



# Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 03:49 pm BST

PDB ID : 3GBX  
Title : Serine hydroxymethyltransferase from *Salmonella typhimurium*  
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Deposited on : 2009-02-20  
Resolution : 1.80 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

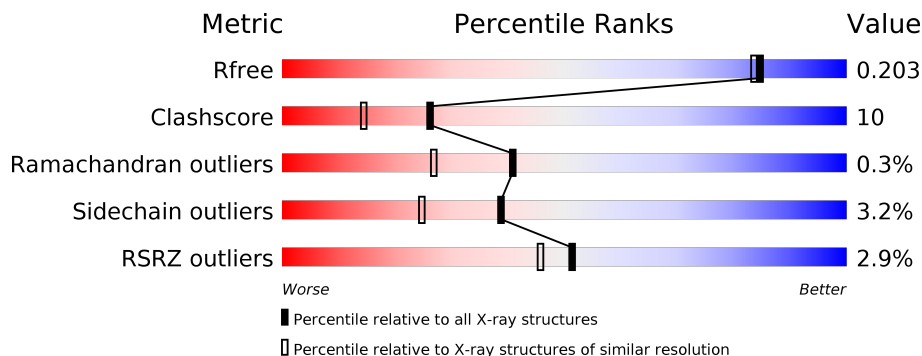
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 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	420	 2% 79% 12% • 7%
1	B	420	 3% 74% 16% • 9%

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6646 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.

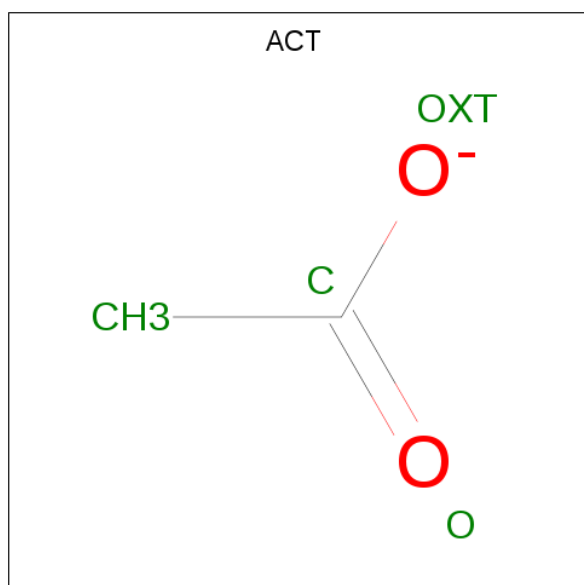
- Molecule 1 is a protein called Serine hydroxymethyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	N	O	S				Se
1	A	390	3119	1988	518	596	2	15	0	24	0
1	B	383	3073	1961	519	575	2	16	0	24	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP P0A2E1
A	-1	ASN	-	EXPRESSION TAG	UNP P0A2E1
A	0	ALA	-	EXPRESSION TAG	UNP P0A2E1
B	-2	SER	-	EXPRESSION TAG	UNP P0A2E1
B	-1	ASN	-	EXPRESSION TAG	UNP P0A2E1
B	0	ALA	-	EXPRESSION TAG	UNP P0A2E1

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C<sub>2</sub>H<sub>3</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			4	2	2		

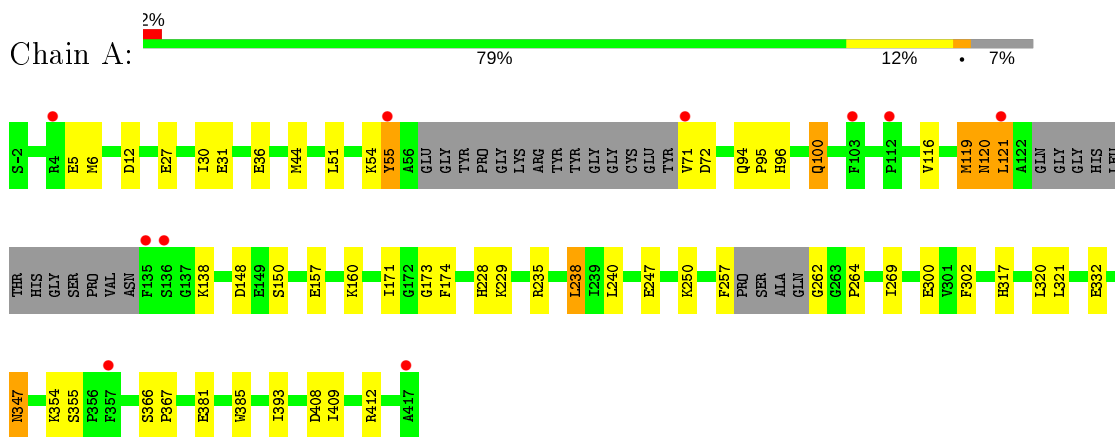
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	211	Total	O	0	3
			212	212		
3	B	238	Total	O	0	1
			238	238		

