



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:16 PM GMT

PDB ID : 4X8E  
Title : Ergothioneine-biosynthetic sulfoxide synthase EgtB in complex with N,N,N-trimethyl-histidine  
Authors : Vit, A.; Goncharenko, K.V.; Blankenfeldt, W.; Seebeck, F.P.  
Deposited on : 2014-12-10  
Resolution : 1.60 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.

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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

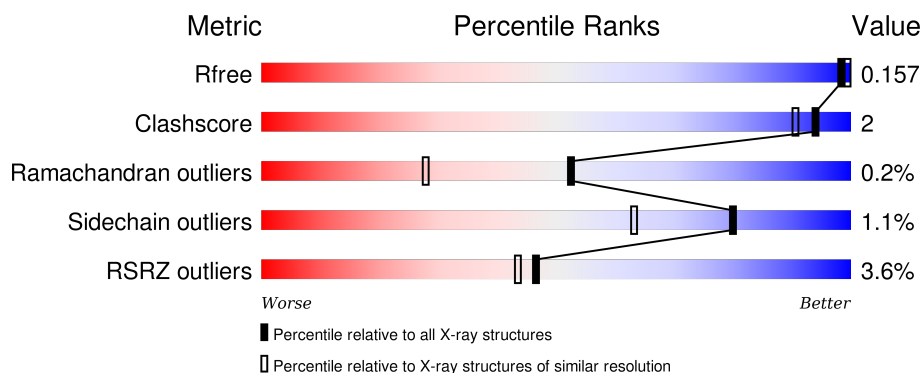
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.60 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	2475 (1.60-1.60)
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)
RSRZ outliers	91569	2479 (1.60-1.60)

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	447	<div> <div></div> <div>90%</div> <div>6%</div> <div>.</div> </div>
1	B	447	<div> <div>6%</div> <div>92%</div> <div>.</div> <div>5%</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
3	GOL	A	502	-	-	-	X
3	GOL	A	504	-	-	-	X
3	GOL	A	505	-	-	-	X
3	GOL	A	506	-	-	-	X
3	GOL	A	507	-	-	-	X
3	GOL	A	508	-	-	-	X
3	GOL	A	509	-	-	-	X
3	GOL	B	502	-	-	-	X
3	GOL	B	503	-	-	-	X
3	GOL	B	504	-	-	-	X
3	GOL	B	505	-	-	-	X
4	CL	A	512	-	-	X	-

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 14685 atoms, of which 6657 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.

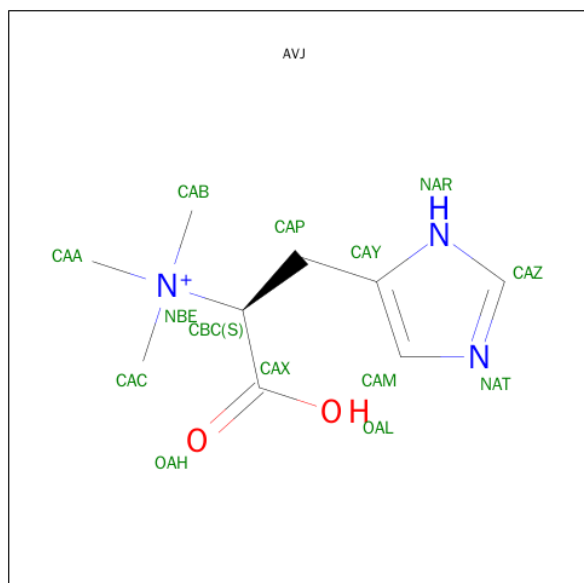
- Molecule 1 is a protein called Sulfoxide synthase EgtB.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	428	Total	C	H	N	O	S	0	14	0
			6759	2188	3314	635	612	10			
1	B	426	Total	C	H	N	O	S	0	3	0
			6583	2138	3219	615	601	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP G7CFI3
A	1	HIS	-	expression tag	UNP G7CFI3
A	2	MET	-	expression tag	UNP G7CFI3
B	0	GLY	-	expression tag	UNP G7CFI3
B	1	HIS	-	expression tag	UNP G7CFI3
B	2	MET	-	expression tag	UNP G7CFI3

- Molecule 2 is N,N,N-trimethyl-histidine (three-letter code: AVJ) (formula: C<sub>9</sub>H<sub>16</sub>N<sub>3</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	0	0
			29	9	15	3	2		
2	B	1	Total	C	H	N	O	0	0
			29	9	15	3	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			13	3	7	3		
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			13	3	7	3		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Cl	0	0
			2	2		
4	A	5	Total	Cl	0	0
			5	5		

- Molecule 5 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Fe	0	0
			1	1		
5	A	1	Total	Fe	0	0
			1	1		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 7 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Mg	0	0
			1	1		
7	A	1	Total	Mg	0	0
			1	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	671	Total	O	0	0
			671	671		

