



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:49 AM UTC

PDB ID : 5OVW / pdb_00005ovw
Title : Nanobody-bound BtuF, the vitamin B12 binding protein in Escherichia coli
Authors : Mireku, S.A.; Sauer, M.M.; Glockshuber, R.; Locher, K.P.
Deposited on : 2017-08-30
Resolution : 2.65 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

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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

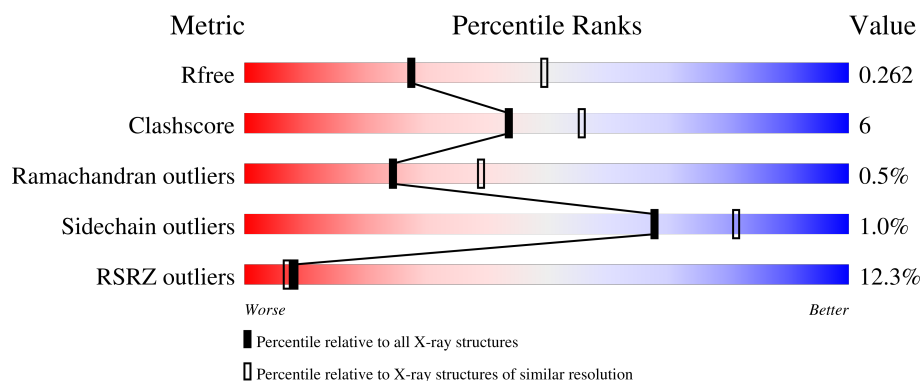
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 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}	180053	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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 ≥ 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	289	
1	B	289	
1	C	289	
1	D	289	
1	E	289	

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Mol	Chain	Length	Quality of chain
1	F	289	
2	G	159	
2	H	159	
2	I	159	
2	J	159	
2	K	159	
2	L	159	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17352 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 Vitamin B12-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total 1907	C 1216	N 332	O 355	S 4	0	0	0
1	B	245	Total 1907	C 1216	N 332	O 355	S 4	0	0	0
1	C	245	Total 1907	C 1216	N 332	O 355	S 4	0	0	0
1	D	245	Total 1907	C 1216	N 332	O 355	S 4	0	0	0
1	E	245	Total 1907	C 1216	N 332	O 355	S 4	0	0	0
1	F	245	Total 1907	C 1216	N 332	O 355	S 4	0	0	0

There are 264 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	MET	-	initiating methionine	UNP P37028
A	-1	LYS	-	expression tag	UNP P37028
A	0	LYS	-	expression tag	UNP P37028
A	1	THR	-	expression tag	UNP P37028
A	2	ALA	-	expression tag	UNP P37028
A	3	ILE	-	expression tag	UNP P37028
A	4	ALA	-	expression tag	UNP P37028
A	5	ILE	-	expression tag	UNP P37028
A	6	ALA	-	expression tag	UNP P37028
A	7	VAL	-	expression tag	UNP P37028
A	8	ALA	-	expression tag	UNP P37028
A	9	LEU	-	expression tag	UNP P37028
A	10	ALA	-	expression tag	UNP P37028
A	11	GLY	-	expression tag	UNP P37028
A	12	PHE	-	expression tag	UNP P37028
A	13	ALA	-	expression tag	UNP P37028
A	14	THR	-	expression tag	UNP P37028

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Chain	Residue	Modelled	Actual	Comment	Reference
A	15	VAL	-	expression tag	UNP P37028
A	16	ALA	-	expression tag	UNP P37028
A	17	GLN	-	expression tag	UNP P37028
A	18	ALA	-	expression tag	UNP P37028
A	19	ALA	-	expression tag	UNP P37028
A	20	SER	-	expression tag	UNP P37028
A	21	MET	-	expression tag	UNP P37028
A	267	SER	-	expression tag	UNP P37028
A	268	GLY	-	expression tag	UNP P37028
A	269	SER	-	expression tag	UNP P37028
A	270	LEU	-	expression tag	UNP P37028
A	271	GLU	-	expression tag	UNP P37028
A	272	VAL	-	expression tag	UNP P37028
A	273	LEU	-	expression tag	UNP P37028
A	274	PHE	-	expression tag	UNP P37028
A	275	GLN	-	expression tag	UNP P37028
A	276	GLY	-	expression tag	UNP P37028
A	277	PRO	-	expression tag	UNP P37028
A	278	GLY	-	expression tag	UNP P37028
A	279	GLY	-	expression tag	UNP P37028
A	280	SER	-	expression tag	UNP P37028
A	281	HIS	-	expression tag	UNP P37028
A	282	HIS	-	expression tag	UNP P37028
A	283	HIS	-	expression tag	UNP P37028
A	284	HIS	-	expression tag	UNP P37028
A	285	HIS	-	expression tag	UNP P37028
A	286	HIS	-	expression tag	UNP P37028
B	-2	MET	-	initiating methionine	UNP P37028
B	-1	LYS	-	expression tag	UNP P37028
B	0	LYS	-	expression tag	UNP P37028
B	1	THR	-	expression tag	UNP P37028
B	2	ALA	-	expression tag	UNP P37028
B	3	ILE	-	expression tag	UNP P37028
B	4	ALA	-	expression tag	UNP P37028
B	5	ILE	-	expression tag	UNP P37028
B	6	ALA	-	expression tag	UNP P37028
B	7	VAL	-	expression tag	UNP P37028
B	8	ALA	-	expression tag	UNP P37028
B	9	LEU	-	expression tag	UNP P37028
B	10	ALA	-	expression tag	UNP P37028
B	11	GLY	-	expression tag	UNP P37028
B	12	PHE	-	expression tag	UNP P37028

