



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 11:58 pm BST

PDB ID : 5JIS  
Title : The Crystal Structure of O-acetyl serine sulphydralase from Brucella abortus  
Authors : Dharavath, S.; Gourinath, S.  
Deposited on : 2016-04-22  
Resolution : 2.20 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

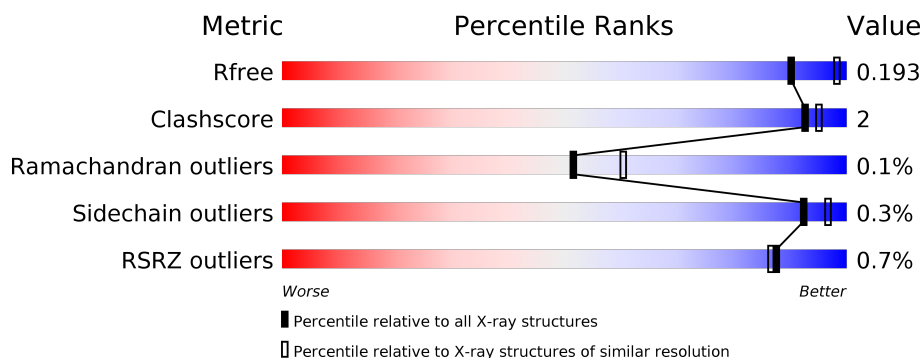
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (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 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	350	<div> <div style="width: 92%;"></div> <div style="width: 5%;"></div> <div style="width: 3%;"></div> </div> <div>92% 5%</div>
1	B	350	<div> <div style="width: 92%;"></div> <div style="width: 5%;"></div> <div style="width: 3%;"></div> </div> <div>92% 5%</div>
1	C	350	<div> <div style="width: 91%;"></div> <div style="width: 6%;"></div> <div style="width: 3%;"></div> </div> <div>91% 6%</div>
1	D	350	<div> <div style="width: 93%;"></div> <div style="width: 5%;"></div> <div style="width: 2%;"></div> </div> <div>93% 5%</div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10645 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 Cysteine synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	P	S	0	0	0
			2513	1581	441	485	1	5			
1	B	331	Total	C	N	O	P	S	0	1	0
			2519	1586	442	485	1	5			
1	C	339	Total	C	N	O	P	S	0	1	0
			2584	1628	450	500	1	5			
1	D	339	Total	C	N	O	P	S	0	1	0
			2586	1628	450	502	1	5			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	343	LEU	-	expression tag	UNP A0A0F6AQU1
A	344	GLU	-	expression tag	UNP A0A0F6AQU1
A	345	HIS	-	expression tag	UNP A0A0F6AQU1
A	346	HIS	-	expression tag	UNP A0A0F6AQU1
A	347	HIS	-	expression tag	UNP A0A0F6AQU1
A	348	HIS	-	expression tag	UNP A0A0F6AQU1
A	349	HIS	-	expression tag	UNP A0A0F6AQU1
A	350	HIS	-	expression tag	UNP A0A0F6AQU1
B	343	LEU	-	expression tag	UNP A0A0F6AQU1
B	344	GLU	-	expression tag	UNP A0A0F6AQU1
B	345	HIS	-	expression tag	UNP A0A0F6AQU1
B	346	HIS	-	expression tag	UNP A0A0F6AQU1
B	347	HIS	-	expression tag	UNP A0A0F6AQU1
B	348	HIS	-	expression tag	UNP A0A0F6AQU1
B	349	HIS	-	expression tag	UNP A0A0F6AQU1
B	350	HIS	-	expression tag	UNP A0A0F6AQU1
C	343	LEU	-	expression tag	UNP A0A0F6AQU1
C	344	GLU	-	expression tag	UNP A0A0F6AQU1
C	345	HIS	-	expression tag	UNP A0A0F6AQU1
C	346	HIS	-	expression tag	UNP A0A0F6AQU1
C	347	HIS	-	expression tag	UNP A0A0F6AQU1

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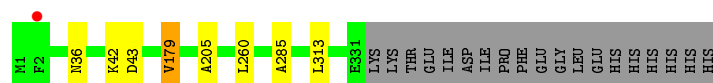
Chain	Residue	Modelled	Actual	Comment	Reference
C	348	HIS	-	expression tag	UNP A0A0F6AQU1
C	349	HIS	-	expression tag	UNP A0A0F6AQU1
C	350	HIS	-	expression tag	UNP A0A0F6AQU1
D	343	LEU	-	expression tag	UNP A0A0F6AQU1
D	344	GLU	-	expression tag	UNP A0A0F6AQU1
D	345	HIS	-	expression tag	UNP A0A0F6AQU1
D	346	HIS	-	expression tag	UNP A0A0F6AQU1
D	347	HIS	-	expression tag	UNP A0A0F6AQU1
D	348	HIS	-	expression tag	UNP A0A0F6AQU1
D	349	HIS	-	expression tag	UNP A0A0F6AQU1
D	350	HIS	-	expression tag	UNP A0A0F6AQU1

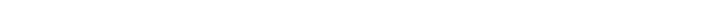
- Molecule 2 is water.

