



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 26, 2025 – 06:05 PM EDT

PDB ID : 9CLG / pdb\_00009clg  
Title : Importin alpha isoform 2 with synthetic zero net-charge nuclear localization signal  
Authors : Zeytuni, N.; Yang, Z.; Odeh-Ahmed, A.; Abdoli, A.; Leyton, V.J.; Bednova, O.; Lavigne, P.; Lemieux, B.  
Deposited on : 2024-07-11  
Resolution : 2.42 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0rc1  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.006 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.45.1

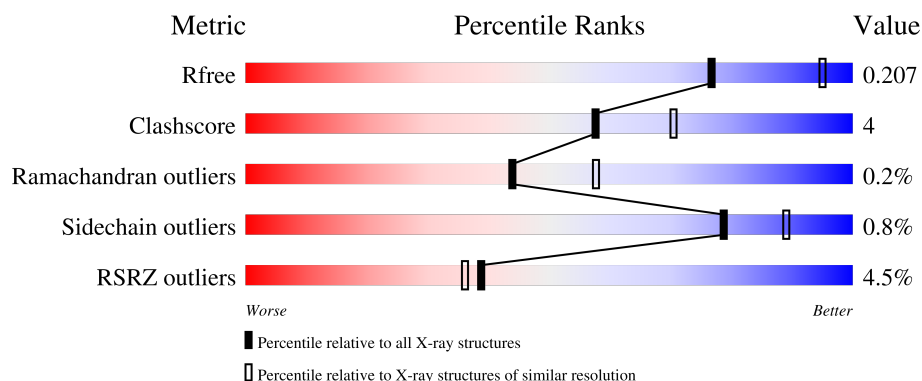
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.42 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))
$R_{free}$	164625	5670 (2.44-2.40)
Clashscore	180529	6299 (2.44-2.40)
Ramachandran outliers	177936	6232 (2.44-2.40)
Sidechain outliers	177891	6233 (2.44-2.40)
RSRZ outliers	164620	5670 (2.44-2.40)

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 of 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	510	 2% 75% 8% 17%
2	B	9	 67% 44% 22% 33%
2	C	9	 44% 89% 11%

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3631 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 Importin subunit alpha-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	425	3264	2080	557	617	10	0	4	0

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	MET	-	initiating methionine	UNP P52293
A	21	HIS	-	expression tag	UNP P52293
A	22	HIS	-	expression tag	UNP P52293
A	23	HIS	-	expression tag	UNP P52293
A	24	HIS	-	expression tag	UNP P52293
A	25	HIS	-	expression tag	UNP P52293
A	26	HIS	-	expression tag	UNP P52293
A	27	SER	-	expression tag	UNP P52293
A	28	SER	-	expression tag	UNP P52293
A	29	GLY	-	expression tag	UNP P52293
A	30	LEU	-	expression tag	UNP P52293
A	31	VAL	-	expression tag	UNP P52293
A	32	PRO	-	expression tag	UNP P52293
A	33	ARG	-	expression tag	UNP P52293
A	34	GLY	-	expression tag	UNP P52293
A	35	SER	-	expression tag	UNP P52293
A	36	GLY	-	expression tag	UNP P52293
A	37	MET	-	expression tag	UNP P52293
A	38	LYS	-	expression tag	UNP P52293
A	39	GLU	-	expression tag	UNP P52293
A	40	THR	-	expression tag	UNP P52293
A	41	ALA	-	expression tag	UNP P52293
A	42	ALA	-	expression tag	UNP P52293
A	43	ALA	-	expression tag	UNP P52293
A	44	LYS	-	expression tag	UNP P52293
A	45	PHE	-	expression tag	UNP P52293
A	46	GLU	-	expression tag	UNP P52293

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Chain	Residue	Modelled	Actual	Comment	Reference
A	47	ARG	-	expression tag	UNP P52293
A	48	GLN	-	expression tag	UNP P52293
A	49	HIS	-	expression tag	UNP P52293
A	50	MET	-	expression tag	UNP P52293
A	51	ASP	-	expression tag	UNP P52293
A	52	SER	-	expression tag	UNP P52293
A	53	PRO	-	expression tag	UNP P52293
A	54	ASP	-	expression tag	UNP P52293
A	55	LEU	-	expression tag	UNP P52293
A	56	GLY	-	expression tag	UNP P52293
A	57	THR	-	expression tag	UNP P52293
A	58	ASP	-	expression tag	UNP P52293
A	59	ASP	-	expression tag	UNP P52293
A	60	ASP	-	expression tag	UNP P52293
A	61	ASP	-	expression tag	UNP P52293
A	62	LYS	-	expression tag	UNP P52293
A	63	ALA	-	expression tag	UNP P52293
A	64	MET	-	expression tag	UNP P52293
A	65	ALA	-	expression tag	UNP P52293
A	66	ASP	-	expression tag	UNP P52293
A	67	ILE	-	expression tag	UNP P52293
A	68	GLY	-	expression tag	UNP P52293
A	69	SER	-	expression tag	UNP P52293

- Molecule 2 is a protein called Zero net-charge NLS.