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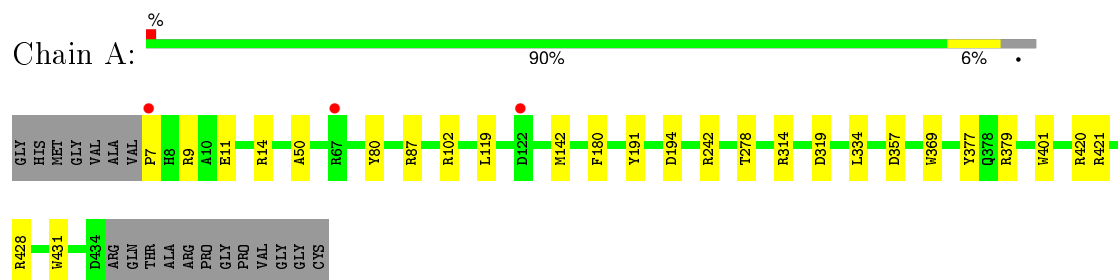
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	435	Total 435	O 435	0	0

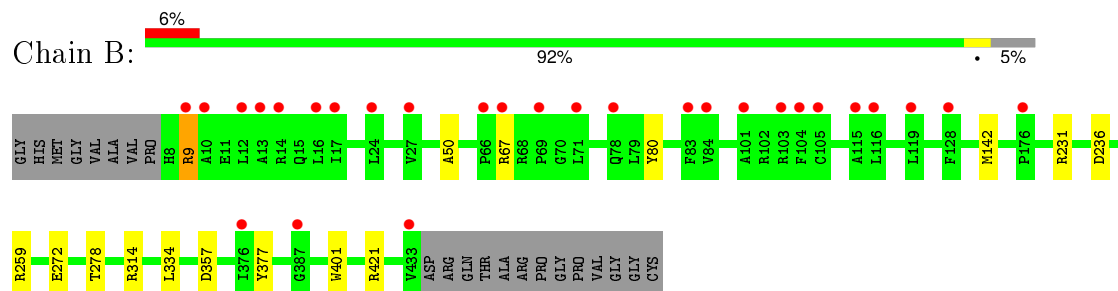
### 3 Residue-property plots [i](#)

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: Sulfoxide synthase EgtB



- Molecule 1: Sulfoxide synthase EgtB





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	135.24Å 135.24Å 142.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.97 – 1.60 48.97 – 1.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.97-1.60) 100.0 (48.97-1.60)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.34 (at 1.60Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.144 , 0.163 0.159 , 0.157	Depositor DCC
$R_{free}$ test set	8755 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	21.5	Xtriage
Anisotropy	0.133	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 51.5	EDS
Estimated twinning fraction	0.013 for -h,l,k 0.009 for -l,-k,-h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 172440 reflections (0.001%)	Xtriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	14685	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, MG, CL, CA, FE, AVJ

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	0.55	0/3602	0.74	9/4924 (0.2%)
1	B	0.44	0/3482	0.59	3/4764 (0.1%)
All	All	0.50	0/7084	0.67	12/9688 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	142	MET	CG-SD-CE	-8.33	86.87	100.20
1	B	9	ARG	NE-CZ-NH1	5.93	123.27	120.30
1	A	420	ARG	NE-CZ-NH2	-5.59	117.51	120.30
1	A	319[A]	ASP	CB-CG-OD1	5.36	123.13	118.30
1	A	319[B]	ASP	CB-CG-OD1	5.36	123.13	118.30
1	A	428	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	A	194	ASP	CB-CG-OD1	5.09	122.88	118.30
1	A	428	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	B	67	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	B	142	MET	CG-SD-CE	-5.06	92.11	100.20
1	A	87	ARG	NE-CZ-NH1	5.00	122.80	120.30
1	A	194	ASP	CB-CG-OD2	-5.00	113.80	118.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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	3445	3314	3316	11	1
1	B	3364	3219	3220	9	0
2	A	14	15	15	0	0
2	B	14	15	15	0	0
3	A	48	63	63	4	1
3	B	24	31	32	3	0
4	A	5	0	0	2	0
4	B	2	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
8	A	671	0	0	4	1
8	B	435	0	0	3	0
All	All	8028	6657	6661	24	2

