



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:10 PM UTC

PDB ID : 6ZNE / pdb_00006zne
Title : MaeB PTA domain R535E mutant
Authors : Lovering, A.L.; Harding, C.J.
Deposited on : 2020-07-06
Resolution : 2.39 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:

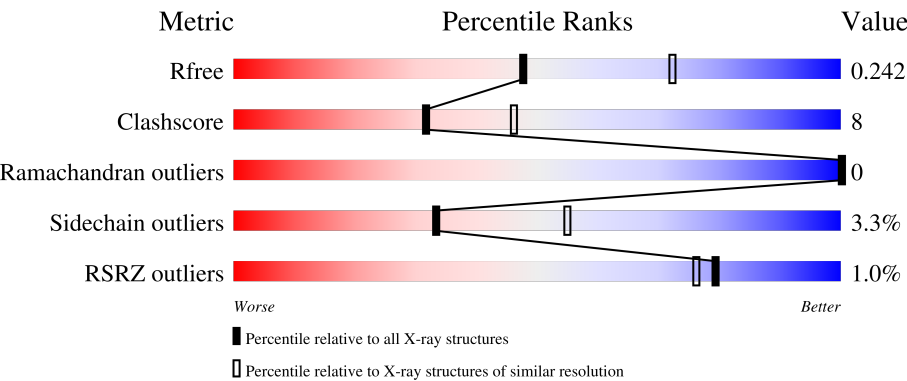
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	362	% 74%18%• 7%
1	B	362	% 74%19%• 7%
1	C	362	2% 79%14%• 6%
1	D	362	2% 69%20%• • 7%
1	E	362	75%17%• 7%

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Mol	Chain	Length	Quality of chain
1	F	362	% 76% 15% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15628 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 Malate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	0	0
			2596	1657	441	487	11			
1	B	338	Total	C	N	O	S	0	0	0
			2596	1657	441	487	11			
1	D	338	Total	C	N	O	S	0	0	0
			2596	1657	441	487	11			
1	F	338	Total	C	N	O	S	0	0	0
			2596	1657	441	487	11			
1	E	338	Total	C	N	O	S	0	0	0
			2596	1657	441	487	11			
1	C	339	Total	C	N	O	S	0	0	0
			2601	1660	442	488	11			

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	419	MET	-	initiating methionine	UNP Q6MM15
A	420	GLY	-	expression tag	UNP Q6MM15
A	421	SER	-	expression tag	UNP Q6MM15
A	422	SER	-	expression tag	UNP Q6MM15
A	423	HIS	-	expression tag	UNP Q6MM15
A	424	HIS	-	expression tag	UNP Q6MM15
A	425	HIS	-	expression tag	UNP Q6MM15
A	426	HIS	-	expression tag	UNP Q6MM15
A	427	HIS	-	expression tag	UNP Q6MM15
A	428	HIS	-	expression tag	UNP Q6MM15
A	429	SER	-	expression tag	UNP Q6MM15
A	430	SER	-	expression tag	UNP Q6MM15
A	431	GLY	-	expression tag	UNP Q6MM15
A	432	LEU	-	expression tag	UNP Q6MM15
A	433	VAL	-	expression tag	UNP Q6MM15
A	434	PRO	-	expression tag	UNP Q6MM15
A	435	ALA	-	expression tag	UNP Q6MM15

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Chain	Residue	Modelled	Actual	Comment	Reference
A	436	GLY	-	expression tag	UNP Q6MM15
A	437	SER	-	expression tag	UNP Q6MM15
A	438	HIS	-	expression tag	UNP Q6MM15
A	535	GLU	ARG	engineered mutation	UNP Q6MM15
B	419	MET	-	initiating methionine	UNP Q6MM15
B	420	GLY	-	expression tag	UNP Q6MM15
B	421	SER	-	expression tag	UNP Q6MM15
B	422	SER	-	expression tag	UNP Q6MM15
B	423	HIS	-	expression tag	UNP Q6MM15
B	424	HIS	-	expression tag	UNP Q6MM15
B	425	HIS	-	expression tag	UNP Q6MM15
B	426	HIS	-	expression tag	UNP Q6MM15
B	427	HIS	-	expression tag	UNP Q6MM15
B	428	HIS	-	expression tag	UNP Q6MM15
B	429	SER	-	expression tag	UNP Q6MM15
B	430	SER	-	expression tag	UNP Q6MM15
B	431	GLY	-	expression tag	UNP Q6MM15
B	432	LEU	-	expression tag	UNP Q6MM15
B	433	VAL	-	expression tag	UNP Q6MM15
B	434	PRO	-	expression tag	UNP Q6MM15
B	435	ALA	-	expression tag	UNP Q6MM15
B	436	GLY	-	expression tag	UNP Q6MM15
B	437	SER	-	expression tag	UNP Q6MM15
B	438	HIS	-	expression tag	UNP Q6MM15
B	535	GLU	ARG	engineered mutation	UNP Q6MM15
D	419	MET	-	initiating methionine	UNP Q6MM15
D	420	GLY	-	expression tag	UNP Q6MM15
D	421	SER	-	expression tag	UNP Q6MM15
D	422	SER	-	expression tag	UNP Q6MM15
D	423	HIS	-	expression tag	UNP Q6MM15
D	424	HIS	-	expression tag	UNP Q6MM15
D	425	HIS	-	expression tag	UNP Q6MM15
D	426	HIS	-	expression tag	UNP Q6MM15
D	427	HIS	-	expression tag	UNP Q6MM15
D	428	HIS	-	expression tag	UNP Q6MM15
D	429	SER	-	expression tag	UNP Q6MM15
D	430	SER	-	expression tag	UNP Q6MM15
D	431	GLY	-	expression tag	UNP Q6MM15
D	432	LEU	-	expression tag	UNP Q6MM15
D	433	VAL	-	expression tag	UNP Q6MM15
D	434	PRO	-	expression tag	UNP Q6MM15
D	435	ALA	-	expression tag	UNP Q6MM15

