74%)	220 (94%)	14 (6%)	19	49
All	All	936/1272 (74%)	882 (94%)	54 (6%)	20	50

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	ARG
1	A	82	MET
1	A	130	ILE
1	A	145	GLN
1	A	193	PRO
1	A	215	ILE
1	A	224	ARG
1	A	241	MET
1	A	247	LEU
1	A	312	SER
1	A	313	VAL
1	A	339	ASP
1	A	365	GLN
1	A	369	ASP
1	B	46	ARG
1	B	82	MET
1	B	130	ILE

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Mol	Chain	Res	Type
1	B	193	PRO
1	B	241	MET
1	B	247	LEU
1	B	262	GLN
1	B	298	ILE
1	B	311	LYS
1	B	312	SER
1	B	313	VAL
1	B	339	ASP
1	B	365	GLN
1	B	369	ASP
1	C	46	ARG
1	C	82	MET
1	C	130	ILE
1	C	145	GLN
1	C	215	ILE
1	C	224	ARG
1	C	241	MET
1	C	247	LEU
1	C	312	SER
1	C	339	ASP
1	C	365	GLN
1	C	369	ASP
1	D	46	ARG
1	D	82	MET
1	D	130	ILE
1	D	145	GLN
1	D	193	PRO
1	D	224	ARG
1	D	241	MET
1	D	247	LEU
1	D	262	GLN
1	D	312	SER
1	D	313	VAL
1	D	339	ASP
1	D	365	GLN
1	D	369	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	HIS

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Mol	Chain	Res	Type
1	A	66	GLN
1	A	148	ASN
1	A	262	GLN
1	A	272	ASN
1	A	295	GLN
1	A	365	GLN
1	B	17	HIS
1	B	66	GLN
1	B	148	ASN
1	B	262	GLN
1	B	272	ASN
1	B	365	GLN
1	C	66	GLN
1	C	148	ASN
1	C	262	GLN
1	C	272	ASN
1	C	365	GLN
1	D	17	HIS
1	D	66	GLN
1	D	262	GLN
1	D	365	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	283/382 (74%)	-0.42	1 (0%) 92 93	10, 32, 57, 67	9 (3%)
1	B	282/382 (73%)	-0.42	0 100 100	9, 32, 57, 67	9 (3%)
1	C	283/382 (74%)	-0.42	0 100 100	9, 32, 56, 67	10 (3%)
1	D	282/382 (73%)	-0.39	2 (0%) 87 87	10, 32, 57, 67	8 (2%)
All	All	1130/1528 (73%)	-0.41	3 (0%) 94 94	9, 32, 57, 67	36 (3%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	17	HIS	3.2
1	A	17	HIS	3.0
1	D	19	LYS	2.2

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 carbohydrates 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.