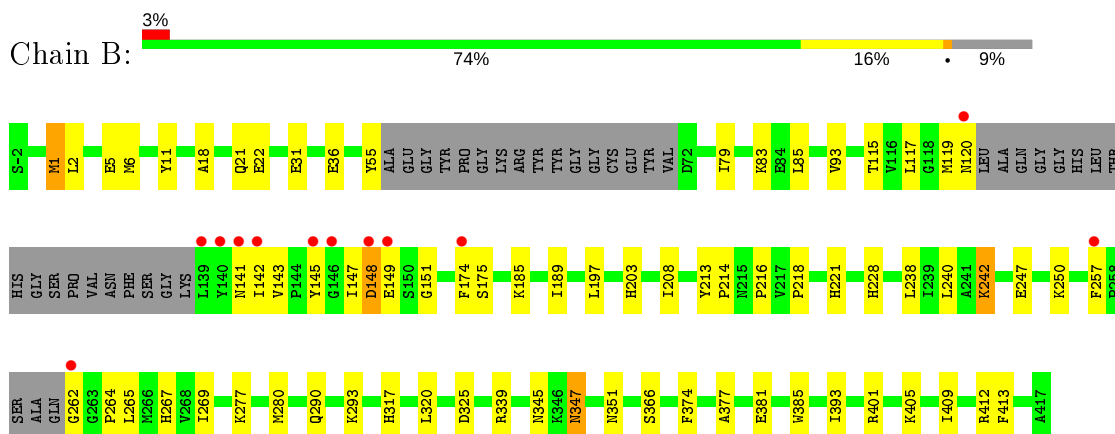
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Serine hydroxymethyltransferase



- Molecule 1: Serine hydroxymethyltransferase



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.95Å 49.73Å 94.74Å 90.00° 110.67° 90.00°	Depositor
Resolution (Å)	45.30 – 1.80 45.28 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.4 (45.30-1.80) 99.3 (45.28-1.80)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.68 (at 1.79Å)	Xtrriage
Refinement program	REFMAC 5.5.0054	Depositor
R, $R_{free}$	0.152 , 0.191 0.164 , 0.203	Depositor DCC
$R_{free}$ test set	3022 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.0	Xtrriage
Anisotropy	0.051	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 51.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6646	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.65% 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](#)

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

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	0.77	0/3242	0.81	3/4351 (0.1%)
1	B	0.83	0/3195	0.82	5/4289 (0.1%)
All	All	0.80	0/6437	0.81	8/8640 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	339	ARG	CG-CD-NE	-6.96	97.18	111.80
1	B	228[A]	HIS	N-CA-C	6.29	128.00	111.00
1	B	228[B]	HIS	N-CA-C	6.29	128.00	111.00
1	A	228[A]	HIS	N-CA-C	-5.58	95.93	111.00
1	A	228[B]	HIS	N-CA-C	-5.58	95.93	111.00
1	B	339	ARG	NE-CZ-NH1	-5.46	117.57	120.30
1	B	325	ASP	CB-CG-OD1	5.27	123.04	118.30
1	A	12	ASP	CB-CG-OD2	-5.18	113.64	118.30

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	3119	0	3137	63	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3073	0	3124	72	0
2	B	4	0	3	0	0
3	A	212	0	0	8	0
3	B	238	0	0	5	0
All	All	6646	0	6264	130	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 10.

