



# wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 22, 2023 – 05:21 PM JST

PDB ID : 7YMP  
Title : Crystal structure of lysoplasmalogen specific phospholipase D  
Authors : Murayama, K.; Kato-Murayama, M.; Sugimori, D.; Shirouzu, M.; Hamana, H.  
Deposited on : 2022-07-29  
Resolution : 2.57 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

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<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>.

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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 : 2.36  
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) : 2.36

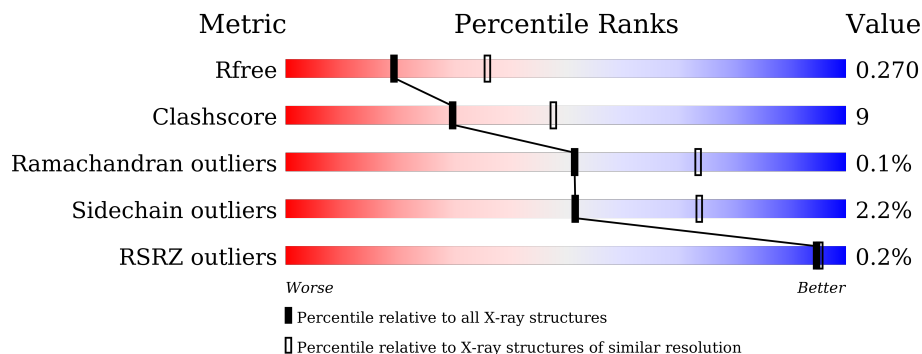
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.57 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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  $\geq 3$ , 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	314	
1	B	314	
1	C	314	
1	D	314	
1	E	314	
1	F	314	

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Mol	Chain	Length	Quality of chain	
1	G	314	 71%	25% ..
1	H	314	 72%	23% ..

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 18857 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 Lysoplasmalogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	306	Total 2356	C 1498	N 411	O 440	S 7	0	0	0
1	B	301	Total 2317	C 1476	N 403	O 431	S 7	0	0	0
1	C	301	Total 2317	C 1476	N 403	O 431	S 7	0	0	0
1	D	301	Total 2317	C 1476	N 403	O 431	S 7	0	0	0
1	E	301	Total 2317	C 1476	N 403	O 431	S 7	0	0	0
1	F	301	Total 2317	C 1476	N 403	O 431	S 7	0	0	0
1	G	301	Total 2317	C 1476	N 403	O 431	S 7	0	0	0
1	H	301	Total 2317	C 1476	N 403	O 431	S 7	0	0	0

