



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:08 PM GMT

PDB ID : 4NVR  
Title : 2.22 Angstrom Resolution Crystal Structure of a Putative Acyltransferase from *Salmonella enterica*  
Authors : Minasov, G.; Wawrzak, Z.; Skarina, T.; Gordon, E.; Stam, J.; Kwon, K.; Savchenko, A.; Anderson, W.F.; Center for Structural Genomics of Infectious Diseases (CSGID)  
Deposited on : 2013-12-05  
Resolution : 2.22 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](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.

---

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

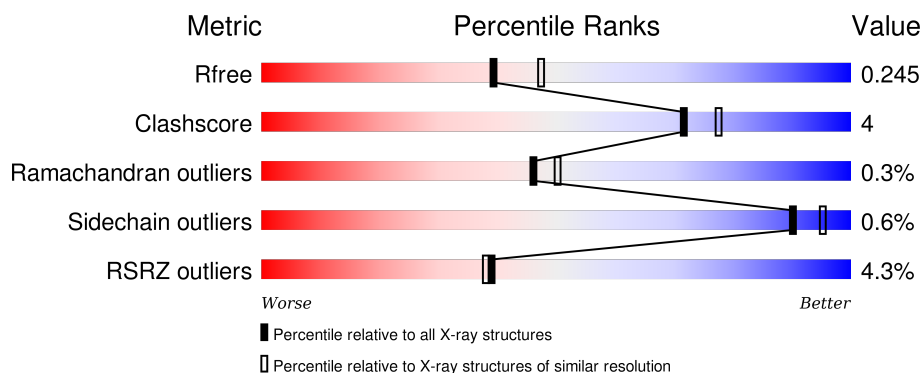
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.22 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	4405 (2.24-2.20)
Clashscore	102246	5146 (2.24-2.20)
Ramachandran outliers	100387	5065 (2.24-2.20)
Sidechain outliers	100360	5066 (2.24-2.20)
RSRZ outliers	91569	4414 (2.24-2.20)

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	307	<div> <div>6%</div> <div> <div></div> <div>88%</div> <div>10%</div> <div></div> </div> </div>
1	B	307	<div> <div>4%</div> <div> <div></div> <div>90%</div> <div>8%</div> <div></div> </div> </div>
1	C	307	<div> <div>4%</div> <div> <div></div> <div>87%</div> <div>10%</div> <div></div> </div> </div>
1	D	307	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>10%</div> <div></div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CA	C	402	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10229 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 Putative acyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	Se	0	10	0
			2458	1564	430	450	8	6			
1	B	302	Total	C	N	O	S	Se	0	2	0
			2386	1522	414	436	8	6			
1	C	301	Total	C	N	O	S	Se	0	4	0
			2401	1533	419	435	8	6			
1	D	301	Total	C	N	O	S	Se	0	0	0
			2367	1512	411	430	8	6			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q8ZRI7
A	-1	ASN	-	EXPRESSION TAG	UNP Q8ZRI7
A	0	ALA	-	EXPRESSION TAG	UNP Q8ZRI7
B	-2	SER	-	EXPRESSION TAG	UNP Q8ZRI7
B	-1	ASN	-	EXPRESSION TAG	UNP Q8ZRI7
B	0	ALA	-	EXPRESSION TAG	UNP Q8ZRI7
C	-2	SER	-	EXPRESSION TAG	UNP Q8ZRI7
C	-1	ASN	-	EXPRESSION TAG	UNP Q8ZRI7
C	0	ALA	-	EXPRESSION TAG	UNP Q8ZRI7
D	-2	SER	-	EXPRESSION TAG	UNP Q8ZRI7
D	-1	ASN	-	EXPRESSION TAG	UNP Q8ZRI7
D	0	ALA	-	EXPRESSION TAG	UNP Q8ZRI7

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

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

*Continued on next page...*

*Continued from previous page...*

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

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Cl 1	0	0
3	A	1	Total 1	Cl 1	0	0

