



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:41 PM UTC

PDB ID : 6TFT / pdb_00006tft
Title : Linalool Dehydratase Isomerase C171A mutant
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Deposited on : 2019-11-14
Resolution : 2.52 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

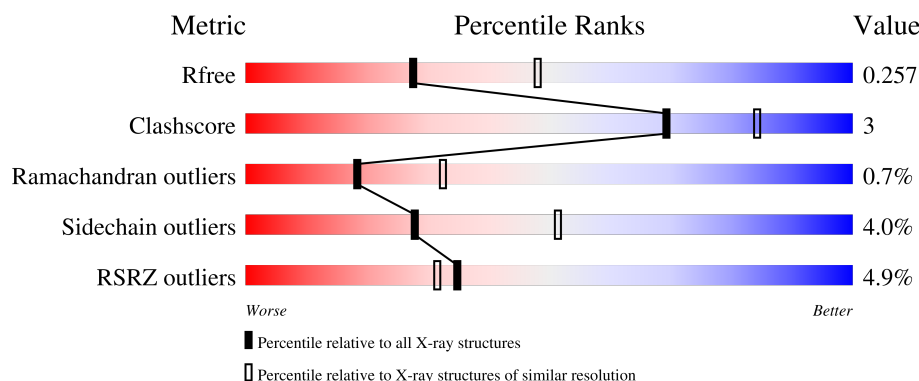
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.52 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	7383 (2.54-2.50)
Clashscore	190562	8079 (2.54-2.50)
Ramachandran outliers	187476	7944 (2.54-2.50)
Sidechain outliers	187428	7946 (2.54-2.50)
RSRZ outliers	180081	7387 (2.54-2.50)

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	372	
1	B	372	
1	C	372	
1	D	372	
1	E	372	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14472 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 Linalool dehydratase-isomerase protein LDI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	365	Total	C	N	O	S	0	1	0
			2908	1877	489	530	12			
1	B	364	Total	C	N	O	S	0	0	0
			2876	1862	477	525	12			
1	C	364	Total	C	N	O	S	0	0	0
			2883	1864	482	525	12			
1	D	363	Total	C	N	O	S	0	1	0
			2844	1844	471	517	12			
1	E	361	Total	C	N	O	S	0	0	0
			2775	1800	454	509	12			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP W8X534
A	171	ALA	CYS	engineered mutation	UNP W8X534
B	1	MET	-	initiating methionine	UNP W8X534
B	171	ALA	CYS	engineered mutation	UNP W8X534
C	1	MET	-	initiating methionine	UNP W8X534
C	171	ALA	CYS	engineered mutation	UNP W8X534
D	1	MET	-	initiating methionine	UNP W8X534
D	171	ALA	CYS	engineered mutation	UNP W8X534
E	1	MET	-	initiating methionine	UNP W8X534
E	171	ALA	CYS	engineered mutation	UNP W8X534

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	59	Total	O	0	0
			59	59		
2	B	47	Total	O	0	0
			47	47		

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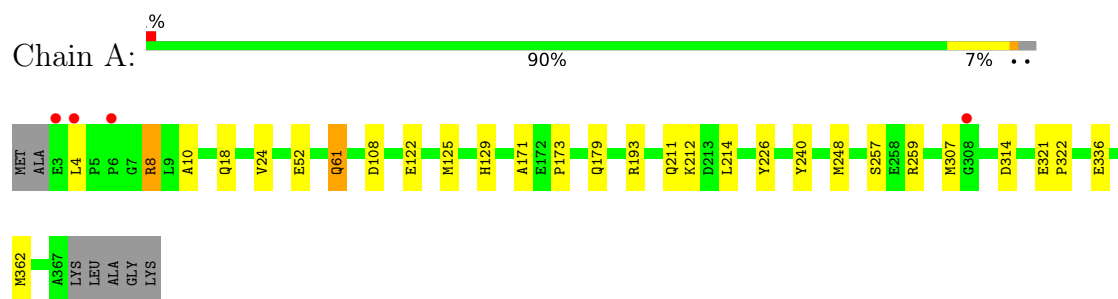
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	28	Total 28	O 28	0	0
2	D	36	Total 36	O 36	0	0
2	E	16	Total 16	O 16	0	0