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Chain	Residue	Modelled	Actual	Comment	Reference
D	436	GLY	-	expression tag	UNP Q6MM15
D	437	SER	-	expression tag	UNP Q6MM15
D	438	HIS	-	expression tag	UNP Q6MM15
D	535	GLU	ARG	engineered mutation	UNP Q6MM15
F	419	MET	-	initiating methionine	UNP Q6MM15
F	420	GLY	-	expression tag	UNP Q6MM15
F	421	SER	-	expression tag	UNP Q6MM15
F	422	SER	-	expression tag	UNP Q6MM15
F	423	HIS	-	expression tag	UNP Q6MM15
F	424	HIS	-	expression tag	UNP Q6MM15
F	425	HIS	-	expression tag	UNP Q6MM15
F	426	HIS	-	expression tag	UNP Q6MM15
F	427	HIS	-	expression tag	UNP Q6MM15
F	428	HIS	-	expression tag	UNP Q6MM15
F	429	SER	-	expression tag	UNP Q6MM15
F	430	SER	-	expression tag	UNP Q6MM15
F	431	GLY	-	expression tag	UNP Q6MM15
F	432	LEU	-	expression tag	UNP Q6MM15
F	433	VAL	-	expression tag	UNP Q6MM15
F	434	PRO	-	expression tag	UNP Q6MM15
F	435	ALA	-	expression tag	UNP Q6MM15
F	436	GLY	-	expression tag	UNP Q6MM15
F	437	SER	-	expression tag	UNP Q6MM15
F	438	HIS	-	expression tag	UNP Q6MM15
F	535	GLU	ARG	engineered mutation	UNP Q6MM15
E	419	MET	-	initiating methionine	UNP Q6MM15
E	420	GLY	-	expression tag	UNP Q6MM15
E	421	SER	-	expression tag	UNP Q6MM15
E	422	SER	-	expression tag	UNP Q6MM15
E	423	HIS	-	expression tag	UNP Q6MM15
E	424	HIS	-	expression tag	UNP Q6MM15
E	425	HIS	-	expression tag	UNP Q6MM15
E	426	HIS	-	expression tag	UNP Q6MM15
E	427	HIS	-	expression tag	UNP Q6MM15
E	428	HIS	-	expression tag	UNP Q6MM15
E	429	SER	-	expression tag	UNP Q6MM15
E	430	SER	-	expression tag	UNP Q6MM15
E	431	GLY	-	expression tag	UNP Q6MM15
E	432	LEU	-	expression tag	UNP Q6MM15
E	433	VAL	-	expression tag	UNP Q6MM15
E	434	PRO	-	expression tag	UNP Q6MM15
E	435	ALA	-	expression tag	UNP Q6MM15

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Chain	Residue	Modelled	Actual	Comment	Reference
E	436	GLY	-	expression tag	UNP Q6MM15
E	437	SER	-	expression tag	UNP Q6MM15
E	438	HIS	-	expression tag	UNP Q6MM15
E	535	GLU	ARG	engineered mutation	UNP Q6MM15
C	419	MET	-	initiating methionine	UNP Q6MM15
C	420	GLY	-	expression tag	UNP Q6MM15
C	421	SER	-	expression tag	UNP Q6MM15
C	422	SER	-	expression tag	UNP Q6MM15
C	423	HIS	-	expression tag	UNP Q6MM15
C	424	HIS	-	expression tag	UNP Q6MM15
C	425	HIS	-	expression tag	UNP Q6MM15
C	426	HIS	-	expression tag	UNP Q6MM15
C	427	HIS	-	expression tag	UNP Q6MM15
C	428	HIS	-	expression tag	UNP Q6MM15
C	429	SER	-	expression tag	UNP Q6MM15
C	430	SER	-	expression tag	UNP Q6MM15
C	431	GLY	-	expression tag	UNP Q6MM15
C	432	LEU	-	expression tag	UNP Q6MM15
C	433	VAL	-	expression tag	UNP Q6MM15
C	434	PRO	-	expression tag	UNP Q6MM15
C	435	ALA	-	expression tag	UNP Q6MM15
C	436	GLY	-	expression tag	UNP Q6MM15
C	437	SER	-	expression tag	UNP Q6MM15
C	438	HIS	-	expression tag	UNP Q6MM15
C	535	GLU	ARG	engineered mutation	UNP Q6MM15

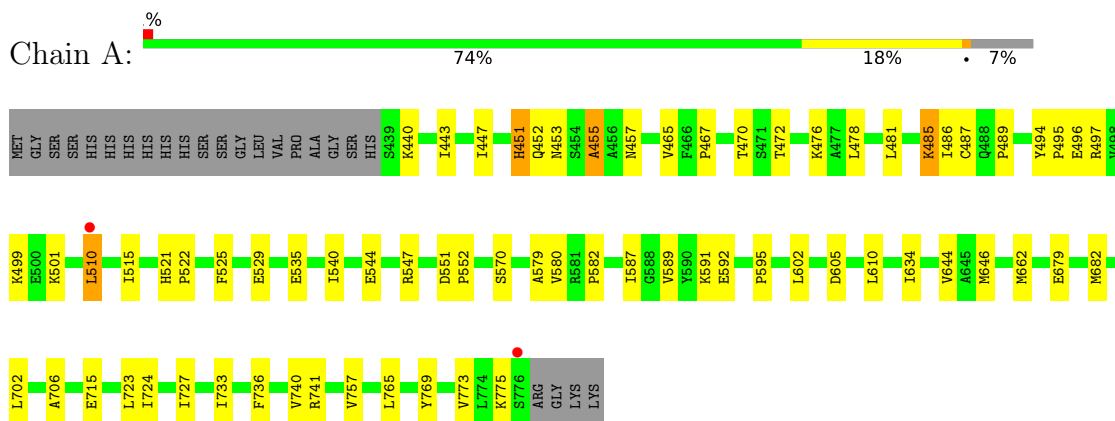
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	10	Total O 10 10	0	0
2	B	9	Total O 9 9	0	0
2	D	10	Total O 10 10	0	0
2	F	2	Total O 2 2	0	0
2	E	9	Total O 9 9	0	0
2	C	7	Total O 7 7	0	0