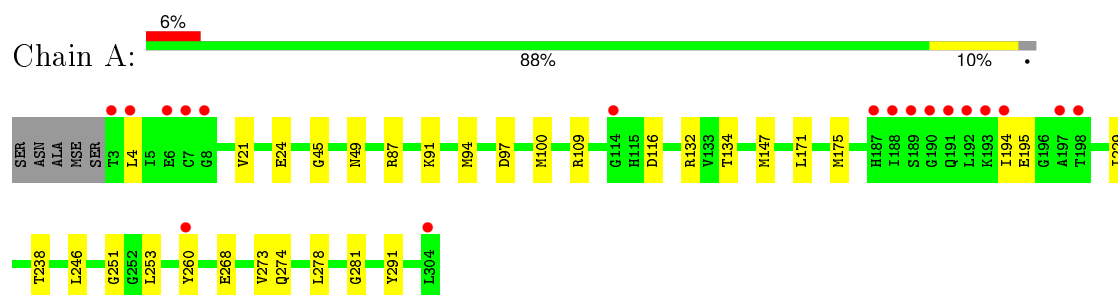
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	147	Total 155	O 155	0	9
4	B	165	Total 166	O 166	0	1
4	C	141	Total 144	O 144	0	4
4	D	144	Total 145	O 145	0	1

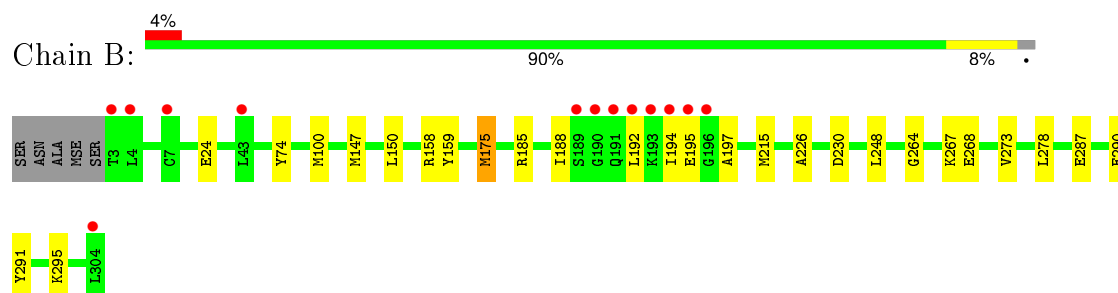
### 3 Residue-property plots [i](#)

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.

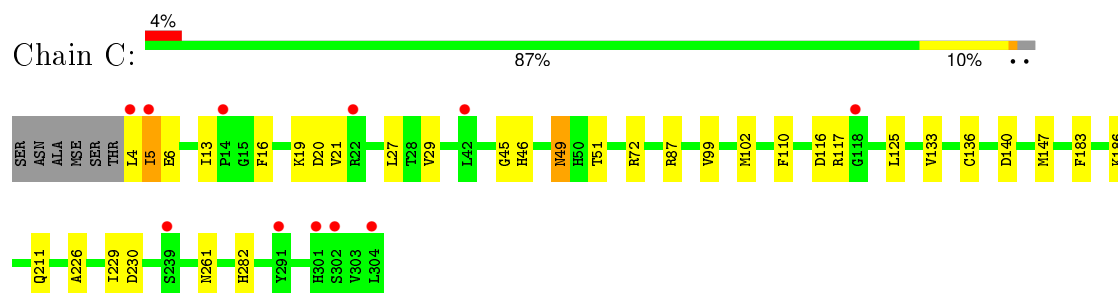
- Molecule 1: Putative acyltransferase



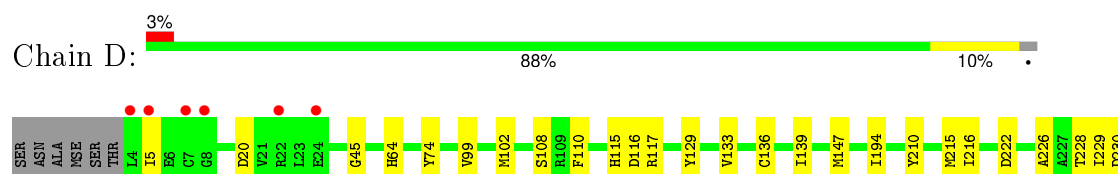
- Molecule 1: Putative acyltransferase

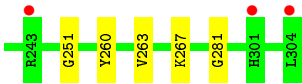


- Molecule 1: Putative acyltransferase



- Molecule 1: Putative acyltransferase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	47.93Å 71.38Å 90.33Å 113.03° 92.41° 90.09°	Depositor
Resolution (Å)	28.63 – 2.22 28.63 – 2.22	Depositor EDS
% Data completeness (in resolution range)	98.1 (28.63-2.22) 85.0 (28.63-2.22)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.64 (at 2.22Å)	Xtriage
Refinement program	REFMAC 5.8.0046	Depositor
R, $R_{free}$	0.181 , 0.241 0.191 , 0.245	Depositor DCC
$R_{free}$ test set	2697 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	22.5	Xtriage
Anisotropy	0.448	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 41.7	EDS
Estimated twinning fraction	0.079 for h,-k,-l 0.017 for -h,k,-k-l 0.006 for -h,-k,k+l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 53363 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10229	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% 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, CL

