



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2020 – 06:04 PM BST

PDB ID : 6TFT
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 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 : 2.14.6
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) : 2.14.6

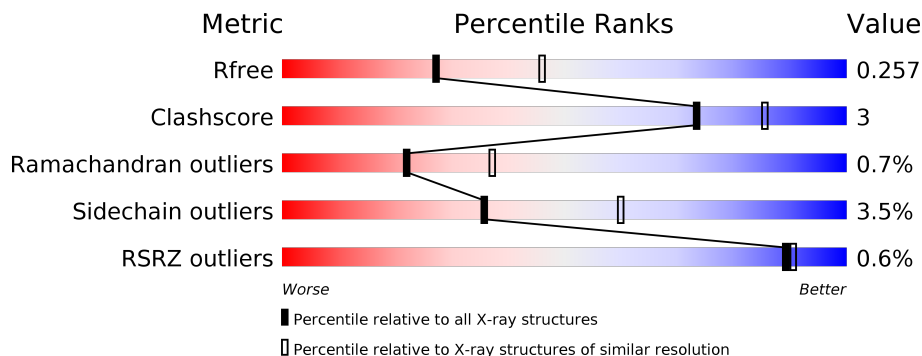
1 Overall quality at a glance

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}	130704	5743 (2.54-2.50)
Clashscore	141614	6463 (2.54-2.50)
Ramachandran outliers	138981	6335 (2.54-2.50)
Sidechain outliers	138945	6337 (2.54-2.50)
RSRZ outliers	127900	5630 (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 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	372	
1	B	372	
1	C	372	
1	D	372	
1	E	372	

2 Entry composition [i](#)

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

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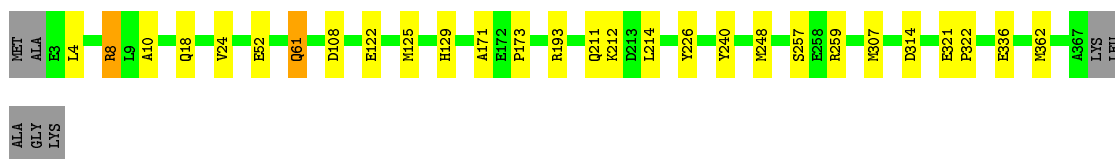
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	28	Total 28	O 28	0	0
2	D	36	Total 36	O 36	0	0
2	E	16	Total 16	O 16	0	0

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.

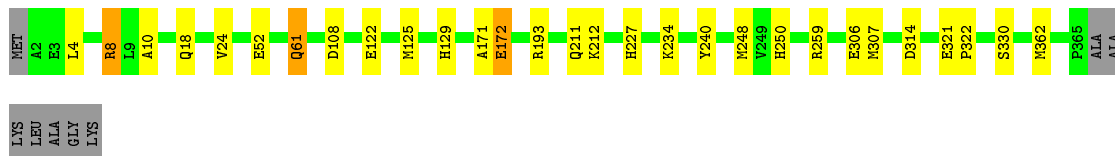
- Molecule 1: Linalool dehydratase-isomerase protein LDI

Chain A: 




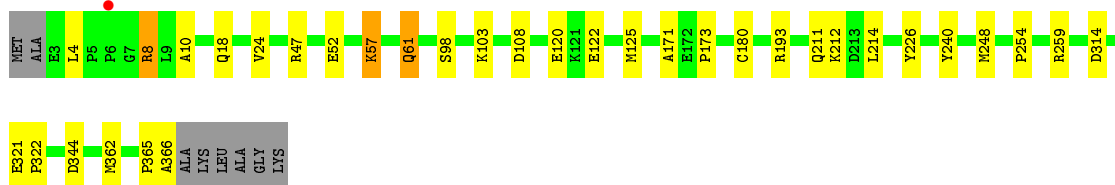
- Molecule 1: Linalool dehydratase-isomerase protein LDI

Chain B: 