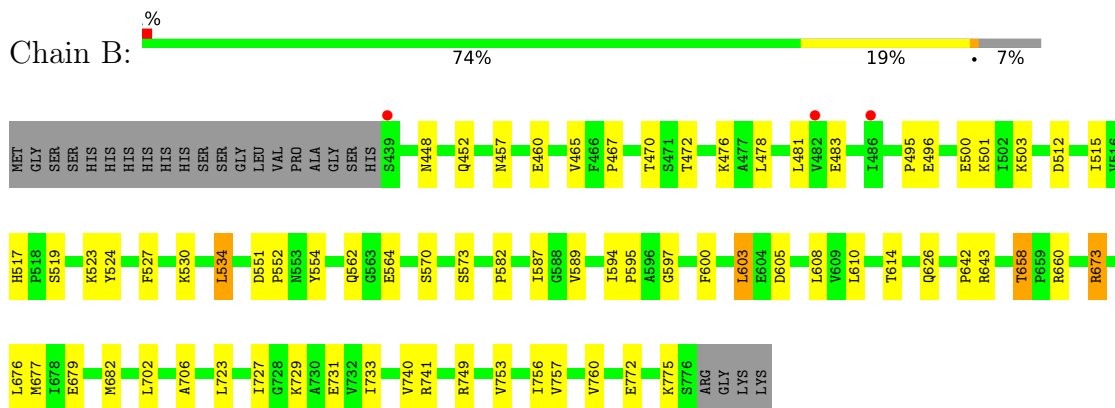
3 Residue-property plots

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: Malate dehydrogenase

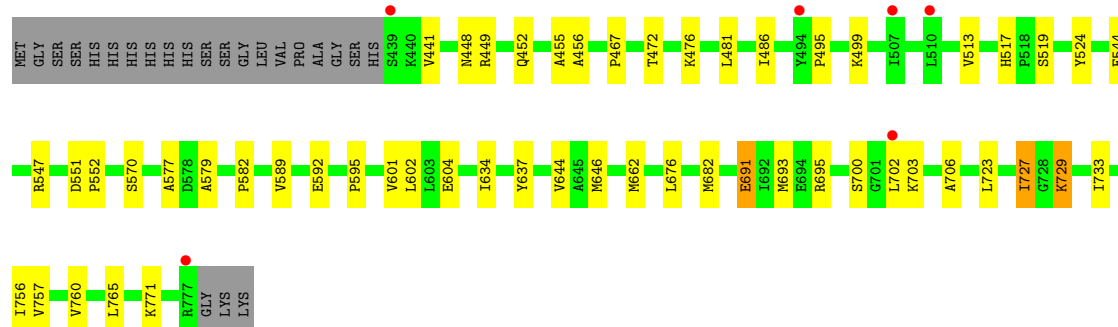


• Molecule 1: Malate dehydrogenase



• Molecule 1: Malate dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	129.42Å 183.40Å 120.24Å 90.00° 117.03° 90.00°	Depositor
Resolution (Å)	69.66 – 2.39 69.66 – 2.39	Depositor EDS
% Data completeness (in resolution range)	98.2 (69.66-2.39) 98.3 (69.66-2.39)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 2.40Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.215 , 0.260 (Not available) , 0.242	Depositor DCC
R_{free} test set	4767 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	56.6	Xtriage
Anisotropy	0.551	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15628	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.87% 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	0.65	2/2640 (0.1%)	1.01	7/3574 (0.2%)
1	B	0.64	0/2640	0.98	6/3574 (0.2%)
1	C	0.61	0/2645	1.04	5/3581 (0.1%)
1	D	0.64	0/2640	1.17	16/3574 (0.4%)
1	E	0.61	0/2640	0.99	8/3574 (0.2%)
1	F	0.61	1/2640 (0.0%)	0.99	5/3574 (0.1%)
All	All	0.63	3/15845 (0.0%)	1.03	47/21451 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	451	HIS	CG-CD2	-8.84	1.26	1.35
1	A	451	HIS	CG-ND1	7.45	1.46	1.38
1	F	451	HIS	CG-CD2	-5.18	1.30	1.35

