



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 04:44 AM GMT

PDB ID : 2O18  
Title : Crystal structure of a Thiamine biosynthesis lipoprotein apbE, NorthEast Structural Genomics target ER559  
Authors : Seetharaman, J.; Su, M.; Wang, D.; Fang, Y.; Cunningham, K.; Ma, L.; Xiao, R.; Liu, J.; Baran, M.C.; Acton, T.B.; Rost, B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2006-11-28  
Resolution : 2.20 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)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

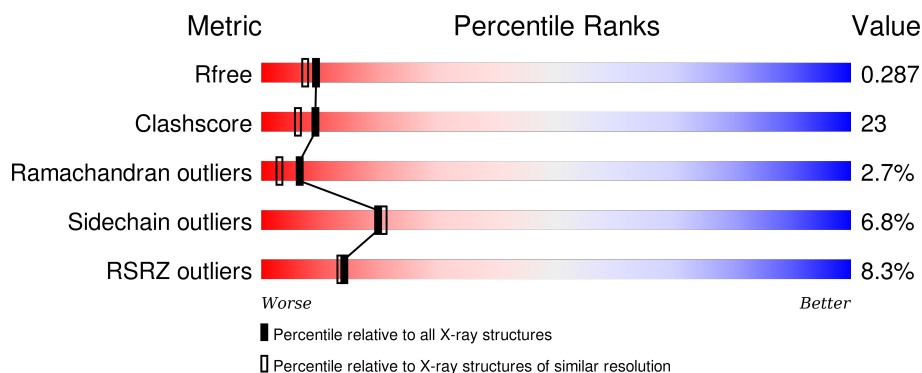
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	340	<div> <div>4%</div> <div>56%</div> <div>29%</div> <div>10%</div> </div>
1	B	340	<div> <div>8%</div> <div>55%</div> <div>29%</div> <div>12%</div> </div>
1	C	340	<div> <div>9%</div> <div>55%</div> <div>30%</div> <div>7%</div> <div>8%</div> </div>
1	D	340	<div> <div>8%</div> <div>58%</div> <div>28%</div> <div>10%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9555 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 Thiamine biosynthesis lipoprotein apbE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	307	Total	C	N	O	Se	0	0	0
			2351	1476	406	459	10			
1	B	300	Total	C	N	O	Se	0	0	0
			2307	1449	396	452	10			
1	C	313	Total	C	N	O	Se	0	0	0
			2374	1491	410	463	10			
1	D	306	Total	C	N	O	Se	0	0	0
			2316	1457	398	451	10			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	INITIATING METHIONINE	UNP P0AB85
A	23	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	67	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	84	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	102	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	133	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	180	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	201	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	287	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	307	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	319	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
A	333	LEU	-	CLONING ARTIFACT	UNP P0AB85
A	334	GLU	-	CLONING ARTIFACT	UNP P0AB85
A	335	HIS	-	EXPRESSION TAG	UNP P0AB85
A	336	HIS	-	EXPRESSION TAG	UNP P0AB85
A	337	HIS	-	EXPRESSION TAG	UNP P0AB85
A	338	HIS	-	EXPRESSION TAG	UNP P0AB85
A	339	HIS	-	EXPRESSION TAG	UNP P0AB85
A	340	HIS	-	EXPRESSION TAG	UNP P0AB85
B	1	MSE	-	INITIATING METHIONINE	UNP P0AB85
B	23	MSE	MET	MODIFIED RESIDUE	UNP P0AB85

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Chain	Residue	Modelled	Actual	Comment	Reference
B	67	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
B	84	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
B	102	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
B	133	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
B	180	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
B	201	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
B	287	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
B	307	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
B	319	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
B	333	LEU	-	CLONING ARTIFACT	UNP P0AB85
B	334	GLU	-	CLONING ARTIFACT	UNP P0AB85
B	335	HIS	-	EXPRESSION TAG	UNP P0AB85
B	336	HIS	-	EXPRESSION TAG	UNP P0AB85
B	337	HIS	-	EXPRESSION TAG	UNP P0AB85
B	338	HIS	-	EXPRESSION TAG	UNP P0AB85
B	339	HIS	-	EXPRESSION TAG	UNP P0AB85
B	340	HIS	-	EXPRESSION TAG	UNP P0AB85
C	1	MSE	-	INITIATING METHIONINE	UNP P0AB85
C	23	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	67	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	84	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	102	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	133	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	180	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	201	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	287	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	307	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	319	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
C	333	LEU	-	CLONING ARTIFACT	UNP P0AB85
C	334	GLU	-	CLONING ARTIFACT	UNP P0AB85
C	335	HIS	-	EXPRESSION TAG	UNP P0AB85
C	336	HIS	-	EXPRESSION TAG	UNP P0AB85
C	337	HIS	-	EXPRESSION TAG	UNP P0AB85
C	338	HIS	-	EXPRESSION TAG	UNP P0AB85
C	339	HIS	-	EXPRESSION TAG	UNP P0AB85
C	340	HIS	-	EXPRESSION TAG	UNP P0AB85
D	1	MSE	-	INITIATING METHIONINE	UNP P0AB85
D	23	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
D	67	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
D	84	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
D	102	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
D	133	MSE	MET	MODIFIED RESIDUE	UNP P0AB85