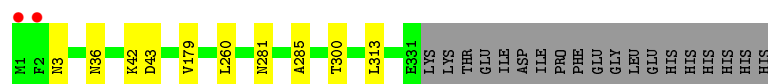
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	125	Total O 125 125	0	0
2	B	126	Total O 126 126	0	0
2	C	98	Total O 98 98	0	0
2	D	94	Total O 94 94	0	0



- Molecule 1: Cysteine synthase



- Chain B:  92% 5%



- Chain C:  91% 6%



- Chain D:  93%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.79Å 104.79Å 241.51Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.07 – 2.20 48.07 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.1 (40.07-2.20) 98.1 (48.07-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.67 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, $R_{free}$	0.184 , 0.226 0.194 , 0.193	Depositor DCC
$R_{free}$ test set	3737 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	24.1	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 16.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.480 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10645	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% 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: LLP

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	0.45	0/2532	0.67	1/3435 (0.0%)
1	B	0.45	0/2541	0.68	0/3446
1	C	0.45	0/2607	0.69	0/3536
1	D	0.45	0/2606	0.68	1/3535 (0.0%)
All	All	0.45	0/10286	0.68	2/13952 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	179	VAL	CB-CA-C	-6.26	99.50	111.40
1	D	308	ARG	NE-CZ-NH2	-5.08	117.76	120.30

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	2513	0	2508	5	0
1	B	2519	0	2521	5	0
1	C	2584	0	2571	14	0
1	D	2586	0	2567	7	0
2	A	125	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	126	0	0	0	0
2	C	98	0	0	1	0
2	D	94	0	0	0	0
All	All	10645	0	10167	31	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 2.

All (31) 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:C:72:ASN:OD1	1:C:308:ARG:NH1	1.95	0.99
1:C:25:GLY:O	1:C:294:PRO:O	2.09	0.70
1:C:114:VAL:HG22	1:C:115:PRO:HD2	1.85	0.58
1:C:59:LEU:HD11	1:C:65:ILE:CD1	2.35	0.57
1:C:96:GLN:O	1:C:101:LYS:NZ	2.38	0.57
1:A:179:VAL:HG21	1:A:285:ALA:HA	1.90	0.53
1:D:59:LEU:HD11	1:D:65:ILE:CD1	2.41	0.51
1:C:42:LLP:H2'1	1:C:72:ASN:ND2	2.29	0.48
1:A:36:ASN:HB3	1:A:43:ASP:OD2	2.14	0.48
1:A:42:LLP:OP2	1:A:42:LLP:H4'1	2.14	0.47
1:C:42:LLP:H4'1	1:C:42:LLP:OP2	2.15	0.47
1:B:179:VAL:HG11	1:B:285:ALA:HA	1.96	0.46
1:D:42:LLP:H2'1	1:D:72:ASN:ND2	2.30	0.46
1:C:179:VAL:HG22	1:C:288:LEU:HD22	1.98	0.46
1:B:42:LLP:H4'1	1:B:42:LLP:OP2	2.16	0.45
1:A:179:VAL:HG13	1:A:205:ALA:HB3	1.99	0.44
1:D:112:ILE:HD13	1:D:132:LEU:HD21	1.99	0.44
1:D:36:ASN:HB3	1:D:43:ASP:OD2	2.17	0.44
1:B:260:LEU:HD22	1:B:313:LEU:HD11	2.01	0.43
1:C:36:ASN:HB3	1:C:43:ASP:OD2	2.19	0.43
1:B:281:ASN:HB3	1:B:300:THR:OG1	2.18	0.43
1:C:114:VAL:HG22	1:C:115:PRO:CD	2.49	0.43
1:D:179:VAL:HG13	1:D:205:ALA:HB3	2.01	0.43
1:C:56:ARG:NH1	2:C:402:HOH:O	2.51	0.42
1:C:59:LEU:HD11	1:C:65:ILE:HD11	2.00	0.42
1:A:260:LEU:HD22	1:A:313:LEU:HD11	2.03	0.41
1:D:42:LLP:OP2	1:D:42:LLP:H4'1	2.20	0.41
1:D:179:VAL:HG22	1:D:288:LEU:HD22	2.02	0.41
1:C:2:PHE:CZ	1:C:51:ARG:HD2	2.56	0.40
1:C:292:LEU:HD12	1:C:298:ILE:HD11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:ASN:HB3	1:B:43:ASP:OD2	2.20	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	328/350 (94%)	323 (98%)	5 (2%)	0	100	100
1	B	329/350 (94%)	323 (98%)	6 (2%)	0	100	100
1	C	335/350 (96%)	330 (98%)	5 (2%)	0	100	100
1	D	335/350 (96%)	326 (97%)	8 (2%)	1 (0%)	41	46
All	All	1327/1400 (95%)	1302 (98%)	24 (2%)	1 (0%)	51	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	71	GLY