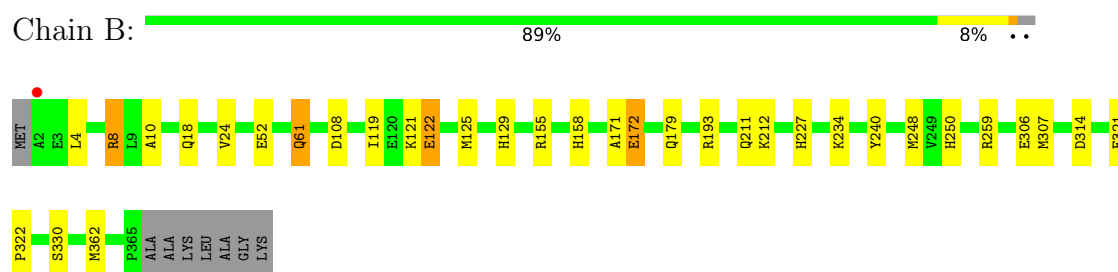
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

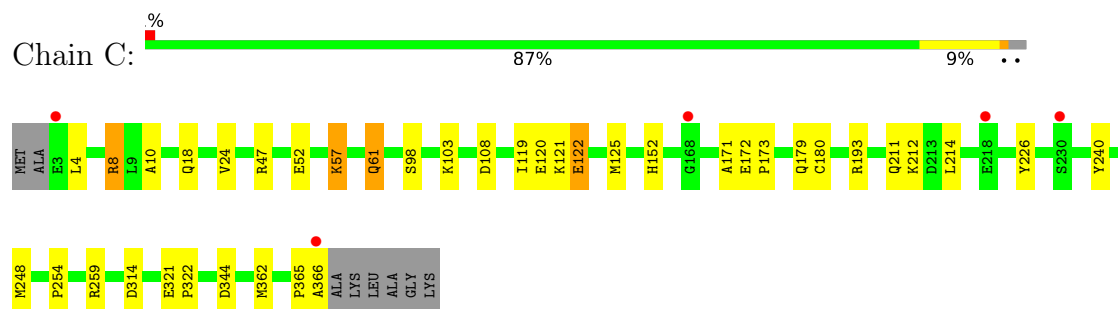
- Molecule 1: Linalool dehydratase-isomerase protein LDI



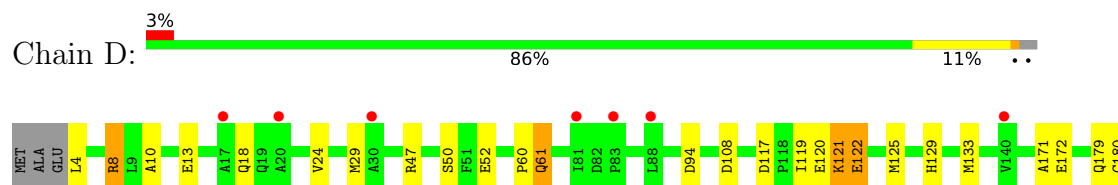
- Molecule 1: Linalool dehydratase-isomerase protein LDI