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Chain	Residue	Modelled	Actual	Comment	Reference
D	180	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
D	201	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
D	287	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
D	307	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
D	319	MSE	MET	MODIFIED RESIDUE	UNP P0AB85
D	333	LEU	-	CLONING ARTIFACT	UNP P0AB85
D	334	GLU	-	CLONING ARTIFACT	UNP P0AB85
D	335	HIS	-	EXPRESSION TAG	UNP P0AB85
D	336	HIS	-	EXPRESSION TAG	UNP P0AB85
D	337	HIS	-	EXPRESSION TAG	UNP P0AB85
D	338	HIS	-	EXPRESSION TAG	UNP P0AB85
D	339	HIS	-	EXPRESSION TAG	UNP P0AB85
D	340	HIS	-	EXPRESSION TAG	UNP P0AB85

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ca 1 1	0	0
2	A	1	Total Ca 1 1	0	0
2	D	1	Total Ca 1 1	0	0
2	C	1	Total Ca 1 1	0	0

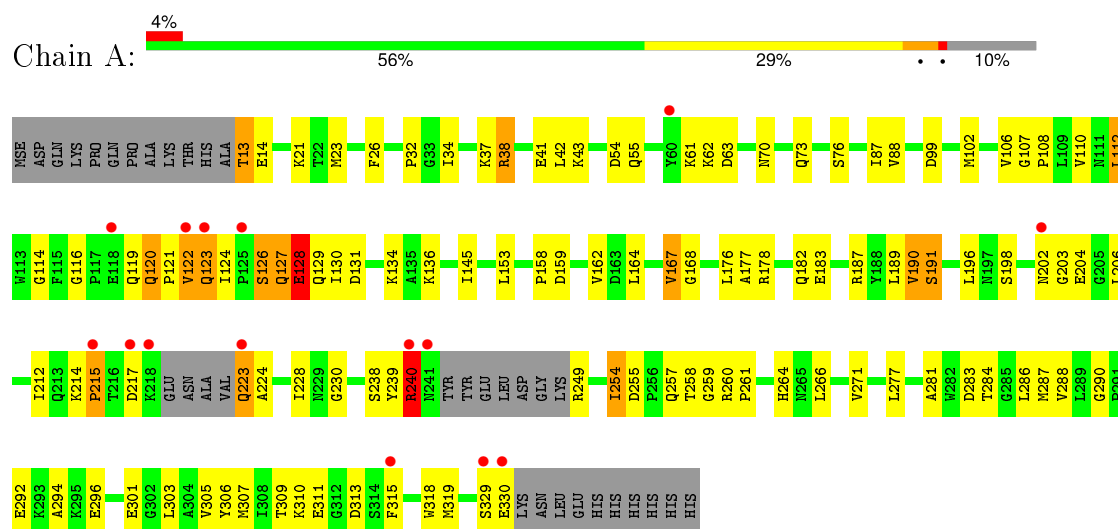
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	52	Total O 52 52	0	0
3	B	47	Total O 47 47	0	0
3	C	48	Total O 48 48	0	0
3	D	56	Total O 56 56	0	0