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Chain	Residue	Modelled	Actual	Comment	Reference
B	13	ALA	-	expression tag	UNP P37028
B	14	THR	-	expression tag	UNP P37028
B	15	VAL	-	expression tag	UNP P37028
B	16	ALA	-	expression tag	UNP P37028
B	17	GLN	-	expression tag	UNP P37028
B	18	ALA	-	expression tag	UNP P37028
B	19	ALA	-	expression tag	UNP P37028
B	20	SER	-	expression tag	UNP P37028
B	21	MET	-	expression tag	UNP P37028
B	267	SER	-	expression tag	UNP P37028
B	268	GLY	-	expression tag	UNP P37028
B	269	SER	-	expression tag	UNP P37028
B	270	LEU	-	expression tag	UNP P37028
B	271	GLU	-	expression tag	UNP P37028
B	272	VAL	-	expression tag	UNP P37028
B	273	LEU	-	expression tag	UNP P37028
B	274	PHE	-	expression tag	UNP P37028
B	275	GLN	-	expression tag	UNP P37028
B	276	GLY	-	expression tag	UNP P37028
B	277	PRO	-	expression tag	UNP P37028
B	278	GLY	-	expression tag	UNP P37028
B	279	GLY	-	expression tag	UNP P37028
B	280	SER	-	expression tag	UNP P37028
B	281	HIS	-	expression tag	UNP P37028
B	282	HIS	-	expression tag	UNP P37028
B	283	HIS	-	expression tag	UNP P37028
B	284	HIS	-	expression tag	UNP P37028
B	285	HIS	-	expression tag	UNP P37028
B	286	HIS	-	expression tag	UNP P37028
C	-2	MET	-	initiating methionine	UNP P37028
C	-1	LYS	-	expression tag	UNP P37028
C	0	LYS	-	expression tag	UNP P37028
C	1	THR	-	expression tag	UNP P37028
C	2	ALA	-	expression tag	UNP P37028
C	3	ILE	-	expression tag	UNP P37028
C	4	ALA	-	expression tag	UNP P37028
C	5	ILE	-	expression tag	UNP P37028
C	6	ALA	-	expression tag	UNP P37028
C	7	VAL	-	expression tag	UNP P37028
C	8	ALA	-	expression tag	UNP P37028
C	9	LEU	-	expression tag	UNP P37028
C	10	ALA	-	expression tag	UNP P37028

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Chain	Residue	Modelled	Actual	Comment	Reference
C	11	GLY	-	expression tag	UNP P37028
C	12	PHE	-	expression tag	UNP P37028
C	13	ALA	-	expression tag	UNP P37028
C	14	THR	-	expression tag	UNP P37028
C	15	VAL	-	expression tag	UNP P37028
C	16	ALA	-	expression tag	UNP P37028
C	17	GLN	-	expression tag	UNP P37028
C	18	ALA	-	expression tag	UNP P37028
C	19	ALA	-	expression tag	UNP P37028
C	20	SER	-	expression tag	UNP P37028
C	21	MET	-	expression tag	UNP P37028
C	267	SER	-	expression tag	UNP P37028
C	268	GLY	-	expression tag	UNP P37028
C	269	SER	-	expression tag	UNP P37028
C	270	LEU	-	expression tag	UNP P37028
C	271	GLU	-	expression tag	UNP P37028
C	272	VAL	-	expression tag	UNP P37028
C	273	LEU	-	expression tag	UNP P37028
C	274	PHE	-	expression tag	UNP P37028
C	275	GLN	-	expression tag	UNP P37028
C	276	GLY	-	expression tag	UNP P37028
C	277	PRO	-	expression tag	UNP P37028
C	278	GLY	-	expression tag	UNP P37028
C	279	GLY	-	expression tag	UNP P37028
C	280	SER	-	expression tag	UNP P37028
C	281	HIS	-	expression tag	UNP P37028
C	282	HIS	-	expression tag	UNP P37028
C	283	HIS	-	expression tag	UNP P37028
C	284	HIS	-	expression tag	UNP P37028
C	285	HIS	-	expression tag	UNP P37028
C	286	HIS	-	expression tag	UNP P37028
D	-2	MET	-	initiating methionine	UNP P37028
D	-1	LYS	-	expression tag	UNP P37028
D	0	LYS	-	expression tag	UNP P37028
D	1	THR	-	expression tag	UNP P37028
D	2	ALA	-	expression tag	UNP P37028
D	3	ILE	-	expression tag	UNP P37028
D	4	ALA	-	expression tag	UNP P37028
D	5	ILE	-	expression tag	UNP P37028
D	6	ALA	-	expression tag	UNP P37028
D	7	VAL	-	expression tag	UNP P37028
D	8	ALA	-	expression tag	UNP P37028