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	455	ALA	N-CA-C	13.68	139.94	110.80
1	C	456	ALA	N-CA-C	-12.45	94.32	111.87
1	D	507	ILE	CA-C-N	-11.99	104.86	119.84
1	D	507	ILE	C-N-CA	-11.99	104.86	119.84
1	D	730	ALA	N-CA-CB	-9.95	94.67	110.69
1	D	729	LYS	N-CA-CB	9.80	125.76	111.54
1	B	517	HIS	CA-C-N	9.07	128.67	119.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	517	HIS	C-N-CA	9.07	128.67	119.24
1	D	728	GLY	N-CA-C	-8.59	100.31	112.60
1	A	510	LEU	CA-C-N	8.03	131.03	120.28
1	A	510	LEU	C-N-CA	8.03	131.03	120.28
1	D	603	LEU	N-CA-C	-7.88	98.20	110.42
1	D	509	LEU	N-CA-C	-7.06	105.03	112.93
1	A	510	LEU	CB-CA-C	-6.70	99.89	110.94
1	D	510	LEU	N-CA-C	-6.70	104.86	113.16
1	F	517	HIS	CA-C-N	6.62	128.12	119.84
1	F	517	HIS	C-N-CA	6.62	128.12	119.84
1	C	441	VAL	N-CA-C	6.41	116.58	110.42
1	E	517	HIS	CA-C-N	6.40	125.90	119.24
1	E	517	HIS	C-N-CA	6.40	125.90	119.24
1	D	729	LYS	N-CA-C	-6.35	102.80	111.56
1	D	509	LEU	CA-CB-CG	6.07	137.56	116.30
1	B	658	THR	CA-C-N	-5.95	112.78	118.97
1	B	658	THR	C-N-CA	-5.95	112.78	118.97
1	D	512	ASP	N-CA-C	-5.93	104.43	112.26
1	F	680	GLY	N-CA-C	-5.89	103.26	111.09
1	A	455	ALA	N-CA-C	-5.74	102.95	110.53
1	C	644	VAL	N-CA-C	5.67	116.11	108.17
1	E	510	LEU	CA-C-N	-5.66	113.25	122.65
1	E	510	LEU	C-N-CA	-5.66	113.25	122.65
1	D	513	VAL	N-CA-C	-5.65	102.30	109.58
1	D	644	VAL	N-CA-C	5.58	115.89	107.80
1	F	774	LEU	N-CA-C	-5.57	105.29	111.36
1	B	673	ARG	CA-C-N	5.51	126.73	119.84
1	B	673	ARG	C-N-CA	5.51	126.73	119.84
1	D	517	HIS	CA-C-N	5.51	126.72	119.84
1	D	517	HIS	C-N-CA	5.51	126.72	119.84
1	D	498	VAL	N-CA-C	-5.36	105.49	110.53
1	F	447	ILE	N-CA-C	5.25	115.46	110.42
1	E	551	ASP	CA-C-N	5.24	125.29	119.32
1	E	551	ASP	C-N-CA	5.24	125.29	119.32
1	C	513	VAL	N-CA-C	5.20	116.28	109.58
1	A	451	HIS	ND1-CE1-NE2	5.18	113.58	108.40
1	A	451	HIS	CE1-NE2-CD2	-5.15	103.85	109.00
1	A	644	VAL	N-CA-C	5.15	115.38	108.17
1	E	510	LEU	N-CA-C	5.02	119.50	113.17
1	E	644	VAL	N-CA-C	5.00	115.18	108.17

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	482	VAL	Peptide
1	D	510	LEU	Peptide