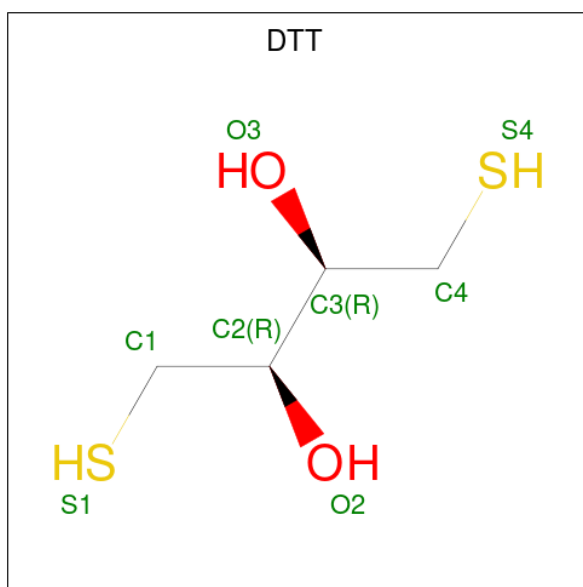
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	6	Total	C	N	O	0	0	0
			50	33	9	8			
2	C	9	Total	C	N	O	0	0	0
			75	48	12	15			

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (CCD ID: DTT) (formula:  $C_4H_{10}O_2S_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			8	4	2	2		
4	A	1	Total	C	O	S	0	0
			8	4	2	2		

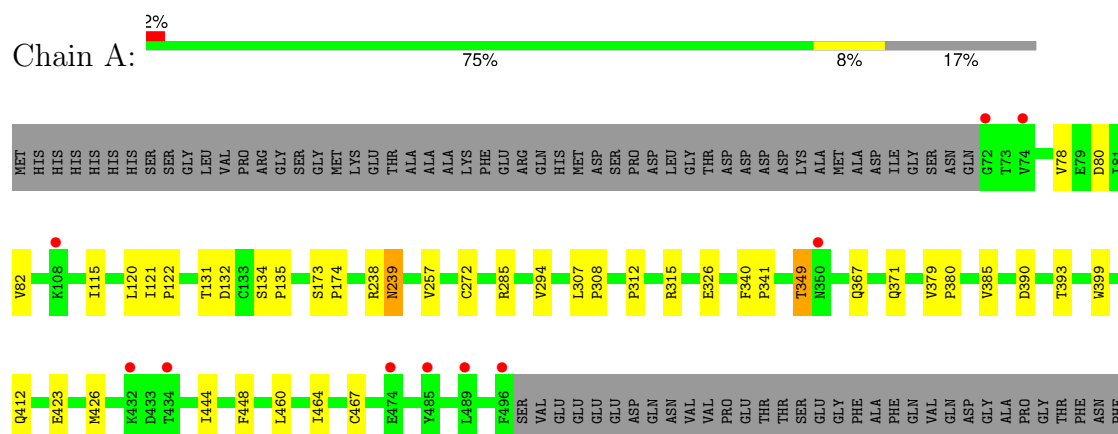
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	187	Total	O	0	0
			187	187		
5	B	1	Total	O	0	0
			1	1		
5	C	2	Total	O	0	0
			2	2		

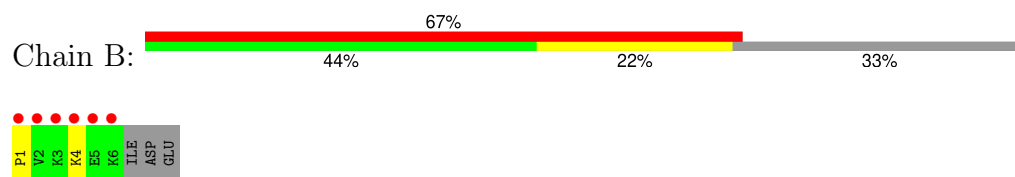
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