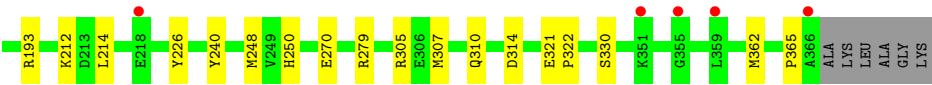


- Molecule 1: Linalool dehydratase-isomerase protein LDI

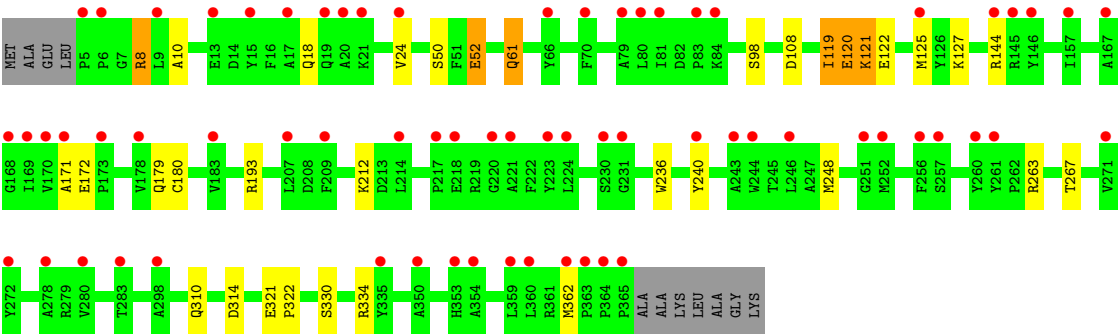
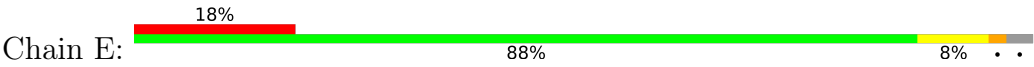


- Molecule 1: Linalool dehydratase-isomerase protein LDI





● Molecule 1: Linalool dehydratase-isomerase protein LDI



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.61Å 111.81Å 120.93Å 90.00° 102.24° 90.00°	Depositor
Resolution (Å)	87.72 – 2.52 87.72 – 2.52	Depositor EDS
% Data completeness (in resolution range)	99.6 (87.72-2.52) 99.6 (87.72-2.52)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.219 , 0.254 0.224 , 0.257	Depositor DCC
R_{free} test set	3898 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	47.9	Xtriage
Anisotropy	0.504	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 33.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14472	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% 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	1.10	0/2997	1.41	1/4079 (0.0%)
1	B	1.15	2/2962 (0.1%)	1.44	4/4034 (0.1%)
1	C	1.10	1/2969 (0.0%)	1.43	2/4043 (0.0%)
1	D	1.13	0/2933	1.43	4/4000 (0.1%)
1	E	1.12	0/2860	1.43	1/3907 (0.0%)
All	All	1.12	3/14721 (0.0%)	1.43	12/20063 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	172	GLU	C-N	14.96	1.51	1.33
1	B	158	HIS	CE1-NE2	7.07	1.39	1.32
1	C	152	HIS	CE1-NE2	6.06	1.38	1.32

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	172	GLU	CA-C-N	12.03	132.80	119.93
1	B	172	GLU	C-N-CA	12.03	132.80	119.93
1	C	179	GLN	CB-CA-C	-5.60	102.34	110.96
1	B	172	GLU	N-CA-C	-5.55	102.78	109.83
1	C	47	ARG	CB-CA-C	-5.43	99.80	110.11
1	B	179	GLN	CB-CA-C	-5.35	102.72	110.96
1	A	179	GLN	CB-CA-C	-5.33	102.75	110.96
1	E	179	GLN	CB-CA-C	-5.31	102.79	110.96
1	D	60	PRO	CA-C-O	-5.16	115.46	121.34
1	D	179	GLN	CB-CA-C	-5.03	102.92	110.92
1	D	119	ILE	CA-C-N	5.02	128.35	120.82
1	D	119	ILE	C-N-CA	5.02	128.35	120.82