All (130) 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:277:LYS:HA	1:B:280[A]:MSE:HE2	1.29	1.13
1:A:30:ILE:HD11	1:A:409[B]:ILE:HD11	1.35	1.03
1:A:31[A]:GLU:HG3	3:A:441:HOH:O	1.57	1.02
1:A:54:LYS:NZ	1:A:71:VAL:HG11	1.81	0.95
1:B:374:PHE:CZ	1:B:409[B]:ILE:HD11	2.02	0.94
1:A:54:LYS:HG2	1:A:71:VAL:HG12	1.52	0.92
1:A:119:MSE:C	1:A:120[A]:ASN:HD22	1.73	0.91
1:A:54:LYS:HZ3	1:A:71:VAL:HG11	1.33	0.89
1:A:30:ILE:CD1	1:A:409[B]:ILE:HD11	2.09	0.82
1:A:36[B]:GLU:CD	1:B:55:TYR:HB2	2.01	0.81
1:B:197:LEU:H	1:B:221:HIS:HD2	1.31	0.78
1:B:31[B]:GLU:HG3	3:B:479:HOH:O	1.87	0.74
1:A:121:LEU:HD13	1:A:174:PHE:HE2	1.54	0.72
1:B:147[B]:ILE:HD11	1:B:151:GLY:C	2.10	0.71
1:A:27[A]:GLU:HG2	3:A:703:HOH:O	1.90	0.71
1:A:235[B]:ARG:HH21	1:A:235[B]:ARG:HB3	1.55	0.70
1:A:30:ILE:HD11	1:A:409[B]:ILE:CD1	2.18	0.69
1:B:93:VAL:HG12	1:B:238[A]:LEU:HD21	1.74	0.69
1:A:320:LEU:HD21	1:A:347:ASN:HB3	1.74	0.69
1:A:54:LYS:HG2	1:A:71:VAL:CG1	2.24	0.68
1:B:119:MSE:HE3	1:B:147[A]:ILE:CG2	2.23	0.68
1:B:374:PHE:CE1	1:B:409[B]:ILE:HD11	2.31	0.65
1:A:55:TYR:HB2	1:B:36[A]:GLU:CD	2.17	0.65
1:B:197:LEU:H	1:B:221:HIS:CD2	2.16	0.64
1:A:121:LEU:HD22	3:A:507:HOH:O	1.97	0.64
1:B:385:TRP:HH2	1:B:409[B]:ILE:HG21	1.63	0.64
1:A:385:TRP:HH2	1:A:409[A]:ILE:HG13	1.63	0.63
1:A:119:MSE:C	1:A:120[A]:ASN:ND2	2.49	0.63
1:A:235[B]:ARG:NH2	1:A:235[B]:ARG:HB3	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:MSE:HE1	1:B:147[A]:ILE:HD13	1.81	0.62
1:B:401[B]:ARG:HG2	1:B:405[B]:LYS:HE3	1.82	0.62
1:A:121:LEU:HD13	1:A:174:PHE:CE2	2.34	0.61
1:B:175:SER:HA	1:B:203[B]:HIS:CD2	2.36	0.61
1:B:6[B]:MSE:HE3	1:B:11:TYR:HA	1.83	0.60
1:B:119:MSE:HE3	1:B:147[A]:ILE:HG23	1.83	0.59
1:A:157[B]:GLU:HA	1:A:157[B]:GLU:OE1	2.03	0.59
1:A:257:PHE:CG	1:A:262:GLY:O	2.56	0.59
1:B:385:TRP:HH2	1:B:409[A]:ILE:HG13	1.66	0.59
1:B:257:PHE:CD2	1:B:262:GLY:O	2.56	0.59
1:A:96:HIS:HB2	1:A:100:GLN:HE21	1.68	0.58
1:A:235[B]:ARG:HH21	1:A:235[B]:ARG:CB	2.17	0.58
1:A:71:VAL:CG2	1:A:72:ASP:N	2.66	0.58
1:A:71:VAL:HG23	1:A:72:ASP:N	2.17	0.58
1:A:385:TRP:CH2	1:A:409[A]:ILE:HG13	2.38	0.58
1:B:148:ASP:OD1	1:B:148:ASP:N	2.29	0.57
1:B:185:LYS:HE2	1:B:189:ILE:HD11	1.86	0.57
1:A:36[B]:GLU:OE2	1:B:55:TYR:HB2	2.04	0.57
1:A:317:HIS:CD2	1:A:317:HIS:H	2.23	0.56
1:A:317:HIS:HE1	3:A:522:HOH:O	1.88	0.56
1:A:385:TRP:HH2	1:A:409[B]:ILE:HG21	1.71	0.56
1:A:408[B]:ASP:OD2	1:A:412:ARG:NH1	2.37	0.56
1:A:5[A]:GLU:HG2	1:A:6:MSE:HG3	1.88	0.54
1:A:36[B]:GLU:OE1	3:A:455:HOH:O	2.18	0.53
1:B:257:PHE:CG	1:B:262:GLY:O	2.61	0.53
1:B:119:MSE:HB2	1:B:145:TYR:CZ	2.45	0.52
1:A:235[B]:ARG:HD3	1:B:265:LEU:HD23	1.92	0.52
1:B:85:LEU:HD13	1:B:280[B]:MSE:HG2	1.92	0.52
1:B:115:THR:HA	1:B:141:ASN:O	2.11	0.51
1:B:117:LEU:HD12	1:B:143:VAL:O	2.10	0.51
1:B:267:HIS:H	1:B:267:HIS:CD2	2.29	0.51
1:A:51:LEU:HD23	1:B:22:GLU:HG2	1.93	0.50
1:B:175:SER:HB3	1:B:203[B]:HIS:NE2	2.26	0.50