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Chain	Residue	Modelled	Actual	Comment	Reference
D	9	LEU	-	expression tag	UNP P37028
D	10	ALA	-	expression tag	UNP P37028
D	11	GLY	-	expression tag	UNP P37028
D	12	PHE	-	expression tag	UNP P37028
D	13	ALA	-	expression tag	UNP P37028
D	14	THR	-	expression tag	UNP P37028
D	15	VAL	-	expression tag	UNP P37028
D	16	ALA	-	expression tag	UNP P37028
D	17	GLN	-	expression tag	UNP P37028
D	18	ALA	-	expression tag	UNP P37028
D	19	ALA	-	expression tag	UNP P37028
D	20	SER	-	expression tag	UNP P37028
D	21	MET	-	expression tag	UNP P37028
D	267	SER	-	expression tag	UNP P37028
D	268	GLY	-	expression tag	UNP P37028
D	269	SER	-	expression tag	UNP P37028
D	270	LEU	-	expression tag	UNP P37028
D	271	GLU	-	expression tag	UNP P37028
D	272	VAL	-	expression tag	UNP P37028
D	273	LEU	-	expression tag	UNP P37028
D	274	PHE	-	expression tag	UNP P37028
D	275	GLN	-	expression tag	UNP P37028
D	276	GLY	-	expression tag	UNP P37028
D	277	PRO	-	expression tag	UNP P37028
D	278	GLY	-	expression tag	UNP P37028
D	279	GLY	-	expression tag	UNP P37028
D	280	SER	-	expression tag	UNP P37028
D	281	HIS	-	expression tag	UNP P37028
D	282	HIS	-	expression tag	UNP P37028
D	283	HIS	-	expression tag	UNP P37028
D	284	HIS	-	expression tag	UNP P37028
D	285	HIS	-	expression tag	UNP P37028
D	286	HIS	-	expression tag	UNP P37028
E	-2	MET	-	initiating methionine	UNP P37028
E	-1	LYS	-	expression tag	UNP P37028
E	0	LYS	-	expression tag	UNP P37028
E	1	THR	-	expression tag	UNP P37028
E	2	ALA	-	expression tag	UNP P37028
E	3	ILE	-	expression tag	UNP P37028
E	4	ALA	-	expression tag	UNP P37028
E	5	ILE	-	expression tag	UNP P37028
E	6	ALA	-	expression tag	UNP P37028

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Chain	Residue	Modelled	Actual	Comment	Reference
E	7	VAL	-	expression tag	UNP P37028
E	8	ALA	-	expression tag	UNP P37028
E	9	LEU	-	expression tag	UNP P37028
E	10	ALA	-	expression tag	UNP P37028
E	11	GLY	-	expression tag	UNP P37028
E	12	PHE	-	expression tag	UNP P37028
E	13	ALA	-	expression tag	UNP P37028
E	14	THR	-	expression tag	UNP P37028
E	15	VAL	-	expression tag	UNP P37028
E	16	ALA	-	expression tag	UNP P37028
E	17	GLN	-	expression tag	UNP P37028
E	18	ALA	-	expression tag	UNP P37028
E	19	ALA	-	expression tag	UNP P37028
E	20	SER	-	expression tag	UNP P37028
E	21	MET	-	expression tag	UNP P37028
E	267	SER	-	expression tag	UNP P37028
E	268	GLY	-	expression tag	UNP P37028
E	269	SER	-	expression tag	UNP P37028
E	270	LEU	-	expression tag	UNP P37028
E	271	GLU	-	expression tag	UNP P37028
E	272	VAL	-	expression tag	UNP P37028
E	273	LEU	-	expression tag	UNP P37028
E	274	PHE	-	expression tag	UNP P37028
E	275	GLN	-	expression tag	UNP P37028
E	276	GLY	-	expression tag	UNP P37028
E	277	PRO	-	expression tag	UNP P37028
E	278	GLY	-	expression tag	UNP P37028
E	279	GLY	-	expression tag	UNP P37028
E	280	SER	-	expression tag	UNP P37028
E	281	HIS	-	expression tag	UNP P37028
E	282	HIS	-	expression tag	UNP P37028
E	283	HIS	-	expression tag	UNP P37028
E	284	HIS	-	expression tag	UNP P37028
E	285	HIS	-	expression tag	UNP P37028
E	286	HIS	-	expression tag	UNP P37028
F	-2	MET	-	initiating methionine	UNP P37028
F	-1	LYS	-	expression tag	UNP P37028
F	0	LYS	-	expression tag	UNP P37028
F	1	THR	-	expression tag	UNP P37028
F	2	ALA	-	expression tag	UNP P37028
F	3	ILE	-	expression tag	UNP P37028
F	4	ALA	-	expression tag	UNP P37028