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	2596	0	2668	41	0
1	B	2596	0	2668	40	0
1	C	2601	0	2670	42	0
1	D	2596	0	2668	69	0
1	E	2596	0	2668	42	0
1	F	2596	0	2666	43	0
2	A	10	0	0	0	0
2	B	9	0	0	1	0
2	C	7	0	0	1	0
2	D	10	0	0	1	0
2	E	9	0	0	1	0
2	F	2	0	0	0	0
All	All	15628	0	16008	257	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 (257) 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:D:729:LYS:HE3	1:C:729:LYS:CB	1.31	1.55
1:D:729:LYS:CE	1:C:729:LYS:HB3	1.50	1.41
1:F:453:ASN:OD1	1:F:768:GLN:NE2	1.58	1.31
1:D:729:LYS:CE	1:C:729:LYS:CB	2.14	1.18
1:D:729:LYS:HE3	1:C:729:LYS:HB2	1.22	1.13
1:D:729:LYS:NZ	1:C:729:LYS:HB3	1.69	1.07
1:F:449:ARG:HH12	1:F:453:ASN:HB3	1.21	1.03
1:D:766:GLU:O	1:D:770:ILE:HG13	1.66	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:449:ARG:HG2	1:E:449:ARG:HH11	1.40	0.86
1:D:729:LYS:HE3	1:C:729:LYS:CA	2.06	0.86
1:D:729:LYS:NZ	1:C:729:LYS:O	2.09	0.84
1:D:509:LEU:O	1:D:511:ASN:ND2	2.14	0.81
1:F:453:ASN:CG	1:F:768:GLN:NE2	2.34	0.81
1:D:591:LYS:HE3	1:D:592:GLU:HG3	1.63	0.81
1:E:676:LEU:O	2:E:801:HOH:O	1.98	0.80
1:A:440:LYS:HA	1:A:443:ILE:HD12	1.67	0.75
1:F:453:ASN:OD1	1:F:768:GLN:CD	2.31	0.74
1:D:511:ASN:OD1	1:D:512:ASP:N	2.21	0.73
1:D:511:ASN:CG	1:D:512:ASP:H	1.93	0.73
1:D:509:LEU:C	1:D:511:ASN:HD22	1.97	0.72
1:C:691:GLU:OE2	1:C:695:ARG:NH2	2.22	0.72
1:D:440:LYS:H	1:D:440:LYS:HD3	1.54	0.71
1:D:482:VAL:HG21	1:D:487:CYS:O	1.90	0.71
1:B:589:VAL:HG11	1:B:595:PRO:HD3	1.72	0.71
1:C:589:VAL:HG11	1:C:595:PRO:HD3	1.73	0.70
1:A:496:GLU:HB2	1:A:497:ARG:NH1	2.06	0.70
1:D:499:LYS:O	1:D:502:ILE:HG22	1.91	0.70
1:D:729:LYS:HZ1	1:C:729:LYS:C	1.99	0.69
1:E:444:ARG:NH2	1:E:484:GLU:OE2	2.25	0.69
1:E:496:GLU:OE1	1:E:496:GLU:N	2.26	0.68
1:F:449:ARG:NH1	1:F:453:ASN:HB3	2.04	0.68
1:B:702:LEU:HD21	1:B:706:ALA:HB2	1.76	0.68
1:B:495:PRO:HA	1:B:515:ILE:HG21	1.76	0.67
1:B:523:LYS:NZ	1:B:564:GLU:OE2	2.29	0.66
1:E:775:LYS:HE2	1:E:776:SER:HB2	1.79	0.64
1:A:589:VAL:HG11	1:A:595:PRO:HD3	1.79	0.64
1:C:481:LEU:HD21	1:C:757:VAL:HG13	1.78	0.64
1:C:577:ALA:N	2:C:801:HOH:O	2.08	0.63
1:C:448:ASN:O	1:C:452:GLN:HG3	1.99	0.63
1:A:715:GLU:OE2	1:B:658:THR:HG21	1.98	0.63
1:A:447:ILE:HG23	1:A:486:ILE:HD11	1.81	0.63
1:D:746:VAL:O	2:D:801:HOH:O	2.16	0.63
1:D:589:VAL:HG11	1:D:595:PRO:HD3	1.80	0.62
1:F:646:MET:HB3	1:F:662:MET:HE3	1.81	0.61
1:E:726:GLN:O	1:E:729:LYS:NZ	2.33	0.61
1:B:496:GLU:O	1:B:500:GLU:HG2	2.01	0.61
1:A:451:HIS:CD2	1:A:485:LYS:HG2	2.35	0.61
1:A:724:ILE:HD13	1:B:727:ILE:HD11	1.82	0.60
1:D:491:LEU:HB2	1:D:515:ILE:HG12	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:724:ILE:CD1	1:B:727:ILE:HD11	2.31	0.60
1:D:702:LEU:HD21	1:D:706:ALA:HB2	1.84	0.60
1:F:470:THR:O	1:F:501:LYS:NZ	2.30	0.60
1:A:646:MET:HB3	1:A:662:MET:HE3	1.83	0.60
1:C:544:GLU:OE1	1:C:547:ARG:NH2	2.31	0.59
1:F:486:ILE:HD11	1:F:760:VAL:HG12	1.85	0.59
1:C:700:SER:O	1:C:703:LYS:NZ	2.28	0.59
1:D:440:LYS:HD3	1:D:440:LYS:N	2.17	0.59
1:E:605:ASP:OD1	1:E:605:ASP:N	2.25	0.58
1:B:723:LEU:O	1:B:727:ILE:HG12	2.03	0.58
1:D:440:LYS:H	1:D:440:LYS:CD	2.09	0.57
1:F:485:LYS:O	1:F:485:LYS:HG2	2.04	0.57
1:E:749:ARG:HG3	1:E:749:ARG:HH11	1.70	0.57
1:D:643:ARG:HB3	1:D:702:LEU:HD11	1.85	0.57
1:A:467:PRO:HG2	1:A:570:SER:HB3	1.88	0.56
1:D:481:LEU:HD13	1:D:760:VAL:HG11	1.87	0.56
1:E:447:ILE:HG23	1:E:486:ILE:HD11	1.87	0.56
1:D:478:LEU:HD23	1:D:481:LEU:HD12	1.87	0.55
1:E:449:ARG:HG2	1:E:449:ARG:NH1	2.18	0.55
1:F:600:PHE:HB2	1:F:733:ILE:HG12	1.87	0.55
1:A:495:PRO:HA	1:A:515:ILE:HG21	1.89	0.55
1:E:448:ASN:O	1:E:452:GLN:HG3	2.05	0.55
1:E:449:ARG:NH1	1:E:637:TYR:OH	2.40	0.55
1:A:481:LEU:HD21	1:A:757:VAL:HG13	1.89	0.55
1:D:509:LEU:HA	1:D:511:ASN:HB3	1.89	0.54
1:E:723:LEU:O	1:E:727:ILE:HG12	2.07	0.54
1:D:509:LEU:O	1:D:509:LEU:HD12	2.07	0.54
1:A:591:LYS:HG2	1:A:592:GLU:N	2.23	0.54
1:A:733:ILE:HG22	1:A:736:PHE:CZ	2.43	0.54
1:F:764:ALA:O	1:F:768:GLN:HG3	2.07	0.53
1:B:448:ASN:O	1:B:452:GLN:HG3	2.08	0.53
1:F:450:VAL:HA	1:F:453:ASN:ND2	2.23	0.53
1:C:723:LEU:O	1:C:727:ILE:HG12	2.09	0.53
1:E:589:VAL:HG11	1:E:595:PRO:HD3	1.90	0.53
1:B:673:ARG:HH11	1:B:676:LEU:HD11	1.74	0.53
1:C:702:LEU:HD11	1:C:706:ALA:HB2	1.91	0.52
1:F:495:PRO:HA	1:F:515:ILE:HG21	1.90	0.52
1:F:531:LEU:HD22	1:F:549:MET:HE2	1.92	0.52
1:F:673:ARG:HG2	1:F:676:LEU:HD12	1.91	0.52
1:E:647:LEU:O	1:E:662:MET:HG3	2.10	0.52
1:B:600:PHE:HB2	1:B:733:ILE:HG12	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:594:ILE:HG22	1:D:626:GLN:HG3	1.90	0.52
1:D:511:ASN:CG	1:D:512:ASP:N	2.58	0.51
1:A:496:GLU:OE1	1:A:496:GLU:N	2.38	0.51
1:E:487:CYS:O	1:E:489:PRO:HD3	2.11	0.51
1:B:587:ILE:O	1:B:741:ARG:HD3	2.10	0.51
1:A:587:ILE:O	1:A:741:ARG:HD3	2.11	0.51
1:B:501:LYS:NZ	2:B:801:HOH:O	2.42	0.50
1:D:741:ARG:HG2	1:D:742:ARG:NH1	2.27	0.50
1:B:600:PHE:HB2	1:B:733:ILE:CG1	2.41	0.50
1:D:479:ALA:HB2	1:D:507:ILE:HG21	1.94	0.50
1:F:485:LYS:O	1:F:485:LYS:CG	2.59	0.50
1:D:478:LEU:HB3	1:D:510:LEU:HD21	1.93	0.50
1:F:600:PHE:HB2	1:F:733:ILE:CG1	2.42	0.50
1:F:473:LYS:HE2	1:F:753:VAL:HG23	1.93	0.50
1:A:723:LEU:O	1:A:727:ILE:HG12	2.11	0.49
1:B:483:GLU:HA	1:B:483:GLU:OE1	2.11	0.49
1:B:589:VAL:HA	1:B:740:VAL:HA	1.92	0.49
1:F:481:LEU:HD21	1:F:757:VAL:HG13	1.94	0.49
1:D:604:GLU:OE1	1:D:604:GLU:N	2.45	0.49
1:A:605:ASP:OD1	1:A:605:ASP:N	2.34	0.49
1:A:496:GLU:HB2	1:A:497:ARG:HH11	1.76	0.49
1:C:604:GLU:H	1:C:604:GLU:CD	2.21	0.49
1:A:455:ALA:C	1:A:457:ASN:H	2.21	0.49
1:F:771:LYS:HE2	1:F:775:LYS:HD3	1.95	0.48
1:C:592:GLU:O	1:C:592:GLU:HG3	2.12	0.48
1:D:600:PHE:HB2	1:D:733:ILE:HG12	1.95	0.48
1:B:753:VAL:O	1:B:757:VAL:HG23	2.13	0.48
1:D:605:ASP:OD2	1:D:606:LYS:NZ	2.44	0.48
1:E:573:SER:HA	1:E:749:ARG:NH1	2.28	0.48
1:C:646:MET:HB3	1:C:662:MET:HE3	1.95	0.48
1:E:702:LEU:HD21	1:E:706:ALA:HB2	1.96	0.48
1:D:729:LYS:HG2	1:C:729:LYS:HB2	1.95	0.48
1:D:630:GLN:NE2	1:D:738:THR:OG1	2.46	0.48
1:D:463:ARG:HD2	1:D:565:ALA:HA	1.95	0.47
1:A:702:LEU:HD11	1:A:706:ALA:HB2	1.95	0.47
1:B:467:PRO:HG2	1:B:570:SER:HB3	1.96	0.47
1:F:495:PRO:O	1:F:499:LYS:HG3	2.15	0.47
1:D:498:VAL:HG21	1:D:515:ILE:HG21	1.96	0.46
1:C:682:MET:SD	1:C:693:MET:HE1	2.56	0.46
1:A:525:PHE:CE2	1:A:529:GLU:OE2	2.68	0.46
1:F:449:ARG:HD2	1:F:449:ARG:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:573:SER:HA	1:B:749:ARG:NH1	2.31	0.46