- Molecule 1: Linalool dehydratase-isomerase protein LDI

Chain C: 



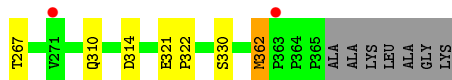
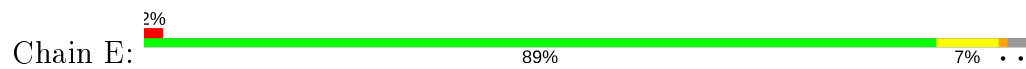
- Molecule 1: Linalool dehydratase-isomerase protein LDI

Chain D: 





- 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 87.57 – 2.52	Depositor EDS
% Data completeness (in resolution range)	99.6 (87.72-2.52) 99.6 (87.57-2.52)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 2.51Å)	Xtrriage
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 (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	47.9	Xtrriage
Anisotropy	0.504	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
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

Xtrriage'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.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.74	0/2997	0.81	0/4079
1	B	0.76	1/2962 (0.0%)	0.81	0/4034
1	C	0.74	0/2969	0.81	1/4043 (0.0%)
1	D	0.76	0/2933	0.82	1/4000 (0.0%)
1	E	0.75	0/2860	0.80	0/3907
All	All	0.75	1/14721 (0.0%)	0.81	2/20063 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	172	GLU	C-N	8.96	1.51	1.34

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	8	ARG	NE-CZ-NH2	-5.64	117.48	120.30
1	C	47	ARG	CB-CA-C	-5.30	99.80	110.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2908	0	2795	18	0
1	B	2876	0	2753	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2883	0	2762	17	0
1	D	2844	0	2707	19	0
1	E	2775	0	2583	17	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	93	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 (93) 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:47:ARG:NH2	1:D:94:ASP:OD2	1.97	0.97
1:E:119:ILE:O	1:E:127:LYS:HE2	1.69	0.91
1:D:47:ARG:HH21	1:D:94:ASP:CG	1.84	0.81
1:A:211:GLN:HE22	1:A:259[A]:ARG:HH11	1.32	0.75
1:B:250:HIS:CG	1:B:307:MET:CE	2.72	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:A:125:MET:HE2	1:A:129:HIS:CD2	2.28	0.68
1:B:211:GLN:HE22	1:B:259:ARG:HH11	1.41	0.68
1:C:18:GLN:HE21	1:C:24:VAL:HA	1.58	0.68
1:B:250:HIS:CG	1:B:307:MET:HE1	2.29	0.68
1:B:250:HIS:CG	1:B:307:MET:HE2	2.31	0.66
1:E:61:GLN:HG3	2:E:401:HOH:O	1.96	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:B:125:MET:HE2	1:B:129:HIS:CD2	2.33	0.63
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:C:125:MET:CE	1:C:180:CYS:SG	2.88	0.61
1:D:125:MET:CE	1:D:180:CYS:SG	2.90	0.59
1:B:250:HIS:CD2	1:B:307:MET:HE2	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:250:HIS:HB2	1:B:307:MET:HE1	1.84	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: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:D:125:MET:HE1	1:D:180:CYS:SG	2.46	0.56
1:A:193:ARG:NH2	1:A:362:MET:O	2.38	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:C:125:MET:HE1	1:C:180:CYS:SG	2.47	0.53
1:B:250:HIS:CB	1:B:307:MET:HE1	2.38	0.53
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:E:125:MET:HE1	1:E:180:CYS:SG	2.49	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:B:125:MET:CE	1:B:129:HIS:CD2	2.93	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:B:125:MET:HE1	1:B:129:HIS:NE2	2.27	0.49
1:A:61:GLN:NE2	2:A:406:HOH:O	2.45	0.49
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: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:A:257:SER:HB3	1:A:307:MET:HE1	1.96	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:E:125:MET:HE2	1:E:180:CYS:SG	2.54	0.46
1:C:8:ARG:NH1	1:C:10:ALA:O	2.50	0.45
1:B:125:MET:O	1:B:125:MET:HE2	2.16	0.45
1:C:8:ARG:NH2	1:C:314:ASP:OD1	2.49	0.45
1:B:306:GLU:HG3	1:B:307:MET:HE3	1.99	0.45
1:E:310:GLN:HE21	1:E:310:GLN:HB3	1.60	0.45
1:A:125:MET:CE	1:A:129:HIS:CD2	2.98	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:125:MET:O	1:A:125:MET:HE2	2.19	0.43
1:B:8:ARG:NH1	1:B:10:ALA:O	2.52	0.43
1:D:47:ARG:NH2	1:D:94:ASP:CG	2.55	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:61:GLN:HE22	1:D:108:ASP:HB3	1.85	0.41
1:A:321:GLU:OE1	2:A:402:HOH:O	2.22	0.41
1:C:125:MET:HE2	1:C:180:CYS:SG	2.60	0.41
1:D:129:HIS:O	1:D:133:MET:HG3	2.22	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