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Chain	Residue	Modelled	Actual	Comment	Reference
F	5	ILE	-	expression tag	UNP P37028
F	6	ALA	-	expression tag	UNP P37028
F	7	VAL	-	expression tag	UNP P37028
F	8	ALA	-	expression tag	UNP P37028
F	9	LEU	-	expression tag	UNP P37028
F	10	ALA	-	expression tag	UNP P37028
F	11	GLY	-	expression tag	UNP P37028
F	12	PHE	-	expression tag	UNP P37028
F	13	ALA	-	expression tag	UNP P37028
F	14	THR	-	expression tag	UNP P37028
F	15	VAL	-	expression tag	UNP P37028
F	16	ALA	-	expression tag	UNP P37028
F	17	GLN	-	expression tag	UNP P37028
F	18	ALA	-	expression tag	UNP P37028
F	19	ALA	-	expression tag	UNP P37028
F	20	SER	-	expression tag	UNP P37028
F	21	MET	-	expression tag	UNP P37028
F	267	SER	-	expression tag	UNP P37028
F	268	GLY	-	expression tag	UNP P37028
F	269	SER	-	expression tag	UNP P37028
F	270	LEU	-	expression tag	UNP P37028
F	271	GLU	-	expression tag	UNP P37028
F	272	VAL	-	expression tag	UNP P37028
F	273	LEU	-	expression tag	UNP P37028
F	274	PHE	-	expression tag	UNP P37028
F	275	GLN	-	expression tag	UNP P37028
F	276	GLY	-	expression tag	UNP P37028
F	277	PRO	-	expression tag	UNP P37028
F	278	GLY	-	expression tag	UNP P37028
F	279	GLY	-	expression tag	UNP P37028
F	280	SER	-	expression tag	UNP P37028
F	281	HIS	-	expression tag	UNP P37028
F	282	HIS	-	expression tag	UNP P37028
F	283	HIS	-	expression tag	UNP P37028
F	284	HIS	-	expression tag	UNP P37028
F	285	HIS	-	expression tag	UNP P37028
F	286	HIS	-	expression tag	UNP P37028

- Molecule 2 is a protein called Nanobody.

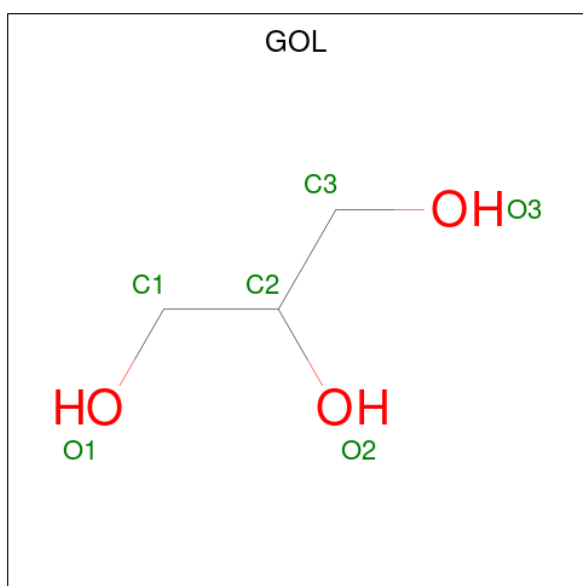
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	126	Total	C	N	O	S	0	0	0
			936	578	165	187	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	126	Total	C	N	O	S	0	0	0
			936	578	165	187	6			
2	I	126	Total	C	N	O	S	0	0	0
			936	578	165	187	6			
2	J	126	Total	C	N	O	S	0	0	0
			936	578	165	187	6			
2	K	126	Total	C	N	O	S	0	0	0
			936	578	165	187	6			
2	L	126	Total	C	N	O	S	0	0	0
			936	578	165	187	6			