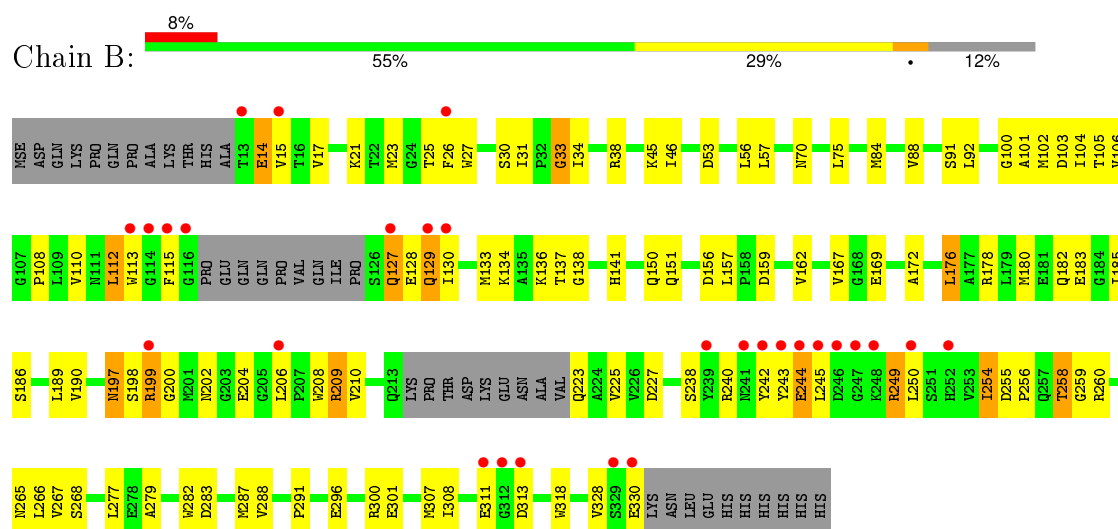
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Thiamine biosynthesis lipoprotein apbE