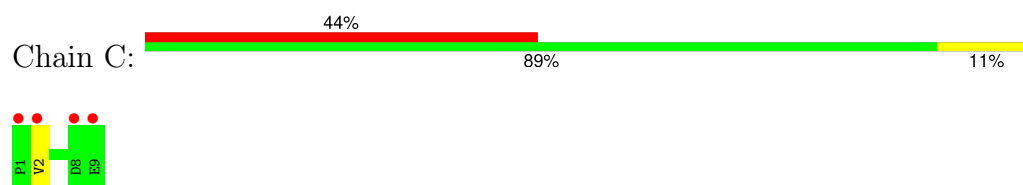
- Molecule 1: Importin subunit alpha-1



- Molecule 2: Zero net-charge NLS



- Molecule 2: Zero net-charge NLS



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.01Å 89.96Å 99.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.63 – 2.42 49.63 – 2.42	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.63-2.42) 99.6 (49.63-2.42)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.67 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.8.0419	Depositor
R, $R_{free}$	0.161 , 0.202 0.169 , 0.207	Depositor DCC
$R_{free}$ test set	1268 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.3	Xtriage
Anisotropy	0.458	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 43.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3631	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.53% 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

### 5.1 Standard geometry

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

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	0.62	0/3335	1.03	1/4543 (0.0%)
2	B	0.76	0/50	1.40	1/63 (1.6%)
2	C	0.69	0/75	1.40	0/97
All	All	0.63	0/3460	1.04	2/4703 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	PRO	CA-N-CD	-5.80	103.88	112.00
1	A	349	THR	CA-CB-OG1	-5.39	101.51	109.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	285	ARG	Sidechain
1	A	315	ARG	Sidechain

## 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	3264	0	3353	29	0
2	B	50	0	63	3	0
2	C	75	0	84	0	0
3	A	36	0	48	3	0
4	A	16	0	20	2	0
5	A	187	0	0	6	0
5	B	1	0	0	0	0
5	C	2	0	0	0	0
All	All	3631	0	3568	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 4.

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:A:371:GLN:HE21	1:A:412:GLN:HE21	1.32	0.77
5:A:827:HOH:O	2:B:4:LYS:HG2	1.96	0.66
1:A:367:GLN:OE1	5:A:701:HOH:O	2.15	0.60
1:A:426:MET:HE3	1:A:444:ILE:HD11	1.85	0.59
1:A:131:THR:HG23	3:A:602:GOL:H31	1.87	0.56
1:A:80:ASP:OD2	5:A:702:HOH:O	2.18	0.55
1:A:371:GLN:NE2	1:A:412:GLN:HE21	2.04	0.52
1:A:173:SER:HA	3:A:602:GOL:H11	1.94	0.50
1:A:174:PRO:HD2	3:A:602:GOL:H32	1.94	0.50
1:A:257:VAL:HG22	1:A:294:VAL:CG1	2.41	0.49
1:A:426:MET:CE	1:A:444:ILE:HD11	2.42	0.49
1:A:340:PHE:N	1:A:341:PRO:CD	2.75	0.49
1:A:272:CYS:HB3	1:A:312:PRO:HB2	1.96	0.48
1:A:448:PHE:CD1	1:A:460:LEU:HD23	2.49	0.47
1:A:78:VAL:O	1:A:82:VAL:HG23	2.14	0.47
1:A:399:TRP:CD1	2:B:4:LYS:HE2	2.49	0.47
1:A:238[A]:ARG:HH22	4:A:608:DTT:HS2	1.61	0.45
1:A:326:GLU:OE1	5:A:703:HOH:O	2.21	0.45
1:A:399:TRP:CG	2:B:4:LYS:HE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:385:VAL:HG13	1:A:393:THR:HG22	1.97	0.45
4:A:608:DTT:S1	5:A:720:HOH:O	2.61	0.45
1:A:132:ASP:OD1	5:A:704:HOH:O	2.21	0.44
1:A:307:LEU:N	1:A:308:PRO:CD	2.81	0.43
1:A:385:VAL:CG1	1:A:393:THR:HG22	2.49	0.43
1:A:423:GLU:HG3	1:A:467:CYS:HB2	2.00	0.42
1:A:134:SER:N	1:A:135:PRO:CD	2.83	0.41
1:A:121:ILE:N	1:A:122:PRO:CD	2.84	0.41
1:A:379:VAL:N	1:A:380:PRO:CD	2.83	0.41
1:A:460:LEU:O	1:A:464:ILE:HG13	2.21	0.40
1:A:115:ILE:HG23	1:A:120:LEU:HD12	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	427/510 (84%)	420 (98%)	6 (1%)	1 (0%)	44	58
2	B	4/9 (44%)	4 (100%)	0	0	100	100
2	C	7/9 (78%)	7 (100%)	0	0	100	100
All	All	438/528 (83%)	431 (98%)	6 (1%)	1 (0%)	44	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	239	ASN