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	364/372 (98%)	356 (98%)	6 (2%)	2 (0%)	29	47
1	B	362/372 (97%)	354 (98%)	6 (2%)	2 (1%)	25	41
1	C	362/372 (97%)	355 (98%)	5 (1%)	2 (1%)	25	41
1	D	362/372 (97%)	354 (98%)	5 (1%)	3 (1%)	19	33
1	E	359/372 (96%)	350 (98%)	6 (2%)	3 (1%)	19	33
All	All	1809/1860 (97%)	1769 (98%)	28 (2%)	12 (1%)	22	37

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%)	49	73
1	B	289/302 (96%)	280 (97%)	9 (3%)	40	65
1	C	290/302 (96%)	279 (96%)	11 (4%)	33	56
1	D	283/302 (94%)	273 (96%)	10 (4%)	36	60
1	E	269/302 (89%)	256 (95%)	13 (5%)	25	45
All	All	1425/1510 (94%)	1375 (96%)	50 (4%)	36	60

All (50) 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	122	GLU
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	120	GLU
1	C	122	GLU
1	C	240	TYR
1	C	248	MET
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	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

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Mol	Chain	Res	Type
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	362	MET

Some 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	211	GLN
1	A	227	HIS
1	A	310	GLN
1	B	18	GLN
1	B	61	GLN
1	B	211	GLN
1	B	227	HIS
1	B	310	GLN
1	C	18	GLN
1	C	61	GLN
1	C	211	GLN
1	C	227	HIS
1	C	250	HIS
1	C	310	GLN
1	D	18	GLN
1	D	61	GLN
1	D	195	HIS
1	D	211	GLN
1	D	227	HIS
1	D	250	HIS
1	E	18	GLN
1	E	61	GLN
1	E	91	HIS
1	E	195	HIS
1	E	227	HIS
1	E	250	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 monosaccharides 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	365/372 (98%)	-0.37	0 100 100	33, 44, 62, 96	0
1	B	364/372 (97%)	-0.42	0 100 100	31, 41, 56, 78	0
1	C	364/372 (97%)	-0.33	1 (0%) 94 94	35, 52, 73, 97	0
1	D	363/372 (97%)	-0.09	2 (0%) 89 90	31, 57, 73, 89	0
1	E	361/372 (97%)	0.11	8 (2%) 62 65	42, 72, 93, 107	0
All	All	1817/1860 (97%)	-0.22	11 (0%) 89 90	31, 52, 82, 107	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	257	SER	2.6
1	E	7	GLY	2.5
1	E	20	ALA	2.5
1	D	4	LEU	2.4
1	E	79	ALA	2.4
1	E	363	PRO	2.3
1	E	271	VAL	2.2
1	E	256	PHE	2.2
1	D	81	ILE	2.2
1	C	6	PRO	2.1
1	E	246	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 monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.