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	2908	0	2795	19	0
1	B	2876	0	2753	21	0
1	C	2883	0	2762	17	0
1	D	2844	0	2707	24	0
1	E	2775	0	2583	18	0
2	A	59	0	0	3	0
2	B	47	0	0	0	0
2	C	28	0	0	2	0
2	D	36	0	0	3	0
2	E	16	0	0	4	0
All	All	14472	0	13600	97	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (97) 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:250:HIS:CG	1:B:307:MET:HE1	1.83	1.12
1:B:125:MET:HE2	1:B:129:HIS:CD2	1.98	0.98
1:D:47:ARG:NH2	1:D:94:ASP:OD2	1.97	0.97
1:A:125:MET:HE2	1:A:129:HIS:CD2	2.03	0.92
1:E:119:ILE:O	1:E:127:LYS:HE2	1.69	0.91
1:C:125:MET:HE2	1:C:180:CYS:SG	2.15	0.87
1:E:125:MET:HE2	1:E:180:CYS:SG	2.14	0.87
1:D:47:ARG:HH21	1:D:94:ASP:CG	1.84	0.85
1:D:125:MET:HE2	1:D:180:CYS:SG	2.17	0.84
1:A:211:GLN:HE22	1:A:259[A]:ARG:HH11	1.32	0.75
1:B:250:HIS:ND1	1:B:307:MET:HE1	2.03	0.72
1:D:310:GLN:CB	2:D:431:HOH:O	2.37	0.72
1:A:18:GLN:HE21	1:A:24:VAL:HA	1.56	0.70
1:C:211:GLN:HE22	1:C:259:ARG:HH11	1.37	0.69
1:E:18:GLN:HE21	1:E:24:VAL:HA	1.57	0.69
1:B:18:GLN:HE21	1:B:24:VAL:HA	1.58	0.69
1:B:211:GLN:HE22	1:B:259:ARG:HH11	1.41	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:GLN:HE21	1:C:24:VAL:HA	1.58	0.68
1:B:306:GLU:HG3	1:B:307:MET:HE2	1.75	0.68
1:E:61:GLN:HG3	2:E:401:HOH:O	1.96	0.65
1:B:250:HIS:CG	1:B:307:MET:CE	2.72	0.65
1:A:257:SER:HB3	1:A:307:MET:HE1	1.77	0.65
1:A:173:PRO:O	2:A:401:HOH:O	2.15	0.64
1:D:13:GLU:CB	2:D:433:HOH:O	2.45	0.64
1:D:18:GLN:HE21	1:D:24:VAL:HA	1.61	0.64
1:E:125:MET:CE	1:E:180:CYS:SG	2.86	0.63
1:D:214:LEU:HD21	1:D:226:TYR:HB2	1.81	0.62
1:B:250:HIS:CB	1:B:307:MET:HE1	2.30	0.62
1:C:125:MET:CE	1:C:180:CYS:SG	2.88	0.61
1:B:125:MET:HE2	1:B:129:HIS:NE2	2.16	0.60
1:D:125:MET:CE	1:D:180:CYS:SG	2.90	0.59
1:A:214:LEU:HD21	1:A:226:TYR:HB2	1.83	0.58
1:B:306:GLU:HG3	1:B:307:MET:CE	2.34	0.58
1:B:125:MET:HA	1:B:125:MET:HE3	1.86	0.58
1:E:267:THR:HG23	2:E:404:HOH:O	2.03	0.57
1:C:365:PRO:O	1:C:366:ALA:HB3	2.03	0.57
1:A:125:MET:HE2	1:A:129:HIS:NE2	2.18	0.56
1:B:193:ARG:NH2	1:B:362:MET:O	2.38	0.56
1:E:193:ARG:NH2	1:E:362:MET:O	2.38	0.56
1:A:193:ARG:NH2	1:A:362:MET:O	2.38	0.56
1:A:125:MET:HE3	1:A:125:MET:HA	1.87	0.56
1:B:61:GLN:NE2	1:B:108:ASP:HB3	2.21	0.55
1:D:117:ASP:OD2	1:D:120:GLU:HB2	2.07	0.55
1:C:214:LEU:HD21	1:C:226:TYR:HB2	1.89	0.55
1:A:61:GLN:NE2	1:A:108:ASP:HB3	2.22	0.55
1:D:61:GLN:NE2	1:D:108:ASP:HB3	2.22	0.55
1:C:193:ARG:NH2	1:C:362:MET:O	2.40	0.54
1:D:193:ARG:NH2	1:D:362:MET:O	2.41	0.54