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

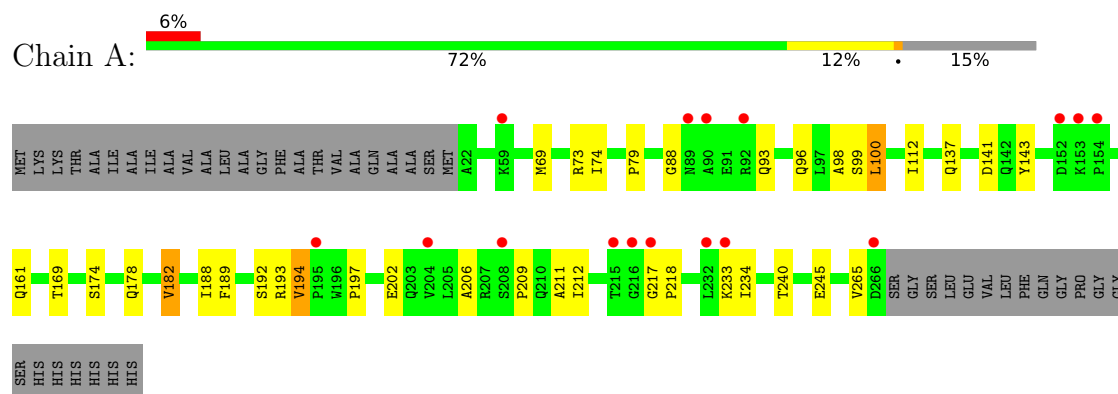
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	42	Total 42	O 42	0	0
4	B	37	Total 37	O 37	0	0
4	C	10	Total 10	O 10	0	0
4	D	66	Total 66	O 66	0	0
4	E	25	Total 25	O 25	0	0
4	F	17	Total 17	O 17	0	0
4	G	4	Total 4	O 4	0	0
4	H	14	Total 14	O 14	0	0
4	I	2	Total 2	O 2	0	0
4	J	27	Total 27	O 27	0	0
4	K	13	Total 13	O 13	0	0
4	L	7	Total 7	O 7	0	0

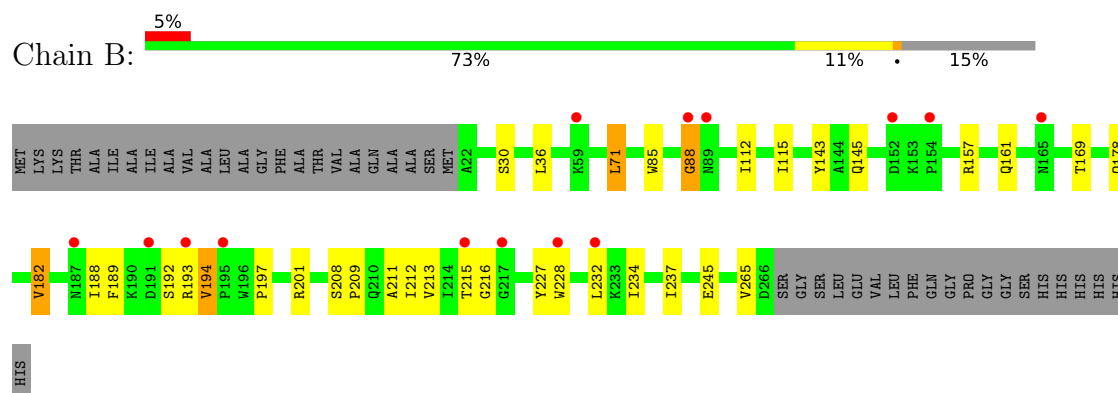
3 Residue-property plots

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.

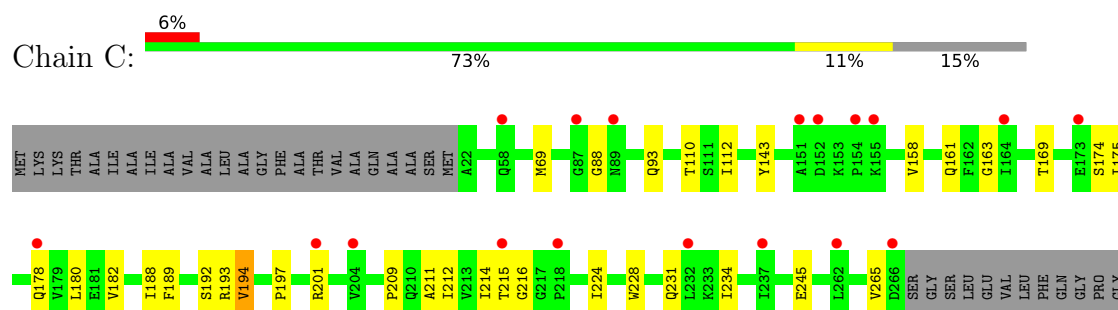
• Molecule 1: Vitamin B12-binding protein