### 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	360/426 (84%)	358 (99%)	2 (1%)	84	92
2	B	6/9 (67%)	6 (100%)	0	100	100
2	C	9/9 (100%)	8 (89%)	1 (11%)	5	6
All	All	375/444 (84%)	372 (99%)	3 (1%)	79	89

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

Mol	Chain	Res	Type
1	A	349	THR
1	A	390	ASP
2	C	2	VAL

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

Mol	Chain	Res	Type
1	A	95	GLN
1	A	98	GLN
1	A	350	ASN
1	A	371	GLN
1	A	375	ASN
1	A	438	GLN
1	A	479	HIS

### 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 [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

8 ligands 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	$\# Z  > 2$
4	DTT	A	607	-	7,7,7	0.14	0	4,8,8	0.96	0
3	GOL	A	601	-	5,5,5	0.14	0	5,5,5	0.47	0
3	GOL	A	603	-	5,5,5	0.14	0	5,5,5	0.44	0
3	GOL	A	606	-	5,5,5	0.15	0	5,5,5	0.36	0
3	GOL	A	605	-	5,5,5	0.16	0	5,5,5	0.44	0
3	GOL	A	604	-	5,5,5	0.42	0	5,5,5	0.63	0
4	DTT	A	608	-	7,7,7	0.76	0	4,8,8	1.49	0
3	GOL	A	602	-	5,5,5	0.24	0	5,5,5	0.72	0

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
4	DTT	A	607	-	-	4/8/8/8	-
3	GOL	A	601	-	-	2/4/4/4	-
3	GOL	A	603	-	-	1/4/4/4	-
3	GOL	A	606	-	-	2/4/4/4	-
3	GOL	A	605	-	-	2/4/4/4	-
3	GOL	A	604	-	-	1/4/4/4	-
4	DTT	A	608	-	-	4/8/8/8	-
3	GOL	A	602	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	605	GOL	C1-C2-C3-O3
4	A	607	DTT	S1-C1-C2-C3
4	A	607	DTT	C2-C3-C4-S4
4	A	607	DTT	O3-C3-C4-S4
4	A	608	DTT	C1-C2-C3-C4
4	A	608	DTT	O2-C2-C3-C4
3	A	601	GOL	O1-C1-C2-C3
3	A	603	GOL	O1-C1-C2-C3
3	A	606	GOL	C1-C2-C3-O3
3	A	601	GOL	O1-C1-C2-O2
3	A	605	GOL	O2-C2-C3-O3
4	A	608	DTT	O2-C2-C3-O3
4	A	608	DTT	C1-C2-C3-O3
4	A	607	DTT	S1-C1-C2-O2
3	A	606	GOL	O2-C2-C3-O3
3	A	604	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	608	DTT	2	0
3	A	602	GOL	3	0

## 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

### 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	425/510 (83%)	-0.26	10 (2%) 59 57	20, 39, 88, 111	4 (0%)
2	B	6/9 (66%)	4.79	6 (100%) 0 0	96, 103, 123, 129	0
2	C	9/9 (100%)	1.71	4 (44%) 1 1	40, 65, 103, 117	0
All	All	440/528 (83%)	-0.15	20 (4%) 39 36	20, 40, 92, 129	4 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	2	VAL	8.0
2	B	1	PRO	7.5
2	C	1	PRO	5.2
2	B	6	LYS	4.6
1	A	74	VAL	4.1
1	A	432	LYS	3.6
1	A	72	GLY	3.5
2	B	4	LYS	3.2
1	A	496	PHE	3.1
1	A	489	LEU	3.0
2	C	2	VAL	2.9
2	B	5	GLU	2.8
1	A	108	LYS	2.7
2	B	3	LYS	2.7
1	A	485	TYR	2.6
1	A	434	THR	2.6
2	C	8	ASP	2.6
2	C	9	GLU	2.5
1	A	474	GLU	2.3
1	A	350	ASN	2.2

## 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 oligosaccharides 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( $\text{\AA}^2$ )	Q<0.9
3	GOL	A	604	6/6	0.74	0.26	59,75,88,94	0
3	GOL	A	602	6/6	0.88	0.15	63,75,79,81	0
4	DTT	A	607	8/8	0.88	0.19	78,91,96,113	0
4	DTT	A	608	8/8	0.89	0.18	92,95,100,109	0
3	GOL	A	606	6/6	0.90	0.16	69,74,82,85	0
3	GOL	A	601	6/6	0.92	0.15	61,69,71,72	0
3	GOL	A	603	6/6	0.92	0.14	64,67,72,75	0
3	GOL	A	605	6/6	0.94	0.12	52,60,64,76	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.