1:F:587:ILE:O	1:F:741:ARG:HD3	2.15	0.46
1:D:729:LYS:CG	1:C:729:LYS:HB2	2.46	0.46
1:E:766:GLU:O	1:E:770:ILE:HG13	2.16	0.46
1:B:554:TYR:CE2	1:B:582:PRO:HG3	2.50	0.46
1:D:442:PHE:CE2	1:D:762:PHE:HZ	2.34	0.46
1:F:472:THR:CG2	1:F:476:LYS:HE3	2.46	0.46
1:D:500:GLU:HA	1:D:503:LYS:HE3	1.96	0.46
1:F:651:ASN:HB3	1:F:697:PHE:CZ	2.51	0.46
1:B:756:ILE:O	1:B:760:VAL:HG23	2.16	0.45
1:E:544:GLU:OE1	1:E:547:ARG:NH2	2.46	0.45
1:D:441:VAL:HA	1:D:444:ARG:HD2	1.97	0.45
1:F:512:ASP:OD1	1:F:512:ASP:N	2.45	0.45
1:E:749:ARG:HG3	1:E:749:ARG:NH1	2.31	0.45
1:C:693:MET:HE3	1:C:702:LEU:HD22	1.97	0.45
1:B:643:ARG:HG2	1:B:677:MET:HE3	1.98	0.45
1:F:723:LEU:O	1:F:727:ILE:HG12	2.15	0.45
1:B:679:GLU:HB3	1:B:682:MET:HE1	1.98	0.45
1:D:495:PRO:HD3	1:D:517:HIS:HB2	1.97	0.45
1:D:729:LYS:CE	1:C:729:LYS:HB2	2.10	0.45
1:F:551:ASP:HA	1:F:552:PRO:HD3	1.89	0.45
1:D:463:ARG:HG3	1:D:566:ASP:OD2	2.17	0.45
1:E:594:ILE:HG22	1:E:626:GLN:HG3	1.98	0.45
1:A:487:CYS:O	1:A:489:PRO:HD3	2.18	0.44
1:F:447:ILE:O	1:F:450:VAL:HB	2.17	0.44
1:F:728:GLY:O	1:E:728:GLY:HA2	2.17	0.44
1:E:646:MET:HB3	1:E:662:MET:HE3	1.98	0.44
1:B:530:LYS:O	1:B:534:LEU:HD23	2.18	0.44
1:D:524:TYR:O	1:D:528:VAL:HG23	2.17	0.44
1:F:506:ASP:O	1:F:508:PRO:HD3	2.18	0.44
1:B:470:THR:O	1:B:501:LYS:HD3	2.18	0.44
1:E:775:LYS:CE	1:E:776:SER:HB2	2.46	0.44
1:C:495:PRO:HD3	1:C:517:HIS:HB2	1.99	0.44
1:B:512:ASP:OD1	1:B:512:ASP:N	2.47	0.44
1:C:519:SER:HA	1:C:524:TYR:CD1	2.53	0.44
1:F:444:ARG:O	1:F:444:ARG:HG3	2.17	0.44
1:D:487:CYS:O	1:D:489:PRO:HD3	2.18	0.44
1:B:527:PHE:CE1	1:B:562:GLN:HG3	2.53	0.43
1:D:574:ILE:HG22	1:E:547:ARG:NH2	2.33	0.43
1:D:478:LEU:HA	1:D:481:LEU:HD12	1.99	0.43
1:A:485:LYS:O	1:A:485:LYS:CG	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:457:ASN:O	1:D:460:GLU:HG2	2.18	0.43
1:C:602:LEU:HD11	1:C:733:ILE:HD12	2.00	0.43
1:E:673:ARG:HG2	1:E:676:LEU:HD12	2.01	0.43
1:A:472:THR:HG23	1:A:476:LYS:HE3	2.00	0.43
1:A:535:GLU:HB3	1:A:540:ILE:HD12	2.00	0.43
1:E:498:VAL:O	1:E:502:ILE:HG13	2.19	0.43
1:E:551:ASP:HA	1:E:552:PRO:HD3	1.90	0.43
1:E:721:TYR:OH	1:E:735:PRO:HD3	2.19	0.43
1:D:467:PRO:HG2	1:D:570:SER:HB3	2.01	0.43
1:D:673:ARG:HG2	1:D:676:LEU:HD12	2.00	0.43
1:C:756:ILE:O	1:C:760:VAL:HG23	2.18	0.43
1:F:451:HIS:CD2	1:F:485:LYS:HG2	2.54	0.43
1:B:605:ASP:OD1	1:B:605:ASP:N	2.35	0.42
1:F:449:ARG:HH12	1:F:453:ASN:CB	2.11	0.42
1:E:581:ARG:HB3	1:E:582:PRO:HD3	2.01	0.42
1:C:449:ARG:HG2	1:C:765:LEU:HD21	2.02	0.42
1:B:457:ASN:O	1:B:460:GLU:HG2	2.19	0.42
1:A:579:ALA:O	1:A:582:PRO:HD2	2.20	0.42
1:B:478:LEU:HD23	1:B:481:LEU:HD12	2.01	0.42
1:E:478:LEU:HD23	1:E:478:LEU:HA	1.80	0.42
1:D:736:PHE:HE1	1:D:759:SER:HA	1.84	0.42
1:D:766:GLU:O	1:D:770:ILE:CG1	2.53	0.42
1:F:478:LEU:CD1	1:F:491:LEU:HD21	2.50	0.42
1:D:463:ARG:NH1	1:D:563:GLY:O	2.53	0.42
1:D:765:LEU:HD23	1:D:765:LEU:HA	1.88	0.42
1:B:478:LEU:HA	1:B:481:LEU:HD12	2.01	0.42
1:B:642:PRO:HG2	1:B:676:LEU:HD22	2.01	0.42
1:F:481:LEU:HD11	1:F:757:VAL:HA	2.00	0.42
1:E:587:ILE:O	1:E:741:ARG:HD3	2.19	0.42
1:C:495:PRO:O	1:C:499:LYS:HG3	2.20	0.42
1:B:603:LEU:HD11	1:B:608:LEU:HG	2.01	0.42
1:D:729:LYS:HZ1	1:C:729:LYS:HB3	1.52	0.42
1:E:506:ASP:O	1:E:508:PRO:HD3	2.20	0.42
1:A:478:LEU:HA	1:A:478:LEU:HD23	1.79	0.42
1:D:727:ILE:HD12	1:C:601:VAL:HG11	2.01	0.42
1:D:457:ASN:C	1:D:459:GLY:H	2.28	0.41
1:F:756:ILE:O	1:F:760:VAL:HG23	2.20	0.41
1:E:457:ASN:O	1:E:460:GLU:HG2	2.20	0.41
1:C:634:ILE:O	1:C:637:TYR:HB3	2.20	0.41
1:A:521:HIS:CG	1:A:522:PRO:HD2	2.55	0.41
1:A:769:TYR:O	1:A:773:VAL:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:591:LYS:C	1:F:592:GLU:HG2	2.45	0.41
1:F:605:ASP:OD1	1:F:605:ASP:N	2.41	0.41
1:E:481:LEU:HD13	1:E:760:VAL:HG11	2.02	0.41
1:C:676:LEU:HD23	1:C:676:LEU:HA	1.83	0.41
1:D:463:ARG:HG3	1:D:463:ARG:H	1.66	0.41
1:C:472:THR:HG23	1:C:476:LYS:HE3	2.01	0.41
1:A:602:LEU:HD11	1:A:733:ILE:HD11	2.01	0.41
1:A:733:ILE:HG22	1:A:736:PHE:HZ	1.84	0.41
1:A:765:LEU:HD23	1:A:765:LEU:HA	1.88	0.41
1:B:519:SER:HA	1:B:524:TYR:CD1	2.54	0.41
1:D:608:LEU:HD23	1:D:707:ASN:HA	2.02	0.41
1:E:628:ALA:HB2	1:E:646:MET:HE1	2.02	0.41
1:A:495:PRO:O	1:A:499:LYS:HG3	2.20	0.41
1:F:521:HIS:CG	1:F:522:PRO:HD2	2.55	0.41
1:F:594:ILE:HG22	1:F:626:GLN:HG3	2.02	0.41
1:A:453:ASN:O	1:A:457:ASN:ND2	2.28	0.41
1:D:729:LYS:NZ	1:C:729:LYS:C	2.68	0.41
1:D:749:ARG:NH1	1:E:541:ASN:HD22	2.19	0.41
1:E:554:TYR:CE1	1:E:582:PRO:HG3	2.56	0.41
1:C:481:LEU:HD22	1:C:486:ILE:CD1	2.51	0.41
1:A:544:GLU:OE1	1:A:547:ARG:NH2	2.49	0.41
1:A:470:THR:O	1:A:501:LYS:HD3	2.21	0.41
1:B:643:ARG:HB3	1:B:702:LEU:HD11	2.03	0.41
1:E:599:ASN:ND2	1:E:721:TYR:CE1	2.89	0.41
1:C:467:PRO:HG2	1:C:570:SER:HB3	2.03	0.41
1:A:589:VAL:HA	1:A:740:VAL:HA	2.03	0.40
1:B:472:THR:HG23	1:B:476:LYS:HE3	2.03	0.40
1:B:551:ASP:HA	1:B:552:PRO:HD3	1.90	0.40
1:D:636:GLU:OE2	1:D:673:ARG:NH2	2.26	0.40
1:E:444:ARG:HH22	1:E:484:GLU:CD	2.27	0.40
1:B:597:GLY:HA3	1:B:614:THR:OG1	2.22	0.40
1:C:551:ASP:HA	1:C:552:PRO:HD3	1.84	0.40
1:A:494:TYR:HA	1:A:495:PRO:HD3	1.93	0.40
1:A:679:GLU:HB3	1:A:682:MET:HE1	2.02	0.40
1:D:647:LEU:O	1:D:662:MET:HG3	2.21	0.40
1:E:729:LYS:HD2	1:E:729:LYS:HA	1.53	0.40
1:B:594:ILE:HG22	1:B:626:GLN:HG3	2.02	0.40
1:D:679:GLU:HB3	1:D:682:MET:HE1	2.02	0.40
1:C:579:ALA:O	1:C:582:PRO:HD2	2.21	0.40
1:A:551:ASP:HA	1:A:552:PRO:HD3	1.91	0.40
1:D:478:LEU:HD23	1:D:478:LEU:HA	1.88	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:591:LYS:O	1:F:592:GLU:HG2	2.21	0.40
1:F:774:LEU:HD23	1:F:774:LEU:HA	1.78	0.40
1:C:472:THR:CG2	1:C:476:LYS:HE3	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	336/362 (93%)	332 (99%)	4 (1%)	0	100	100
1	B	336/362 (93%)	330 (98%)	6 (2%)	0	100	100
1	C	337/362 (93%)	333 (99%)	4 (1%)	0	100	100
1	D	336/362 (93%)	329 (98%)	7 (2%)	0	100	100
1	E	336/362 (93%)	332 (99%)	4 (1%)	0	100	100
1	F	336/362 (93%)	331 (98%)	5 (2%)	0	100	100
All	All	2017/2172 (93%)	1987 (98%)	30 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	282/301 (94%)	274 (97%)	8 (3%)	38	60
1	B	282/301 (94%)	272 (96%)	10 (4%)	32	53
1	C	282/301 (94%)	278 (99%)	4 (1%)	59	79
1	D	282/301 (94%)	265 (94%)	17 (6%)	17	31
1	E	282/301 (94%)	275 (98%)	7 (2%)	42	64
1	F	282/301 (94%)	273 (97%)	9 (3%)	34	56
All	All	1692/1806 (94%)	1637 (97%)	55 (3%)	33	55