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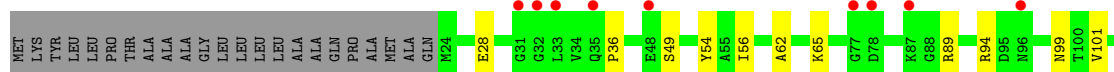


• Molecule 1: Vitamin B12-binding protein

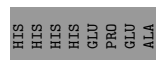
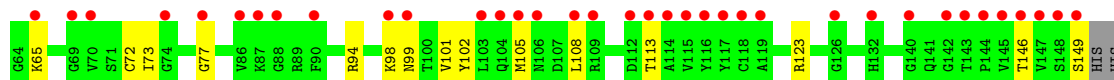
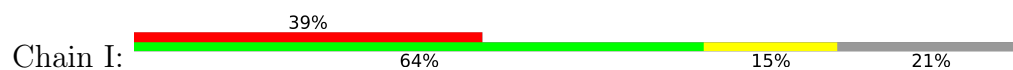




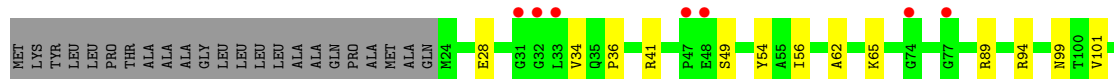
• Molecule 2: Nanobody



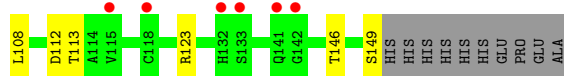
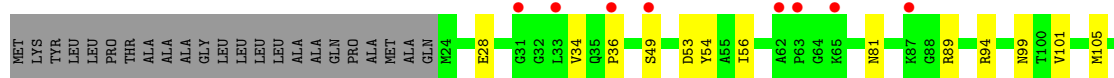
• Molecule 2: Nanobody



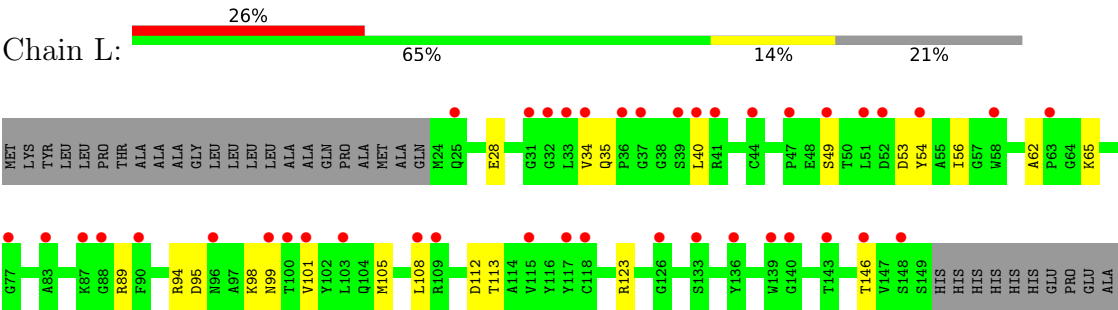
• Molecule 2: Nanobody



• Molecule 2: Nanobody



● Molecule 2: Nanobody



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.80Å 141.00Å 216.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.65 20.00 – 2.65	Depositor EDS
% Data completeness (in resolution range)	85.3 (20.00-2.65) 85.1 (20.00-2.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.95 (at 2.67Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: ???)	Depositor
R, R_{free}	0.223 , 0.259 0.228 , 0.262	Depositor DCC
R_{free} test set	3239 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	47.6	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 78.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	17352	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.07% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: 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	0.13	0/1951	0.34	0/2661
1	B	0.14	0/1951	0.33	0/2661
1	C	0.15	0/1951	0.33	0/2661
1	D	0.14	0/1951	0.33	0/2661
1	E	0.14	0/1951	0.32	0/2661
1	F	0.13	0/1951	0.33	0/2661
2	G	0.10	0/955	0.31	0/1295
2	H	0.12	0/955	0.31	0/1295
2	I	0.10	0/955	0.30	0/1295
2	J	0.14	0/955	0.33	0/1295
2	K	0.11	0/955	0.32	0/1295
2	L	0.12	0/955	0.33	0/1295
All	All	0.13	0/17436	0.33	0/23736

There are no bond length outliers.

There are no bond angle outliers.