1:B:385:TRP:CH2	1:B:409[B]:ILE:HG21	2.47	0.49
1:A:385:TRP:HH2	1:A:409[A]:ILE:CG1	2.26	0.49
1:B:119:MSE:HE2	1:B:174:PHE:CD2	2.48	0.49
1:A:54:LYS:HZ2	1:A:71:VAL:HG11	1.72	0.48
1:A:238[A]:LEU:HD22	1:A:240:LEU:HG	1.94	0.48
1:B:393[B]:ILE:HG23	3:B:464:HOH:O	2.12	0.48
1:A:257:PHE:CD2	1:A:262:GLY:O	2.67	0.48
1:A:44:MSE:SE	3:B:535:HOH:O	2.81	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203[B]:HIS:ND1	3:B:598:HOH:O	2.22	0.48
1:A:302:PHE:CE1	1:A:321:LEU:HD11	2.49	0.48
1:B:293[B]:LYS:HB2	1:B:293[B]:LYS:HE2	1.55	0.48
1:B:320:LEU:HD11	1:B:347:ASN:HB3	1.96	0.47
1:B:149[A]:GLU:CG	1:B:351:ASN:HD22	2.27	0.47
1:B:290:GLN:HA	1:B:293[B]:LYS:HE2	1.97	0.47
1:B:147[B]:ILE:HD11	1:B:151:GLY:O	2.13	0.47
1:B:119:MSE:CE	1:B:147[A]:ILE:HD13	2.43	0.47
1:A:54:LYS:NZ	1:A:71:VAL:CG1	2.67	0.47
1:A:54:LYS:CG	1:A:71:VAL:CG1	2.93	0.46
1:A:366:SER:N	1:A:367:PRO:CD	2.79	0.46
3:A:673:HOH:O	1:B:1[A]:MSE:SE	2.84	0.46
1:B:385:TRP:CH2	1:B:409[A]:ILE:HG13	2.49	0.46
1:A:300[A]:GLU:HG3	3:A:561:HOH:O	2.15	0.46
1:B:218:PRO:HA	1:B:242:LYS:HE3	1.98	0.46
1:B:409[B]:ILE:HG22	3:B:441:HOH:O	2.16	0.46
1:A:54:LYS:CG	1:A:71:VAL:HG12	2.35	0.45
1:B:409[A]:ILE:HD13	1:B:412[A]:ARG:HE	1.81	0.45
1:A:409[B]:ILE:HG23	3:A:514:HOH:O	2.16	0.45
1:B:119:MSE:HB2	1:B:145:TYR:CE1	2.52	0.45
1:B:149[A]:GLU:CD	1:B:351:ASN:HD22	2.20	0.45
1:A:119:MSE:HE2	1:A:173:GLY:H	1.82	0.45
1:B:149[A]:GLU:CG	1:B:351:ASN:ND2	2.79	0.45
1:A:116:VAL:HG11	1:A:171:ILE:HD11	1.99	0.44
1:A:366:SER:N	1:A:367:PRO:HD3	2.33	0.44
1:B:185:LYS:CE	1:B:189:ILE:HD11	2.47	0.44
1:B:18:ALA:HA	1:B:21[A]:GLN:HE21	1.82	0.44
1:A:96:HIS:HB2	1:A:100:GLN:NE2	2.30	0.44
1:B:409[B]:ILE:CD1	1:B:413:PHE:CD2	3.00	0.44
1:B:409[B]:ILE:CD1	1:B:413:PHE:HD2	2.30	0.43
1:A:71:VAL:CG2	1:A:72:ASP:H	2.31	0.43
1:A:332[B]:GLU:HG2	1:A:393:ILE:HD13	2.01	0.43
1:B:119:MSE:HE3	1:B:147[A]:ILE:HG21	2.00	0.43
1:B:377:ALA:O	1:B:381[B]:GLU:HG3	2.18	0.43
1:B:79:ILE:HG22	1:B:83:LYS:HE2	2.01	0.43
1:A:409[A]:ILE:HD13	1:A:412:ARG:HH22	1.85	0.42
1:A:116:VAL:HG11	1:A:171:ILE:CD1	2.50	0.42
1:A:148[A]:ASP:HB2	1:A:150:SER:OG	2.20	0.42
1:B:115:THR:HG23	1:B:141:ASN:HD22	1.85	0.41
1:B:409[B]:ILE:HD13	1:B:413:PHE:CD2	2.55	0.41
1:A:94:GLN:N	1:A:95:PRO:CD	2.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1[B]:MSE:O	1:B:2:LEU:HD23	2.20	0.41
1:B:119:MSE:CE	1:B:147[A]:ILE:HG21	2.51	0.41
1:B:238[B]:LEU:HD23	1:B:240:LEU:HD11	2.03	0.41
1:A:264:PRO:HB2	1:A:269:ILE:CD1	2.51	0.41
1:A:381[B]:GLU:HG2	1:A:385:TRP:CZ2	2.56	0.40
1:B:247:GLU:HG2	1:B:250[A]:LYS:NZ	2.36	0.40
1:B:264:PRO:HB2	1:B:269:ILE:HD12	2.03	0.40
1:B:213:TYR:CD2	1:B:214:PRO:HD2	2.57	0.40
1:B:208:ILE:HD13	1:B:216:PRO:HD3	2.03	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	406/420 (97%)	391 (96%)	13 (3%)	2 (0%)	29	15
1	B	399/420 (95%)	381 (96%)	17 (4%)	1 (0%)	41	27
All	All	805/840 (96%)	772 (96%)	30 (4%)	3 (0%)	41	21