1:E:61:GLN:NE2	1:E:108:ASP:HB3	2.23	0.53
1:C:61:GLN:NE2	1:C:108:ASP:HB3	2.23	0.53
1:C:57:LYS:HE2	1:C:344:ASP:OD2	2.09	0.53
1:C:211:GLN:NE2	1:C:259:ARG:HH11	2.07	0.51
1:E:61:GLN:NE2	2:E:401:HOH:O	2.43	0.51
1:C:321:GLU:HB3	1:C:322:PRO:HD3	1.93	0.50
1:A:321:GLU:HB3	1:A:322:PRO:HD3	1.93	0.50
1:D:29:MET:HE3	1:E:236:TRP:NE1	2.27	0.50
1:D:29:MET:HE3	1:E:236:TRP:CE2	2.46	0.50
1:A:61:GLN:NE2	2:A:406:HOH:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:321:GLU:HB3	1:B:322:PRO:HD3	1.95	0.48
1:D:321:GLU:HB3	1:D:322:PRO:HD3	1.95	0.48
1:A:8:ARG:NH2	1:A:314:ASP:OD1	2.46	0.48
1:C:254:PRO:HD2	2:C:404:HOH:O	2.11	0.48
1:E:8:ARG:NH2	1:E:314:ASP:OD1	2.46	0.48
1:C:173:PRO:O	2:C:401:HOH:O	2.20	0.48
1:D:47:ARG:NH2	1:D:94:ASP:CG	2.55	0.48
1:E:120:GLU:O	1:E:121:LYS:O	2.32	0.48
1:D:8:ARG:NH2	1:D:314:ASP:OD1	2.46	0.47
1:D:305:ARG:HD2	2:D:401:HOH:O	2.14	0.47
1:E:321:GLU:HB3	1:E:322:PRO:HD3	1.96	0.47
1:B:211:GLN:NE2	1:B:259:ARG:HH11	2.09	0.47
1:E:8:ARG:NH1	1:E:10:ALA:O	2.48	0.47
1:D:270:GLU:CB	1:D:279[A]:ARG:NH2	2.79	0.46
1:A:336:GLU:HG2	1:A:336:GLU:O	2.16	0.46
1:C:8:ARG:NH1	1:C:10:ALA:O	2.50	0.45
1:C:8:ARG:NH2	1:C:314:ASP:OD1	2.49	0.45
1:E:310:GLN:HE21	1:E:310:GLN:HB3	1.60	0.45
1:B:8:ARG:NH2	1:B:314:ASP:OD1	2.49	0.45
1:E:61:GLN:CG	2:E:401:HOH:O	2.59	0.44
1:D:8:ARG:NH1	1:D:10:ALA:O	2.51	0.44
1:B:61:GLN:HE22	1:B:108:ASP:HB3	1.82	0.44
1:E:52:GLU:H	1:E:52:GLU:HG2	1.56	0.43
1:A:8:ARG:NH1	1:A:10:ALA:O	2.51	0.43
1:A:257:SER:CB	1:A:307:MET:HE1	2.47	0.43
1:B:8:ARG:NH1	1:B:10:ALA:O	2.52	0.43
1:A:18:GLN:HE21	1:A:24:VAL:CA	2.29	0.42
1:B:227:HIS:CE1	1:B:234:LYS:HE3	2.55	0.42
1:A:61:GLN:HE22	1:A:108:ASP:HB3	1.84	0.42
1:D:250:HIS:ND1	1:D:307:MET:HE3	2.35	0.41
1:D:61:GLN:HE22	1:D:108:ASP:HB3	1.85	0.41
1:D:250:HIS:CE1	1:D:307:MET:HE3	2.56	0.41
1:D:121:LYS:C	1:D:122:GLU:CG	2.93	0.41
1:B:250:HIS:HB2	1:B:307:MET:HE1	2.02	0.41
1:A:321:GLU:OE1	2:A:402:HOH:O	2.22	0.41
1:C:121:LYS:C	1:C:122:GLU:CG	2.94	0.41
1:D:129:HIS:O	1:D:133:MET:HG3	2.22	0.40
1:B:121:LYS:C	1:B:122:GLU:CG	2.94	0.40
1:C:61:GLN:HE22	1:C:108:ASP:HB3	1.85	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	364/372 (98%)	356 (98%)	6 (2%)	2 (0%)	24	41
1	B	362/372 (97%)	354 (98%)	6 (2%)	2 (1%)	21	36
1	C	362/372 (97%)	355 (98%)	5 (1%)	2 (1%)	21	36
1	D	362/372 (97%)	354 (98%)	5 (1%)	3 (1%)	16	29
1	E	359/372 (96%)	350 (98%)	6 (2%)	3 (1%)	16	29
All	All	1809/1860 (97%)	1769 (98%)	28 (2%)	12 (1%)	18	32