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	27	MET	-	initiating methionine	UNP A0A0U4VTN7
A	335	HIS	-	expression tag	UNP A0A0U4VTN7
A	336	HIS	-	expression tag	UNP A0A0U4VTN7
A	337	HIS	-	expression tag	UNP A0A0U4VTN7
A	338	HIS	-	expression tag	UNP A0A0U4VTN7
A	339	HIS	-	expression tag	UNP A0A0U4VTN7
A	340	HIS	-	expression tag	UNP A0A0U4VTN7
B	27	MET	-	initiating methionine	UNP A0A0U4VTN7
B	335	HIS	-	expression tag	UNP A0A0U4VTN7
B	336	HIS	-	expression tag	UNP A0A0U4VTN7
B	337	HIS	-	expression tag	UNP A0A0U4VTN7
B	338	HIS	-	expression tag	UNP A0A0U4VTN7
B	339	HIS	-	expression tag	UNP A0A0U4VTN7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	340	HIS	-	expression tag	UNP A0A0U4VTN7
C	27	MET	-	initiating methionine	UNP A0A0U4VTN7
C	335	HIS	-	expression tag	UNP A0A0U4VTN7
C	336	HIS	-	expression tag	UNP A0A0U4VTN7
C	337	HIS	-	expression tag	UNP A0A0U4VTN7
C	338	HIS	-	expression tag	UNP A0A0U4VTN7
C	339	HIS	-	expression tag	UNP A0A0U4VTN7
C	340	HIS	-	expression tag	UNP A0A0U4VTN7
D	27	MET	-	initiating methionine	UNP A0A0U4VTN7
D	335	HIS	-	expression tag	UNP A0A0U4VTN7
D	336	HIS	-	expression tag	UNP A0A0U4VTN7
D	337	HIS	-	expression tag	UNP A0A0U4VTN7
D	338	HIS	-	expression tag	UNP A0A0U4VTN7
D	339	HIS	-	expression tag	UNP A0A0U4VTN7
D	340	HIS	-	expression tag	UNP A0A0U4VTN7
E	27	MET	-	initiating methionine	UNP A0A0U4VTN7
E	335	HIS	-	expression tag	UNP A0A0U4VTN7
E	336	HIS	-	expression tag	UNP A0A0U4VTN7
E	337	HIS	-	expression tag	UNP A0A0U4VTN7
E	338	HIS	-	expression tag	UNP A0A0U4VTN7
E	339	HIS	-	expression tag	UNP A0A0U4VTN7
E	340	HIS	-	expression tag	UNP A0A0U4VTN7
F	27	MET	-	initiating methionine	UNP A0A0U4VTN7
F	335	HIS	-	expression tag	UNP A0A0U4VTN7
F	336	HIS	-	expression tag	UNP A0A0U4VTN7
F	337	HIS	-	expression tag	UNP A0A0U4VTN7
F	338	HIS	-	expression tag	UNP A0A0U4VTN7
F	339	HIS	-	expression tag	UNP A0A0U4VTN7
F	340	HIS	-	expression tag	UNP A0A0U4VTN7
G	27	MET	-	initiating methionine	UNP A0A0U4VTN7
G	335	HIS	-	expression tag	UNP A0A0U4VTN7
G	336	HIS	-	expression tag	UNP A0A0U4VTN7
G	337	HIS	-	expression tag	UNP A0A0U4VTN7
G	338	HIS	-	expression tag	UNP A0A0U4VTN7
G	339	HIS	-	expression tag	UNP A0A0U4VTN7
G	340	HIS	-	expression tag	UNP A0A0U4VTN7
H	27	MET	-	initiating methionine	UNP A0A0U4VTN7
H	335	HIS	-	expression tag	UNP A0A0U4VTN7
H	336	HIS	-	expression tag	UNP A0A0U4VTN7
H	337	HIS	-	expression tag	UNP A0A0U4VTN7
H	338	HIS	-	expression tag	UNP A0A0U4VTN7
H	339	HIS	-	expression tag	UNP A0A0U4VTN7

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Chain	Residue	Modelled	Actual	Comment	Reference
H	340	HIS	-	expression tag	UNP A0A0U4VTN7

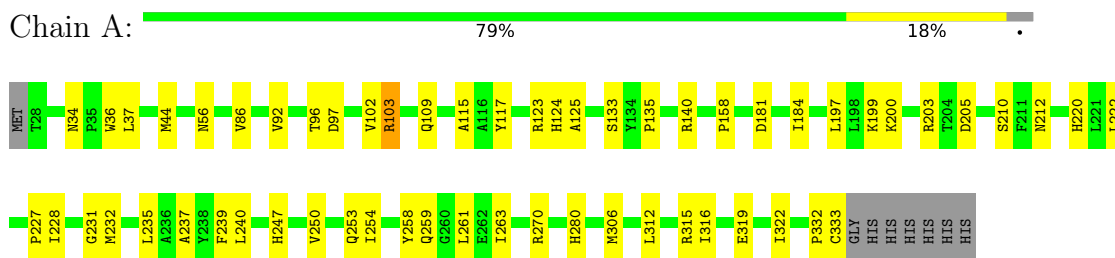
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	35	Total O 35 35	0	0
2	B	45	Total O 45 45	0	0
2	C	40	Total O 40 40	0	0
2	D	29	Total O 29 29	0	0
2	E	36	Total O 36 36	0	0
2	F	30	Total O 30 30	0	0
2	G	42	Total O 42 42	0	0
2	H	25	Total O 25 25	0	0