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	452	GLN
1	A	465	VAL
1	A	485	LYS
1	A	510	LEU
1	A	580	VAL
1	A	610	LEU
1	A	634	ILE
1	A	775	LYS
1	B	465	VAL
1	B	503	LYS
1	B	534	LEU
1	B	603	LEU
1	B	610	LEU
1	B	660	ARG
1	B	729	LYS
1	B	731	GLU
1	B	772	GLU
1	B	775	LYS
1	D	440	LYS
1	D	441	VAL
1	D	444	ARG
1	D	465	VAL
1	D	476	LYS
1	D	482	VAL
1	D	499	LYS
1	D	502	ILE
1	D	509	LEU
1	D	538	LYS
1	D	609	VAL
1	D	729	LYS

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Mol	Chain	Res	Type
1	D	731	GLU
1	D	770	ILE
1	D	771	LYS
1	D	772	GLU
1	D	775	LYS
1	F	444	ARG
1	F	529	GLU
1	F	592	GLU
1	F	610	LEU
1	F	721	TYR
1	F	729	LYS
1	F	771	LYS
1	F	773	VAL
1	F	775	LYS
1	E	449	ARG
1	E	465	VAL
1	E	511	ASN
1	E	606	LYS
1	E	609	VAL
1	E	772	GLU
1	E	775	LYS
1	C	691	GLU
1	C	727	ILE
1	C	729	LYS
1	C	771	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	453	ASN
1	A	562	GLN
1	B	562	GLN
1	B	718	ASN
1	F	561	ASN
1	F	562	GLN
1	E	448	ASN
1	E	536	GLN
1	E	562	GLN
1	E	718	ASN
1	C	448	ASN
1	C	585	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