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	121	LYS
1	B	171	ALA
1	C	212	LYS
1	D	365	PRO
1	E	212	LYS
1	A	171	ALA
1	A	212	LYS
1	B	212	LYS
1	C	171	ALA
1	D	171	ALA
1	D	212	LYS
1	E	171	ALA

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	294/302 (97%)	287 (98%)	7 (2%)	43	69
1	B	289/302 (96%)	278 (96%)	11 (4%)	29	53
1	C	290/302 (96%)	277 (96%)	13 (4%)	24	46
1	D	283/302 (94%)	272 (96%)	11 (4%)	28	52
1	E	269/302 (89%)	254 (94%)	15 (6%)	19	37
All	All	1425/1510 (94%)	1368 (96%)	57 (4%)	28	51

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	8	ARG
1	A	52	GLU
1	A	61	GLN
1	A	122	GLU
1	A	240	TYR
1	A	248	MET
1	B	4	LEU
1	B	8	ARG
1	B	52	GLU
1	B	61	GLN
1	B	119	ILE
1	B	122	GLU
1	B	155	ARG
1	B	172	GLU
1	B	240	TYR
1	B	248	MET
1	B	330	SER
1	C	4	LEU
1	C	8	ARG
1	C	52	GLU
1	C	57	LYS
1	C	61	GLN
1	C	98	SER
1	C	103	LYS
1	C	119	ILE
1	C	120	GLU
1	C	122	GLU
1	C	172	GLU
1	C	240	TYR
1	C	248	MET

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Mol	Chain	Res	Type
1	D	4	LEU
1	D	8	ARG
1	D	50	SER
1	D	52	GLU
1	D	61	GLN
1	D	121	LYS
1	D	122	GLU
1	D	172	GLU
1	D	240	TYR
1	D	248	MET
1	D	330	SER
1	E	8	ARG
1	E	50	SER
1	E	52	GLU
1	E	61	GLN
1	E	98	SER
1	E	119	ILE
1	E	120	GLU
1	E	122	GLU
1	E	144	ARG
1	E	172	GLU
1	E	240	TYR
1	E	248	MET
1	E	263	ARG
1	E	330	SER
1	E	334	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	61	GLN
1	A	152	HIS
1	A	211	GLN
1	A	227	HIS
1	A	250	HIS
1	A	310	GLN
1	B	18	GLN
1	B	36	ASN
1	B	211	GLN
1	B	227	HIS
1	B	310	GLN

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Mol	Chain	Res	Type
1	C	18	GLN
1	C	36	ASN
1	C	211	GLN
1	C	227	HIS
1	C	250	HIS
1	C	310	GLN
1	D	18	GLN
1	D	36	ASN
1	D	91	HIS
1	D	211	GLN
1	D	227	HIS
1	E	18	GLN
1	E	36	ASN
1	E	61	GLN
1	E	91	HIS
1	E	227	HIS
1	E	310	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 ⓘ