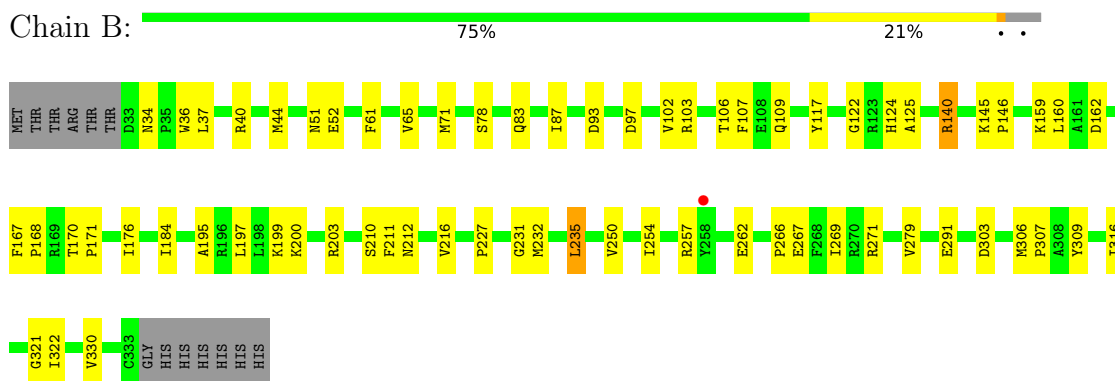
### 3 Residue-property plots [i](#)

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.