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	0.59	0/2518	0.78	1/3414 (0.0%)
1	B	0.59	0/2446	0.78	3/3319 (0.1%)
1	C	0.57	0/2461	0.75	2/3335 (0.1%)
1	D	0.56	0/2427	0.74	0/3291
All	All	0.58	0/9852	0.76	6/13359 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	268	GLU	OE1-CD-OE2	8.23	133.18	123.30
1	A	268	GLU	OE1-CD-OE2	7.57	132.39	123.30
1	B	185	ARG	NE-CZ-NH1	5.52	123.06	120.30
1	C	147	MSE	CA-CB-CG	-5.23	104.41	113.30
1	B	185	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	C	261	ASN	N-CA-C	-5.02	97.45	111.00

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	2458	0	2391	23	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2386	0	2327	15	0
1	C	2401	0	2350	26	0
1	D	2367	0	2311	18	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	155	0	0	2	0
4	B	166	0	0	2	0
4	C	144	0	0	2	0
4	D	145	0	0	3	0
All	All	10229	0	9379	81	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 4.

All (81) 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:147:MSE:HE1	1:A:260:TYR:CE2	2.23	0.73
1:D:5:ILE:HG23	1:D:20:ASP:HB2	1.72	0.69
1:A:147:MSE:HA	1:A:147:MSE:HE2	1.74	0.69
1:C:4:LEU:HD22	1:C:21:VAL:HG22	1.77	0.67
1:D:64:HIS:NE2	4:D:556:HOH:O	2.28	0.67
1:A:24:GLU:HG2	1:A:100:MSE:HE2	1.77	0.67
1:C:4:LEU:CD2	1:C:21:VAL:HG22	2.27	0.65
1:D:133:VAL:HG11	1:D:136:CYS:SG	2.41	0.60
1:C:116:ASP:OD1	1:C:117:ARG:HG2	2.03	0.59
1:A:238:THR:HG23	4:A:631[B]:HOH:O	2.03	0.59
1:A:4:LEU:HD22	1:A:21:VAL:HG22	1.83	0.59
1:D:194:ILE:HG22	4:D:628:HOH:O	2.03	0.56
1:C:49:ASN:ND2	1:C:51:THR:H	2.04	0.55
1:C:5:ILE:HG23	1:C:20:ASP:HB2	1.88	0.55
1:A:147:MSE:HE1	1:A:260:TYR:CD2	2.42	0.55
1:A:109[A]:ARG:NH1	1:A:134:THR:HG22	2.22	0.55
1:A:45:GLY:HA3	1:A:116:ASP:HB3	1.88	0.54
1:C:45:GLY:HA3	1:C:116:ASP:HB3	1.88	0.54
1:A:175:MSE:HE1	1:C:183:PHE:CE1	2.43	0.53
1:D:74:TYR:CD2	1:D:215:MSE:HE3	2.45	0.51