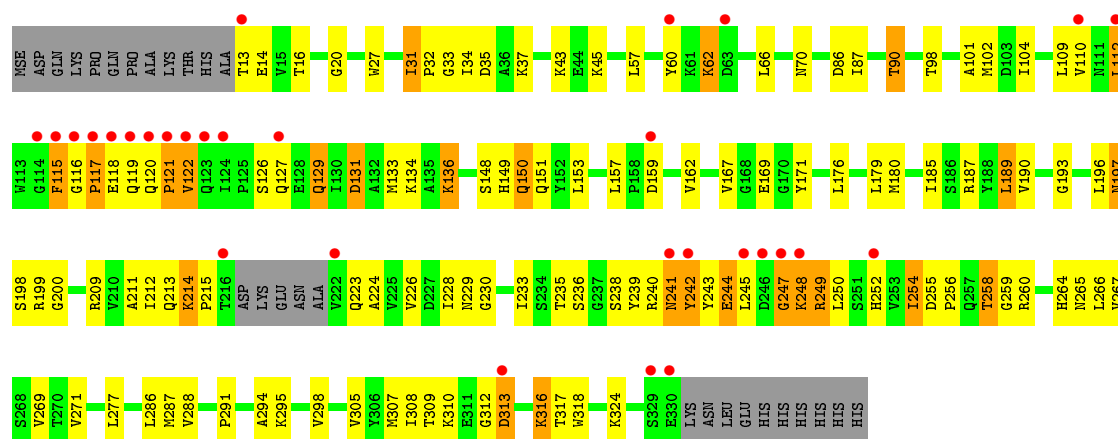


- Molecule 1: Thiamine biosynthesis lipoprotein apbE

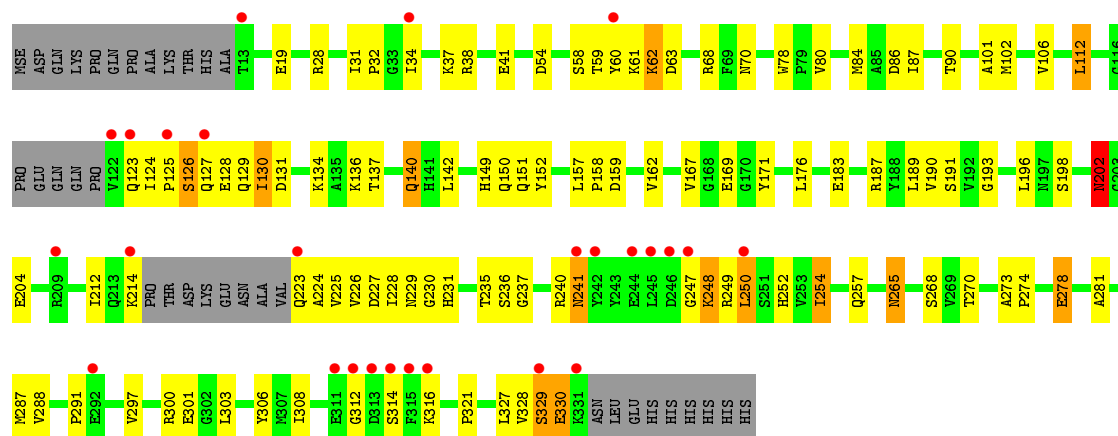


- Molecule 1: Thiamine biosynthesis lipoprotein apbE





• Molecule 1: Thiamine biosynthesis lipoprotein apbE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.32Å 70.65Å 86.64Å 75.76° 71.67° 69.46°	Depositor
Resolution (Å)	49.74 – 2.20 49.74 – 2.20	Depositor EDS
% Data completeness (in resolution range)	78.0 (49.74-2.20) 90.3 (49.74-2.20)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.86 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.233 , 0.284 0.237 , 0.287	Depositor DCC
$R_{free}$ test set	2486 reflections (4.85%)	DCC
Wilson B-factor (Å <sup>2</sup> )	20.9	Xtriage
Anisotropy	0.358	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 43.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 114432 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	9555	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.59% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	0.35	0/2383	0.65	1/3215 (0.0%)
1	B	0.34	0/2338	0.63	1/3149 (0.0%)
1	C	0.36	0/2407	0.63	1/3250 (0.0%)
1	D	0.36	0/2346	0.62	0/3165
All	All	0.35	0/9474	0.64	3/12779 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	121	PRO	N-CA-CB	5.60	110.03	103.30
1	B	243	TYR	N-CA-C	5.51	125.89	111.00
1	A	215	PRO	N-CA-CB	5.22	109.57	103.30

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	2351	0	2321	106	0
1	B	2307	0	2275	113	0
1	C	2374	0	2333	129	0
1	D	2316	0	2277	89	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	52	0	0	7	0
3	B	47	0	0	9	0
3	C	48	0	0	5	0
3	D	56	0	0	4	0
All	All	9555	0	9206	427	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 23.

The worst 5 of 427 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:A:202:ASN:HB3	3:A:343:HOH:O	1.71	0.89
1:B:189:LEU:HD12	1:B:208:TRP:HB3	1.53	0.89
1:B:46:ILE:HG23	1:B:176:LEU:HD21	1.56	0.86
1:A:112:LEU:HD22	1:A:126:SER:HB3	1.56	0.86
1:A:120:GLN:HB2	1:A:121:PRO:C	1.95	0.86

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	301/340 (88%)	275 (91%)	18 (6%)	8 (3%)	<b>6</b> <b>3</b>
1	B	294/340 (86%)	268 (91%)	21 (7%)	5 (2%)	<b>11</b> <b>7</b>
1	C	309/340 (91%)	273 (88%)	24 (8%)	12 (4%)	<b>4</b> <b>1</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	300/340 (88%)	273 (91%)	20 (7%)	7 (2%)	8	4
All	All	1204/1360 (88%)	1089 (90%)	83 (7%)	32 (3%)	6	3

5 of 32 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	GLN
1	C	35	ASP
1	C	117	PRO
1	C	241	ASN
1	C	242	TYR

### 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	254/277 (92%)	238 (94%)	16 (6%)	22	24
1	B	248/277 (90%)	232 (94%)	16 (6%)	21	23
1	C	253/277 (91%)	234 (92%)	19 (8%)	17	17
1	D	246/277 (89%)	229 (93%)	17 (7%)	19	20
All	All	1001/1108 (90%)	933 (93%)	68 (7%)	20	21

5 of 68 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	313	ASP
1	C	129	GLN
1	D	227	ASP
1	B	330	GLU
1	C	62	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 29 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	197	ASN
1	C	70	ASN
1	D	202	ASN
1	B	241	ASN
1	C	147	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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

### 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	297/340 (87%)	0.32	15 (5%) 32 31	7, 20, 52, 61	0
1	B	290/340 (85%)	0.65	28 (9%) 10 9	9, 25, 52, 73	0
1	C	303/340 (89%)	0.63	30 (9%) 9 8	9, 22, 56, 72	0
1	D	296/340 (87%)	0.53	26 (8%) 12 11	6, 22, 54, 67	0
All	All	1186/1360 (87%)	0.53	99 (8%) 14 13	6, 22, 53, 73	0

The worst 5 of 99 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	246	ASP	12.0
1	C	122	VAL	10.6
1	C	116	GLY	9.6
1	A	122	VAL	8.5
1	C	121	PRO	7.6

### 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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors

of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	CA	A	341	1/1	0.87	0.15	1.70	23,23,23,23	0
2	CA	C	341	1/1	0.97	0.17	1.09	13,13,13,13	0
2	CA	B	341	1/1	0.95	0.15	-0.25	24,24,24,24	0
2	CA	D	341	1/1	0.95	0.05	-3.92	20,20,20,20	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.