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	338/362 (93%)	-0.04	2 (0%) 85 83	53, 72, 99, 114	0
1	B	338/362 (93%)	-0.10	3 (0%) 81 78	50, 68, 94, 114	0
1	C	339/362 (93%)	-0.03	6 (1%) 67 63	51, 71, 97, 107	0
1	D	338/362 (93%)	0.08	7 (2%) 63 59	55, 74, 107, 129	0
1	E	338/362 (93%)	-0.08	1 (0%) 90 88	57, 71, 98, 116	0
1	F	338/362 (93%)	0.06	2 (0%) 85 83	58, 79, 111, 129	0
All	All	2029/2172 (93%)	-0.02	21 (1%) 79 76	50, 72, 102, 129	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	439	SER	4.4
1	F	481	LEU	3.6
1	D	507	ILE	3.0
1	C	777	ARG	3.0
1	F	439	SER	3.0
1	D	439	SER	2.8
1	C	510	LEU	2.6
1	B	439	SER	2.5
1	E	439	SER	2.5
1	D	481	LEU	2.5
1	D	776	SER	2.4
1	D	509	LEU	2.4
1	D	774	LEU	2.2
1	C	702	LEU	2.2
1	B	486	ILE	2.1
1	B	482	VAL	2.1
1	A	776	SER	2.1
1	C	494	TYR	2.1
1	D	482	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	507	ILE	2.0
1	A	510	LEU	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.