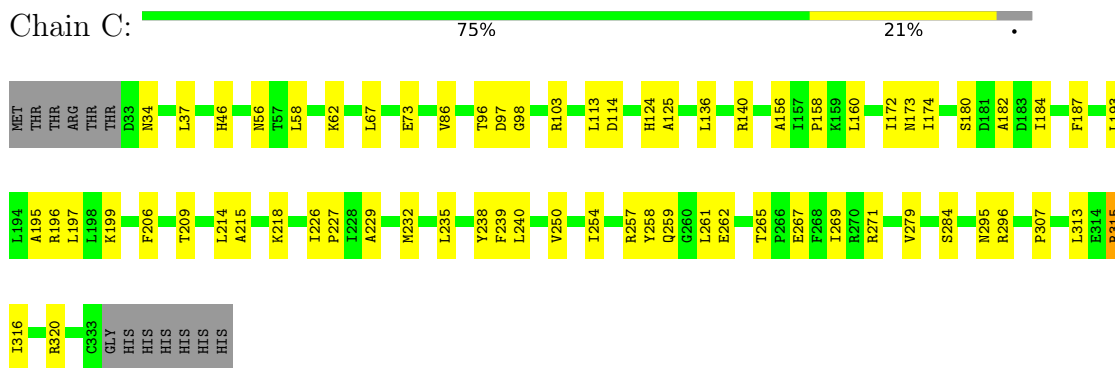
- Molecule 1: Lysoplasmalogenase



- Molecule 1: Lysoplasmalogenase

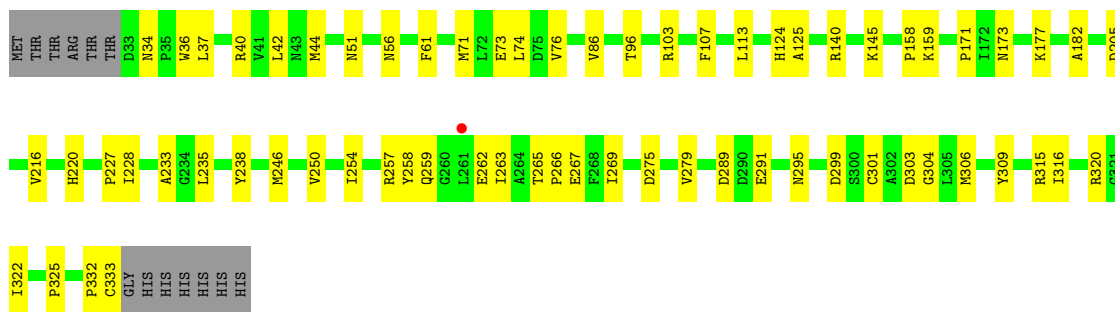


- Molecule 1: Lysoplasmalogenase




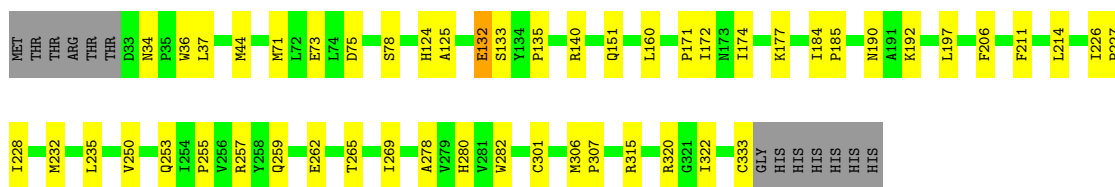
- Molecule 1: Lysoplasmalogenase

Chain D:  75% 21%



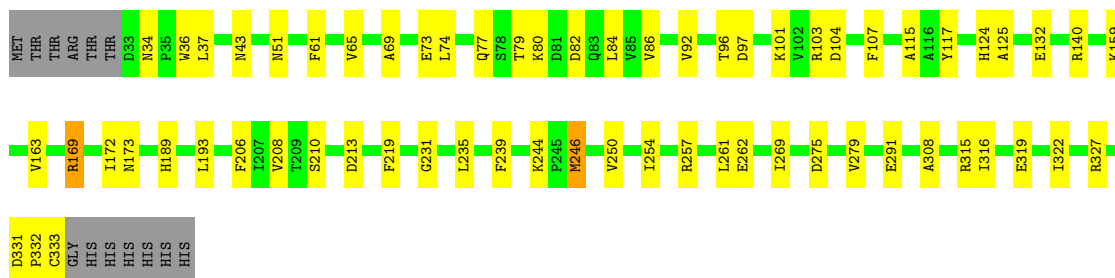
• Molecule 1: Lysoplasmalogenase

Chain E:  80% 16%



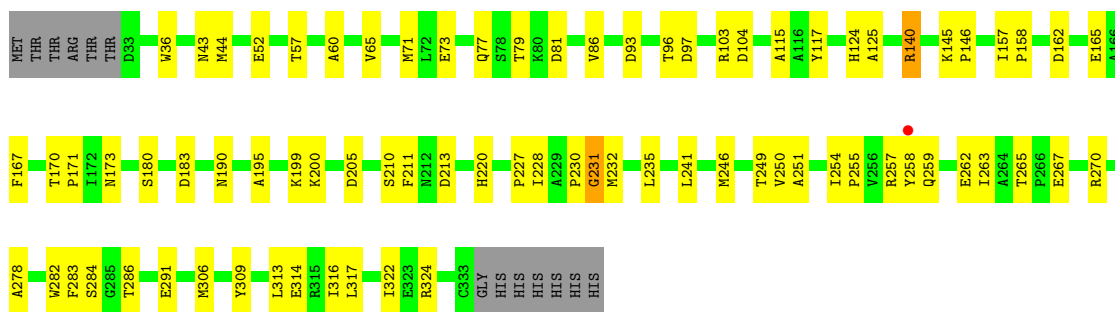
• Molecule 1: Lysoplasmalogenase

Chain F:  75% 20%



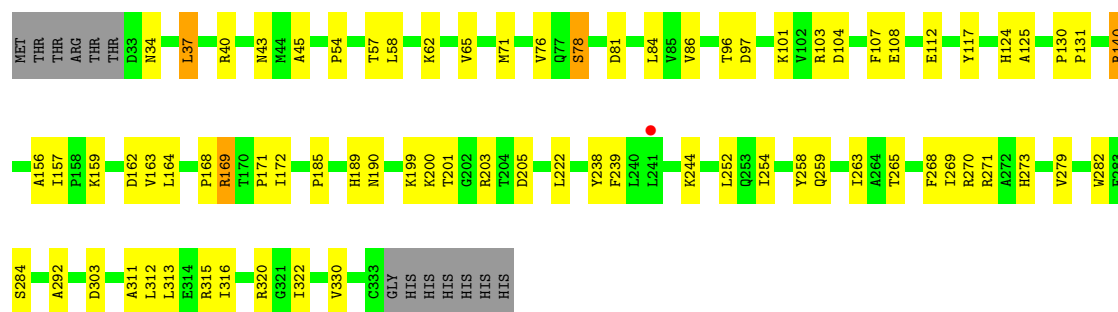
• Molecule 1: Lysoplasmalogenase

Chain G:  71% 25%



• Molecule 1: Lysoplasmalogenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	112.61Å 94.94Å 147.35Å 90.00° 90.24° 90.00°	Depositor
Resolution (Å)	47.47 – 2.57 48.43 – 2.57	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.47-2.57) 98.1 (48.43-2.57)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.22 (at 2.58Å)	Xtrriage
Refinement program	PHENIX 1.13_2998	Depositor
R, $R_{free}$	0.210 , 0.272 0.209 , 0.270	Depositor DCC
$R_{free}$ test set	4777 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	42.3	Xtrriage
Anisotropy	0.866	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 24.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtrriage
Estimated twinning fraction	0.237 for h,-k,-l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	18857	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.83% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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	0.47	0/2416	0.63	0/3297
1	B	0.47	0/2377	0.64	0/3243
1	C	0.44	0/2377	0.63	0/3243
1	D	0.45	0/2377	0.61	0/3243
1	E	0.44	0/2377	0.61	0/3243
1	F	0.42	0/2377	0.60	0/3243
1	G	0.43	0/2377	0.60	0/3243
1	H	0.39	0/2377	0.57	0/3243
All	All	0.44	0/19055	0.61	0/25998

There are no bond length outliers.