### 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	256/274 (93%)	256 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	257/274 (94%)	256 (100%)	1 (0%)	91	96
1	C	264/274 (96%)	264 (100%)	0	100	100
1	D	264/274 (96%)	262 (99%)	2 (1%)	81	90
All	All	1041/1096 (95%)	1038 (100%)	3 (0%)	92	97

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	3	ASN
1	D	60	ARG
1	D	228	ASP

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

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are 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
1	LLP	B	42	1	23,24,25	2.38	4 (17%)	25,32,34	1.71	4 (16%)
1	LLP	C	42	1	23,24,25	2.51	4 (17%)	25,32,34	1.58	4 (16%)
1	LLP	A	42	1	23,24,25	2.43	4 (17%)	25,32,34	1.44	3 (12%)
1	LLP	D	42	1	23,24,25	2.52	4 (17%)	25,32,34	1.53	3 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	LLP	B	42	1	-	1/16/17/19	0/1/1/1
1	LLP	C	42	1	-	2/16/17/19	0/1/1/1
1	LLP	A	42	1	-	2/16/17/19	0/1/1/1
1	LLP	D	42	1	-	2/16/17/19	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	42	LLP	C3-C2	8.02	1.48	1.40
1	D	42	LLP	C3-C2	7.76	1.48	1.40
1	A	42	LLP	C3-C2	7.74	1.48	1.40
1	B	42	LLP	C3-C2	7.46	1.48	1.40
1	D	42	LLP	C4-C5	5.49	1.48	1.42
1	C	42	LLP	C4-C3	5.21	1.48	1.40
1	B	42	LLP	C4-C5	4.82	1.48	1.42
1	A	42	LLP	C4'-NZ	4.82	1.43	1.27
1	D	42	LLP	C4'-NZ	4.79	1.43	1.27
1	D	42	LLP	C4-C3	4.77	1.48	1.40
1	C	42	LLP	C4'-NZ	4.72	1.43	1.27
1	C	42	LLP	C4-C5	4.70	1.47	1.42
1	A	42	LLP	C4-C5	4.68	1.47	1.42
1	B	42	LLP	C4'-NZ	4.61	1.42	1.27
1	B	42	LLP	C4-C3	4.59	1.47	1.40
1	A	42	LLP	C4-C3	4.53	1.47	1.40

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	42	LLP	C4-C3-C2	-5.83	116.58	120.19
1	D	42	LLP	C4-C3-C2	-5.23	116.95	120.19
1	C	42	LLP	C4-C3-C2	-4.96	117.12	120.19
1	A	42	LLP	C4-C3-C2	-4.22	117.58	120.19
1	B	42	LLP	C6-N1-C2	2.61	124.00	119.17
1	C	42	LLP	OP4-C5'-C5	2.55	114.20	109.35
1	C	42	LLP	C6-N1-C2	2.54	123.87	119.17
1	B	42	LLP	O3-C3-C2	2.53	123.01	117.49
1	A	42	LLP	C6-N1-C2	2.33	123.47	119.17
1	D	42	LLP	O3-C3-C2	2.28	122.47	117.49
1	D	42	LLP	C6-N1-C2	2.28	123.38	119.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	42	LLP	OP4-C5'-C5	2.20	113.55	109.35
1	A	42	LLP	O3-C3-C2	2.12	122.11	117.49
1	C	42	LLP	O3-C3-C2	2.10	122.06	117.49

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	42	LLP	C4-C4'-NZ-CE
1	D	42	LLP	C4-C4'-NZ-CE
1	B	42	LLP	C4-C4'-NZ-CE
1	C	42	LLP	C4-C4'-NZ-CE
1	D	42	LLP	CE-CD-CG-CB
1	A	42	LLP	CE-CD-CG-CB
1	C	42	LLP	CE-CD-CG-CB

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	42	LLP	1	0
1	C	42	LLP	2	0
1	A	42	LLP	1	0
1	D	42	LLP	2	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates 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 ⓘ

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	330/350 (94%)	-0.68	1 (0%) 94 93	15, 23, 38, 63	0
1	B	330/350 (94%)	-0.69	2 (0%) 89 88	16, 22, 36, 71	0
1	C	338/350 (96%)	-0.47	4 (1%) 79 77	16, 24, 52, 59	0
1	D	338/350 (96%)	-0.47	3 (0%) 84 83	15, 24, 51, 64	0
All	All	1336/1400 (95%)	-0.58	10 (0%) 87 86	15, 23, 47, 71	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	MET	4.5
1	C	114	VAL	4.2
1	D	70	ALA	3.6
1	A	2	PHE	3.3
1	D	117	ALA	2.4
1	C	92	ILE	2.3
1	C	112	ILE	2.2
1	C	115	PRO	2.1
1	B	2	PHE	2.0
1	D	97	SER	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [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
1	LLP	C	42	24/25	0.97	0.10	16,17,18,18	0
1	LLP	D	42	24/25	0.97	0.09	16,17,17,18	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
1	LLP	A	42	24/25	0.98	0.09	15,16,17,17	0
1	LLP	B	42	24/25	0.98	0.10	15,16,17,17	0

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.