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	229[A]	LYS
1	A	229[B]	LYS
1	B	366	SER

#### 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	336/321 (105%)	322 (96%)	14 (4%)	30	15
1	B	332/321 (103%)	323 (97%)	9 (3%)	44	31
All	All	668/642 (104%)	645 (97%)	23 (3%)	39	22

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	TYR
1	A	100	GLN
1	A	119	MSE
1	A	120[A]	ASN
1	A	120[B]	ASN
1	A	121	LEU
1	A	138	LYS
1	A	160	LYS
1	A	238[A]	LEU
1	A	238[B]	LEU
1	A	247	GLU
1	A	347	ASN
1	A	354	LYS
1	A	355	SER
1	B	1[A]	MSE
1	B	1[B]	MSE
1	B	5	GLU
1	B	120	ASN
1	B	142	ILE
1	B	148	ASP
1	B	242	LYS
1	B	345	ASN
1	B	347	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	GLN
1	A	317	HIS
1	A	345	ASN
1	A	351	ASN

*Continued on next page...*

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Mol	Chain	Res	Type
1	B	96	HIS
1	B	120	ASN
1	B	141	ASN
1	B	221	HIS
1	B	267	HIS
1	B	351	ASN

### 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](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	ACT	B	501	-	1,3,3	1.46	0	0,3,3	0.00	-

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	377/420 (89%)	-0.37	10 (2%) 54 49	9, 15, 27, 38	0
1	B	370/420 (88%)	-0.43	12 (3%) 47 41	8, 14, 27, 38	0
All	All	747/840 (88%)	-0.40	22 (2%) 51 46	8, 15, 27, 38	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	139	LEU	4.5
1	B	174	PHE	3.5
1	A	71	VAL	3.2
1	A	136	SER	3.2
1	B	141	ASN	3.1
1	B	148	ASP	3.0
1	B	120	ASN	2.9
1	A	112	PRO	2.9
1	A	103[A]	PHE	2.9
1	A	55	TYR	2.8
1	A	135	PHE	2.8
1	B	146	GLY	2.7
1	B	142	ILE	2.7
1	A	417	ALA	2.7
1	A	357	PHE	2.5
1	B	262	GLY	2.3
1	B	145	TYR	2.1
1	A	121	LEU	2.1
1	B	140	TYR	2.1
1	B	257	PHE	2.1
1	A	4	ARG	2.0
1	B	149[A]	GLU	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 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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	ACT	B	501	4/4	0.95	0.19	42,42,43,44	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.