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	365/372 (98%)	-0.09	4 (1%) 78 76	23, 44, 62, 96	1 (0%)
1	B	364/372 (97%)	-0.21	1 (0%) 90 88	31, 41, 56, 78	0
1	C	364/372 (97%)	0.24	5 (1%) 73 71	35, 52, 73, 97	0
1	D	363/372 (97%)	0.68	12 (3%) 49 46	31, 57, 73, 89	1 (0%)
1	E	361/372 (97%)	1.20	67 (18%) 3 2	42, 72, 93, 107	0
All	All	1817/1860 (97%)	0.36	89 (4%) 35 31	23, 52, 82, 107	2 (0%)

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	251	GLY	3.9
1	E	79	ALA	3.7
1	E	365	PRO	3.6
1	C	366	ALA	3.5
1	E	80	LEU	3.3
1	E	363	PRO	3.3
1	E	359	LEU	3.3
1	D	366	ALA	3.3
1	E	256	PHE	3.2
1	D	81	ILE	3.2
1	E	21	LYS	3.2
1	E	125	MET	3.2
1	D	83	PRO	3.1
1	E	231	GLY	3.0
1	E	244	TRP	3.0
1	E	19	GLN	3.0
1	E	173	PRO	3.0
1	E	220	GLY	3.0
1	E	360	LEU	3.0
1	E	144	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	3	GLU	2.8
1	E	81	ILE	2.8
1	E	169	ILE	2.8
1	E	5	PRO	2.8
1	E	217	PRO	2.7
1	E	178	VAL	2.7
1	E	15	TYR	2.7
1	E	207	LEU	2.6
1	E	13	GLU	2.6
1	E	257	SER	2.6
1	E	9	LEU	2.6
1	E	283	THR	2.6
1	E	364	PRO	2.6
1	D	218	GLU	2.6
1	E	224	LEU	2.5
1	E	209	PHE	2.5
1	E	17	ALA	2.5
1	E	278	ALA	2.5
1	E	223	TYR	2.5
1	E	240	TYR	2.5
1	D	351	LYS	2.5
1	E	183	VAL	2.5
1	B	2	ALA	2.5
1	E	243	ALA	2.5
1	E	354	ALA	2.5
1	E	20	ALA	2.5
1	E	167	ALA	2.5
1	E	335	TYR	2.4
1	E	171	ALA	2.4
1	A	308	GLY	2.4
1	E	260	TYR	2.4
1	E	24	VAL	2.4
1	E	170	VAL	2.4
1	E	246	LEU	2.4
1	D	20	ALA	2.4
1	E	218	GLU	2.4
1	E	298	ALA	2.4
1	E	146	TYR	2.3
1	E	261	TYR	2.3
1	C	230	SER	2.3
1	E	272	TYR	2.3
1	A	6	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	252	MET	2.3
1	C	218	GLU	2.3
1	D	30	ALA	2.3
1	E	6	PRO	2.3
1	E	230	SER	2.3
1	A	4	LEU	2.3
1	D	17	ALA	2.2
1	E	157	ILE	2.2
1	E	83	PRO	2.2
1	D	355	GLY	2.2
1	E	271	VAL	2.2
1	E	353	HIS	2.2
1	E	66	TYR	2.1
1	A	3	GLU	2.1
1	D	88	LEU	2.1
1	D	359	LEU	2.1
1	E	70	PHE	2.1
1	E	221	ALA	2.1
1	E	362	MET	2.1
1	D	140	VAL	2.1
1	E	350	ALA	2.1
1	E	145	ARG	2.1
1	E	168	GLY	2.0
1	E	280	VAL	2.0
1	E	214	LEU	2.0
1	C	168	GLY	2.0
1	E	84	LYS	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.