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	1907	0	1924	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1907	0	1924	26	0
1	C	1907	0	1924	26	0
1	D	1907	0	1924	25	0
1	E	1907	0	1924	18	0
1	F	1907	0	1924	25	0
2	G	936	0	888	12	0
2	H	936	0	888	12	0
2	I	936	0	888	13	0
2	J	936	0	888	12	0
2	K	936	0	888	12	0
2	L	936	0	888	13	0
3	B	6	0	8	1	0
3	D	6	0	8	0	0
3	E	6	0	8	0	0
3	F	6	0	8	1	0
3	H	6	0	8	0	0
4	A	42	0	0	0	0
4	B	37	0	0	0	0
4	C	10	0	0	0	0
4	D	66	0	0	1	0
4	E	25	0	0	0	0
4	F	17	0	0	0	0
4	G	4	0	0	0	0
4	H	14	0	0	0	0
4	I	2	0	0	0	0
4	J	27	0	0	2	0
4	K	13	0	0	0	0
4	L	7	0	0	0	0
All	All	17352	0	16912	207	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 6.

All (207) 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:F:156:LYS:HZ1	1:F:266:ASP:CG	1.71	0.97
1:E:156:LYS:NZ	1:E:265:VAL:O	2.06	0.88
2:J:41:ARG:NH1	4:J:201:HOH:O	2.12	0.77
1:C:201:ARG:NH1	1:C:231:GLN:OE1	2.24	0.70
2:L:54:TYR:O	2:L:94:ARG:NH2	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:ARG:HH12	1:C:228:TRP:HA	1.57	0.67
2:H:36:PRO:HD2	2:H:149:SER:HB3	1.79	0.64
2:L:34:VAL:HG21	2:L:40:LEU:HG	1.81	0.63
1:A:211:ALA:HB2	1:A:265:VAL:HG11	1.81	0.62
2:H:56:ILE:HG13	2:H:101:VAL:HG21	1.81	0.62
1:E:143:TYR:OH	1:E:178:GLN:NE2	2.32	0.61
2:K:56:ILE:HG13	2:K:101:VAL:HG21	1.83	0.60
1:F:245:GLU:OE2	2:L:123:ARG:HD2	2.01	0.59
1:C:245:GLU:OE1	2:G:123:ARG:NH1	2.34	0.59
1:F:156:LYS:NZ	1:F:266:ASP:OD1	2.36	0.59
1:C:201:ARG:NH1	1:C:228:TRP:HA	2.18	0.59
1:D:143:TYR:CE1	1:D:182:VAL:HG11	2.38	0.59
2:K:36:PRO:HD2	2:K:149:SER:HB3	1.85	0.59
2:J:56:ILE:HG13	2:J:101:VAL:HG21	1.85	0.58
1:B:201:ARG:NH1	1:B:228:TRP:H	2.01	0.58
2:G:54:TYR:O	2:G:94:ARG:NH2	2.37	0.58
1:A:178:GLN:O	1:A:182:VAL:HG13	2.05	0.57
2:H:105:MET:HB3	2:H:108:LEU:HD21	1.87	0.57
2:K:105:MET:HB3	2:K:108:LEU:HD21	1.87	0.57
1:B:209:PRO:O	1:B:234:ILE:HG21	2.06	0.56
2:L:105:MET:HB3	2:L:108:LEU:HD21	1.87	0.56
1:D:245:GLU:OE1	2:J:123:ARG:NH1	2.38	0.55
1:E:211:ALA:HB2	1:E:265:VAL:HG11	1.87	0.55
2:L:28:GLU:N	2:L:28:GLU:OE1	2.39	0.55
2:I:36:PRO:HD2	2:I:149:SER:HB3	1.88	0.55
1:C:174:SER:O	1:C:178:GLN:NE2	2.40	0.55
2:I:113:THR:HG23	2:I:146:THR:HA	1.89	0.55
2:J:105:MET:HB3	2:J:108:LEU:HD21	1.90	0.54
1:F:137:GLN:NE2	1:F:141:ASP:OD1	2.39	0.54
1:B:169:THR:HG21	1:B:189:PHE:CD1	2.43	0.54
1:C:211:ALA:HB2	1:C:265:VAL:HG11	1.91	0.53
1:C:209:PRO:O	1:C:234:ILE:HG21	2.08	0.53
2:H:54:TYR:O	2:H:94:ARG:NH2	2.41	0.53
2:L:56:ILE:HG13	2:L:101:VAL:HG21	1.90	0.53