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:251:GLY:HA3	1:D:281:GLY:O	2.10	0.50
1:C:27:LEU:CD2	1:C:29:VAL:HG13	2.41	0.50
1:B:147:MSE:HE1	1:B:159:TYR:CZ	2.47	0.50
1:A:94:MSE:O	1:A:97:ASP:HB2	2.11	0.49
1:D:102:MSE:HE2	1:D:110:PHE:CG	2.47	0.49
1:A:246:LEU:HD23	1:A:273[B]:VAL:HG12	1.94	0.49
1:C:102:MSE:HG3	1:C:110:PHE:CD2	2.48	0.48
1:C:99:VAL:HG13	1:C:125:LEU:HD11	1.95	0.48
1:C:211:GLN:NE2	4:C:524:HOH:O	2.46	0.47
1:B:290:GLU:HB2	4:B:645:HOH:O	2.14	0.47
1:C:186[B]:LYS:HD3	1:C:186[B]:LYS:HA	1.70	0.47
1:C:49:ASN:HD22	1:C:49:ASN:C	2.18	0.47
1:B:24:GLU:CG	1:B:100:MSE:HE2	2.46	0.46
1:B:264:GLY:HA2	1:B:267:LYS:HD3	1.97	0.46
1:A:4:LEU:HD23	1:A:21:VAL:HG13	1.97	0.46
1:C:6:GLU:OE2	1:C:19:LYS:NZ	2.48	0.46
1:C:140:ASP:OD2	1:C:282:HIS:ND1	2.38	0.46
1:A:278:LEU:HD22	1:A:291:TYR:CD1	2.51	0.46
1:A:273[A]:VAL:HG12	1:A:274:GLN:N	2.32	0.45
1:B:150:LEU:HB2	4:B:529:HOH:O	2.17	0.45
1:B:194:ILE:HD11	1:B:287:GLU:HB3	1.98	0.45
1:C:27:LEU:HD13	1:C:72:ARG:NH2	2.32	0.45
1:B:226:ALA:HA	1:B:230:ASP:HB2	1.99	0.45
1:B:188:ILE:O	1:B:192:LEU:HB2	2.16	0.45
1:C:226:ALA:HA	1:C:230:ASP:HB2	1.99	0.45
1:C:133:VAL:HG11	1:C:136:CYS:SG	2.57	0.45
1:B:192:LEU:HD21	1:B:197:ALA:HB3	1.99	0.44
1:A:24:GLU:CG	1:A:100:MSE:HE2	2.47	0.44
1:D:210:TYR:O	1:D:216:ILE:HD11	2.18	0.44
1:D:117:ARG:HD2	1:D:222:ASP:OD2	2.17	0.44
1:D:147:MSE:HG2	1:D:260:TYR:CZ	2.51	0.44
1:A:251:GLY:HA3	1:A:281:GLY:O	2.18	0.44
1:C:87:ARG:NH2	1:C:229:ILE:HD11	2.33	0.44
1:A:49:ASN:HB3	4:A:506:HOH:O	2.17	0.44
1:B:194:ILE:O	1:B:194:ILE:HG13	2.18	0.44
1:C:102:MSE:HG3	1:C:110:PHE:CE2	2.52	0.43
1:C:46:HIS:CE1	4:C:616:HOH:O	2.70	0.43
1:A:171:LEU:O	1:A:171:LEU:HD12	2.17	0.43
1:D:117:ARG:CD	1:D:222:ASP:OD2	2.66	0.43
1:D:45:GLY:HA3	1:D:116:ASP:HB3	1.99	0.43
1:B:24:GLU:HG2	1:B:100:MSE:HE2	2.01	0.43

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:175:MSE:HE1	4:D:599:HOH:O	2.18	0.43
1:B:248:LEU:CD2	1:B:273:VAL:HG13	2.48	0.43
1:C:116:ASP:OD1	1:C:117:ARG:N	2.49	0.42
1:D:226:ALA:HA	1:D:230:ASP:OD2	2.19	0.42
1:A:171:LEU:C	1:A:171:LEU:HD12	2.40	0.42
1:D:116:ASP:OD1	1:D:117:ARG:HG2	2.19	0.42
1:C:87:ARG:CZ	1:C:229:ILE:HD11	2.49	0.42
1:B:278:LEU:HD22	1:B:291:TYR:CD1	2.55	0.42
1:B:74:TYR:CD2	1:B:215:MSE:HE3	2.55	0.42
1:C:13:ILE:HG22	1:C:16:PHE:CD1	2.55	0.42
1:C:49:ASN:ND2	1:C:49:ASN:C	2.74	0.41
1:D:263:VAL:HG12	1:D:267:LYS:HE2	2.01	0.41
1:A:253:LEU:HD22	1:A:253:LEU:N	2.35	0.41
1:C:5:ILE:HD13	1:C:5:ILE:C	2.41	0.41
1:D:228:THR:OG1	1:D:229:ILE:N	2.53	0.41
1:D:115:HIS:HA	1:D:139:ILE:HB	2.04	0.40
1:A:4:LEU:CD2	1:A:21:VAL:HG13	2.51	0.40
1:A:109[A]:ARG:HA	1:A:132:ARG:O	2.22	0.40
1:D:99:VAL:HG21	1:D:129:TYR:CG	2.57	0.40
1:A:87:ARG:NH2	1:A:229:ILE:HD11	2.37	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	310/307 (101%)	300 (97%)	8 (3%)	2 (1%)	30	29
1	B	302/307 (98%)	291 (96%)	10 (3%)	1 (0%)	46	50
1	C	303/307 (99%)	295 (97%)	8 (3%)	0	100	100
1	D	299/307 (97%)	292 (98%)	6 (2%)	1 (0%)	46	50
All	All	1214/1228 (99%)	1178 (97%)	32 (3%)	4 (0%)	46	50