There are no bond angle outliers.

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	2356	0	2334	43	0
1	B	2317	0	2293	45	0
1	C	2317	0	2293	41	0
1	D	2317	0	2292	43	0
1	E	2317	0	2293	36	0
1	F	2317	0	2293	42	0
1	G	2317	0	2293	59	0
1	H	2317	0	2293	47	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	35	0	0	2	0
2	B	45	0	0	1	0
2	C	40	0	0	2	0
2	D	29	0	0	0	0
2	E	36	0	0	1	0
2	F	30	0	0	2	0
2	G	42	0	0	3	0
2	H	25	0	0	1	0
All	All	18857	0	18384	348	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 9.

The worst 5 of 348 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:291:GLU:HG3	1:B:316:ILE:HD11	1.56	0.88
1:C:257:ARG:HH22	1:C:296:ARG:HH22	1.22	0.83
1:G:103:ARG:HD2	2:G:423:HOH:O	1.76	0.83
1:F:169:ARG:O	1:F:169:ARG:HD3	1.83	0.79
1:A:197:LEU:HA	1:A:200:LYS:HE3	1.65	0.77

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	304/314 (97%)	288 (95%)	16 (5%)	0	100 100
1	B	299/314 (95%)	282 (94%)	17 (6%)	0	100 100
1	C	299/314 (95%)	283 (95%)	15 (5%)	1 (0%)	41 62

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	299/314 (95%)	283 (95%)	16 (5%)	0	100	100
1	E	299/314 (95%)	280 (94%)	19 (6%)	0	100	100
1	F	299/314 (95%)	282 (94%)	17 (6%)	0	100	100
1	G	299/314 (95%)	284 (95%)	14 (5%)	1 (0%)	41	62
1	H	299/314 (95%)	282 (94%)	16 (5%)	1 (0%)	41	62
All	All	2397/2512 (95%)	2264 (94%)	130 (5%)	3 (0%)	51	73

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	231	GLY
1	H	45	ALA
1	C	46	HIS

### 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	248/255 (97%)	242 (98%)	6 (2%)	49	72
1	B	243/255 (95%)	239 (98%)	4 (2%)	62	81
1	C	243/255 (95%)	235 (97%)	8 (3%)	38	61
1	D	243/255 (95%)	239 (98%)	4 (2%)	62	81
1	E	243/255 (95%)	237 (98%)	6 (2%)	47	70
1	F	243/255 (95%)	238 (98%)	5 (2%)	53	75
1	G	243/255 (95%)	239 (98%)	4 (2%)	62	81
1	H	243/255 (95%)	237 (98%)	6 (2%)	47	70
All	All	1949/2040 (96%)	1906 (98%)	43 (2%)	52	74

5 of 43 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	101	LYS
1	G	200	LYS
1	F	103	ARG
1	F	333	CYS
1	H	37	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	306/314 (97%)	-0.24	0 100   100	25, 35, 52, 78	0
1	B	301/314 (95%)	-0.24	1 (0%) 94   94	24, 35, 51, 71	0
1	C	301/314 (95%)	-0.21	0 100   100	24, 42, 62, 82	0
1	D	301/314 (95%)	-0.16	1 (0%) 94   94	26, 44, 66, 83	0
1	E	301/314 (95%)	-0.22	0 100   100	25, 41, 59, 81	0
1	F	301/314 (95%)	-0.10	0 100   100	34, 49, 64, 79	0
1	G	301/314 (95%)	-0.26	1 (0%) 94   94	29, 43, 60, 80	0
1	H	301/314 (95%)	-0.02	1 (0%) 94   94	34, 54, 71, 83	0
All	All	2413/2512 (96%)	-0.18	4 (0%) 95   95	24, 43, 64, 83	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	258	TYR	3.2
1	B	258	TYR	2.9
1	H	241	LEU	2.2
1	D	261	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.