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 (24) 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:272[B]:GLU:OE2	8:B:917:HOH:O	2.05	0.73
1:A:119:LEU:O	8:A:1019:HOH:O	2.14	0.65
1:A:379[A]:ARG:NH2	3:A:506:GOL:O2	2.30	0.65
3:A:508:GOL:O3	8:A:754:HOH:O	2.15	0.64
1:B:259:ARG:NH1	1:B:272[B]:GLU:OE2	2.33	0.61
1:B:231:ARG:NH2	1:B:236:ASP:OD1	2.34	0.60
1:B:314:ARG:NH2	8:B:929:HOH:O	2.36	0.57
1:A:7:PRO:HD3	1:A:191:TYR:OH	2.06	0.56
3:A:507:GOL:C3	4:A:512:CL:CL	2.93	0.54
1:A:7:PRO:CD	1:A:191:TYR:OH	2.55	0.54
1:A:314:ARG:NH2	8:A:1121:HOH:O	2.43	0.52
1:B:421:ARG:HE	3:B:505:GOL:H31	1.75	0.51
3:B:505:GOL:H32	8:B:628:HOH:O	2.10	0.51
1:B:334:LEU:HA	1:B:401:TRP:CZ3	2.46	0.50
1:B:421:ARG:NE	3:B:505:GOL:H31	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:ALA:HB3	1:B:80:TYR:CE1	2.49	0.48
3:A:507:GOL:H31	4:A:512:CL:CL	2.51	0.47
1:B:278:THR:HG23	1:B:421:ARG:HD3	1.98	0.46
1:A:11:GLU:OE2	1:A:14[A]:ARG:NH2	2.50	0.44
1:A:278:THR:HG23	1:A:421:ARG:HD3	1.99	0.44
1:A:50:ALA:HB3	1:A:80:TYR:CE1	2.53	0.43
1:A:102[A]:ARG:HD3	8:A:1221:HOH:O	2.18	0.42
1:A:334:LEU:HA	1:A:401:TRP:CZ3	2.55	0.41
1:A:7:PRO:HD3	1:A:191:TYR:HH	1.86	0.41

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:431:TRP:HE1	3:A:503:GOL:HO2[8_665]	1.12	0.48
8:A:688:HOH:O	8:A:688:HOH:O[8_665]	2.14	0.06

## 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	441/447 (99%)	436 (99%)	4 (1%)	1 (0%)	52	28
1	B	427/447 (96%)	423 (99%)	3 (1%)	1 (0%)	52	28
All	All	868/894 (97%)	859 (99%)	7 (1%)	2 (0%)	52	28

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	357	ASP
1	B	357	ASP

### 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	347/348 (100%)	342 (99%)	5 (1%)	74	53
1	B	334/348 (96%)	332 (99%)	2 (1%)	90	82
All	All	681/696 (98%)	674 (99%)	7 (1%)	80	67

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ARG
1	A	180	PHE
1	A	242	ARG
1	A	369	TRP
1	A	377	TYR
1	B	9	ARG
1	B	377	TYR

Some 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 carbohydrates in this entry.

## 5.6 Ligand geometry

Of 27 ligands modelled in this entry, 13 are monoatomic - leaving 14 for Mogul analysis.

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	AVJ	A	501	5	6,14,14	1.54	1 (16%)	8,20,20	1.26	1 (12%)
3	GOL	A	502	-	5,5,5	0.38	0	5,5,5	0.37	0
3	GOL	A	503	-	5,5,5	0.35	0	5,5,5	0.15	0
3	GOL	A	504	-	5,5,5	0.34	0	5,5,5	0.30	0
3	GOL	A	505	-	5,5,5	0.25	0	5,5,5	0.72	0
3	GOL	A	506	-	5,5,5	0.53	0	5,5,5	0.32	0
3	GOL	A	507	-	5,5,5	0.47	0	5,5,5	0.54	0
3	GOL	A	508	-	5,5,5	1.02	0	5,5,5	1.17	1 (20%)
3	GOL	A	509	-	5,5,5	0.31	0	5,5,5	0.65	0
2	AVJ	B	501	5	6,14,14	2.13	2 (33%)	8,20,20	0.89	0
3	GOL	B	502	-	5,5,5	0.34	0	5,5,5	0.30	0
3	GOL	B	503	-	5,5,5	0.36	0	5,5,5	0.44	0
3	GOL	B	504	-	5,5,5	0.38	0	5,5,5	0.25	0
3	GOL	B	505	-	5,5,5	0.81	0	5,5,5	2.27	3 (60%)

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
2	AVJ	A	501	5	-	0/9/14/14	0/1/1/1
3	GOL	A	502	-	-	0/4/4/4	0/0/0/0
3	GOL	A	503	-	-	0/4/4/4	0/0/0/0
3	GOL	A	504	-	-	0/4/4/4	0/0/0/0
3	GOL	A	505	-	-	0/4/4/4	0/0/0/0
3	GOL	A	506	-	-	0/4/4/4	0/0/0/0
3	GOL	A	507	-	-	0/4/4/4	0/0/0/0
3	GOL	A	508	-	-	0/4/4/4	0/0/0/0
3	GOL	A	509	-	-	0/4/4/4	0/0/0/0
2	AVJ	B	501	5	-	0/9/14/14	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	502	-	-	0/4/4/4	0/0/0/0
3	GOL	B	503	-	-	0/4/4/4	0/0/0/0
3	GOL	B	504	-	-	0/4/4/4	0/0/0/0
3	GOL	B	505	-	-	0/4/4/4	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	AVJ	CBC-NBE	-3.80	1.46	1.54
2	A	501	AVJ	CBC-NBE	-2.56	1.49	1.54
2	B	501	AVJ	CAC-NBE	-2.10	1.44	1.50