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	ILE
1	A	195	GLU
1	B	195	GLU
1	D	108	SER

### 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	262/249 (105%)	261 (100%)	1 (0%)	93	97
1	B	254/249 (102%)	251 (99%)	3 (1%)	78	88
1	C	255/249 (102%)	253 (99%)	2 (1%)	86	93
1	D	251/249 (101%)	251 (100%)	0	100	100
All	All	1022/996 (103%)	1016 (99%)	6 (1%)	90	95

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	LYS
1	B	158	ARG
1	B	175	MSE
1	B	295	LYS
1	C	5	ILE
1	C	49	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	HIS
1	A	187	HIS
1	B	63	ASN
1	B	187	HIS
1	C	46	HIS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	C	187	HIS
1	C	211	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 7 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 ⓘ

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, 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	296/307 (96%)	0.09	18 (6%) 25 24	15, 25, 56, 96	0
1	B	296/307 (96%)	0.06	13 (4%) 38 37	14, 25, 48, 97	0
1	C	295/307 (96%)	0.14	11 (3%) 45 44	15, 28, 47, 58	0
1	D	295/307 (96%)	0.05	9 (3%) 52 51	15, 27, 48, 76	0
All	All	1182/1228 (96%)	0.08	51 (4%) 39 38	14, 26, 49, 97	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	194	ILE	9.4
1	A	194	ILE	6.8
1	D	4	LEU	6.5
1	C	4	LEU	5.5
1	C	301	HIS	4.2
1	A	187	HIS	4.1
1	B	3	THR	4.0
1	B	192	LEU	3.8
1	A	193	LYS	3.7
1	D	301	HIS	3.5
1	C	5	ILE	3.5
1	B	193	LYS	3.4
1	A	3	THR	3.4
1	D	5	ILE	3.2
1	A	188	ILE	3.2
1	A	191	GLN	3.1
1	B	4	LEU	3.1
1	A	4	LEU	3.0
1	A	304	LEU	3.0
1	C	14	PRO	3.0
1	A	7	CYS	2.9

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	304	LEU	2.9
1	B	191	GLN	2.8
1	A	8	GLY	2.7
1	C	291	TYR	2.7
1	B	190	GLY	2.6
1	B	7	CYS	2.6
1	A	192	LEU	2.6
1	C	302	SER	2.5
1	A	197	ALA	2.5
1	C	118	GLY	2.4
1	D	8	GLY	2.4
1	A	198	THR	2.4
1	B	43	LEU	2.4
1	D	243	ARG	2.4
1	D	22	ARG	2.4
1	B	195	GLU	2.3
1	A	114	GLY	2.3
1	D	24	GLU	2.3
1	A	190	GLY	2.3
1	A	260	TYR	2.3
1	C	239	SER	2.2
1	A	189	SER	2.1
1	D	7	CYS	2.1
1	C	304	LEU	2.1
1	C	22	ARG	2.1
1	B	196	GLY	2.1
1	C	42	LEU	2.1
1	D	304	LEU	2.1
1	A	6[A]	GLU	2.0
1	B	189	SER	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

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(Å <sup>2</sup> )	Q<0.9
2	CA	C	402	1/1	0.94	0.31	5.56	52,52,52,52	0
2	CA	B	401	1/1	0.98	0.11	-1.84	28,28,28,28	0
2	CA	A	401	1/1	0.97	0.09	-2.22	28,28,28,28	0
3	CL	A	402	1/1	0.97	0.19	-	30,30,30,30	0
3	CL	B	402	1/1	0.97	0.18	-	30,30,30,30	0
2	CA	C	401	1/1	0.89	0.32	-	51,51,51,51	0
2	CA	D	401	1/1	0.84	0.27	-	57,57,57,57	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.