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	505	GOL	O2-C2-C3	-3.37	93.21	108.65
3	B	505	GOL	O3-C3-C2	-2.73	96.93	110.18
3	A	508	GOL	C3-C2-C1	-2.25	102.29	111.12
3	B	505	GOL	O2-C2-C1	2.29	119.14	108.65
2	A	501	AVJ	CAM-NAT-CAZ	2.91	110.22	105.68

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	GOL	0	1
3	A	506	GOL	1	0
3	A	507	GOL	2	0
3	A	508	GOL	1	0
3	B	505	GOL	3	0

## 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	428/447 (95%)	-0.20	3 (0%) 89 89	14, 22, 40, 63	0
1	B	426/447 (95%)	0.22	28 (6%) 22 19	22, 36, 57, 72	0
All	All	854/894 (95%)	0.01	31 (3%) 46 43	14, 28, 53, 72	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	9	ARG	4.9
1	B	12	LEU	4.6
1	B	69	PRO	4.2
1	B	17	ILE	4.0
1	A	67	ARG	3.9
1	B	10	ALA	3.8
1	B	119	LEU	3.7
1	B	67	ARG	3.4
1	B	116	LEU	3.2
1	B	115	ALA	3.0
1	B	433	VAL	2.8
1	B	128	PHE	2.8
1	B	24	LEU	2.8
1	B	14	ARG	2.7
1	B	13	ALA	2.7
1	B	387	GLY	2.7
1	B	104	PHE	2.6
1	A	7	PRO	2.6
1	B	27	VAL	2.5
1	B	84	VAL	2.5
1	B	103	ARG	2.5
1	B	105	CYS	2.4
1	B	16	LEU	2.3
1	B	101	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	78	GLN	2.2
1	B	71	LEU	2.2
1	B	176	PRO	2.2
1	B	376	ILE	2.2
1	A	122	ASP	2.1
1	B	66	PRO	2.1
1	B	83	PHE	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. 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
3	GOL	B	505	6/6	0.91	0.14	11.75	26,57,71,74	0
3	GOL	A	509	6/6	0.73	0.38	11.27	37,58,70,70	0
3	GOL	A	505	6/6	0.85	0.18	8.26	22,31,56,56	0
3	GOL	A	508	6/6	0.85	0.17	7.91	35,46,66,77	0
3	GOL	A	504	6/6	0.97	0.13	7.27	29,37,43,44	0
3	GOL	A	506	6/6	0.90	0.22	6.76	34,51,61,72	0
3	GOL	A	507	6/6	0.84	0.23	5.18	32,55,63,66	0
3	GOL	B	504	6/6	0.92	0.22	4.75	32,55,70,70	0
3	GOL	B	503	6/6	0.85	0.15	4.63	48,58,68,71	0
3	GOL	A	502	6/6	0.94	0.13	2.80	21,43,56,56	0
3	GOL	B	502	6/6	0.88	0.16	2.21	41,58,66,69	0
6	CA	A	516	1/1	0.99	0.14	1.87	17,17,17,17	0
6	CA	B	509	1/1	0.99	0.08	1.26	27,27,27,27	0
4	CL	A	512	1/1	0.98	0.10	0.27	44,44,44,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	AVJ	B	501	14/14	0.94	0.08	-0.08	25,31,33,33	0
2	AVJ	A	501	14/14	0.96	0.08	-0.15	15,19,21,23	0
4	CL	A	514	1/1	0.96	0.08	-0.48	36,36,36,36	0
3	GOL	A	503	6/6	0.97	0.10	-0.84	18,22,24,24	0
4	CL	A	510	1/1	0.99	0.05	-1.09	24,24,24,24	0
7	MG	B	510	1/1	0.95	0.07	-1.39	50,50,50,50	0
4	CL	B	506	1/1	0.98	0.06	-1.46	41,41,41,41	0
5	FE	B	508	1/1	0.99	0.05	-	24,24,24,24	0
4	CL	A	513	1/1	0.94	0.14	-	51,51,51,51	0
5	FE	A	515	1/1	1.00	0.13	-	15,15,15,15	0
7	MG	A	517	1/1	0.27	0.29	-	102,102,102,102	0
4	CL	A	511	1/1	0.99	0.08	-	29,29,29,29	0
4	CL	B	507	1/1	0.98	0.12	-	38,38,38,38	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.