2:G:56:ILE:HG13	2:G:101:VAL:HG21	1.90	0.53
1:E:137:GLN:NE2	1:E:141:ASP:OD1	2.40	0.53
2:J:28:GLU:OE1	2:J:28:GLU:N	2.42	0.53
1:B:112:ILE:HD12	1:B:143:TYR:HD1	1.74	0.53
1:A:245:GLU:OE1	2:H:123:ARG:NH1	2.41	0.53
1:D:56:GLN:NE2	4:D:403:HOH:O	2.42	0.53
1:E:143:TYR:CE1	1:E:182:VAL:HG11	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:169:THR:HG21	1:E:189:PHE:CD1	2.44	0.52
1:D:85:TRP:CZ2	1:D:88:GLY:HA3	2.45	0.52
1:F:161:GLN:OE1	1:F:212:ILE:HD11	2.10	0.52
1:F:198:GLN:HE21	3:F:301:GOL:H11	1.75	0.52
2:G:28:GLU:N	2:G:28:GLU:OE1	2.42	0.52
2:G:49:SER:O	2:G:99:ASN:ND2	2.41	0.52
1:B:178:GLN:O	1:B:182:VAL:HG13	2.10	0.52
2:H:28:GLU:N	2:H:28:GLU:OE1	2.43	0.52
1:D:201:ARG:NH1	1:D:228:TRP:H	2.08	0.52
1:A:143:TYR:CE1	1:A:182:VAL:HG11	2.45	0.51
1:F:209:PRO:O	1:F:234:ILE:HG21	2.11	0.51
2:K:28:GLU:N	2:K:28:GLU:OE1	2.43	0.51
1:B:201:ARG:HH12	1:B:228:TRP:H	1.58	0.51
1:C:143:TYR:CE1	1:C:182:VAL:HG11	2.46	0.51
2:K:113:THR:HG23	2:K:146:THR:HA	1.92	0.51
1:E:209:PRO:O	1:E:234:ILE:HG21	2.11	0.51
1:B:211:ALA:HB2	1:B:265:VAL:HG11	1.93	0.51
2:I:73:ILE:HD11	2:I:77:GLY:HA2	1.93	0.51
2:I:56:ILE:HG13	2:I:101:VAL:HG21	1.92	0.50
1:B:85:TRP:CZ2	1:B:88:GLY:HA3	2.47	0.50
2:I:105:MET:HB3	2:I:108:LEU:HD21	1.94	0.50
1:F:86:ARG:O	1:F:89:ASN:HB3	2.11	0.50
2:G:105:MET:HB3	2:G:108:LEU:HD21	1.94	0.50
1:F:163:GLY:HA2	2:L:53:ASP:OD2	2.12	0.50
1:C:175:ILE:HA	1:C:178:GLN:HE21	1.76	0.49
1:B:143:TYR:CE1	1:B:182:VAL:HG11	2.47	0.49
1:A:112:ILE:HD12	1:A:143:TYR:HD1	1.77	0.49
1:D:211:ALA:HB2	1:D:265:VAL:HG11	1.94	0.49
1:B:161:GLN:OE1	1:B:212:ILE:HD11	2.12	0.49
1:E:66:TRP:CE2	1:E:67:GLN:HG3	2.47	0.49
1:B:201:ARG:HH11	1:B:228:TRP:HA	1.78	0.49
1:D:66:TRP:CE2	1:D:67:GLN:HG3	2.48	0.48
1:D:209:PRO:O	1:D:234:ILE:HG21	2.13	0.48
2:I:28:GLU:N	2:I:28:GLU:OE1	2.47	0.48
2:J:94:ARG:NH1	4:J:203:HOH:O	2.40	0.48
1:A:209:PRO:O	1:A:234:ILE:HG21	2.14	0.48
2:H:49:SER:O	2:H:99:ASN:ND2	2.47	0.48
1:E:188:ILE:HD13	1:E:209:PRO:HG3	1.96	0.48
1:B:112:ILE:HD12	1:B:143:TYR:CD1	2.49	0.48
1:F:36:LEU:HD21	1:F:115:ILE:HG23	1.96	0.48
2:J:36:PRO:HD2	2:J:149:SER:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:158:VAL:HB	1:F:180:LEU:HD21	1.96	0.48
2:H:130:VAL:HB	2:H:134:SER:HB2	1.96	0.47
1:C:169:THR:HG21	1:C:189:PHE:CD1	2.49	0.47
1:C:175:ILE:HA	1:C:178:GLN:NE2	2.29	0.47
2:G:62:ALA:HB3	2:G:65:LYS:HE3	1.96	0.47
1:B:157:ARG:NE	1:B:208:SER:O	2.44	0.47
1:C:161:GLN:OE1	1:C:212:ILE:HD11	2.15	0.47
2:J:54:TYR:O	2:J:94:ARG:NH2	2.48	0.47
1:A:98:ALA:C	1:A:100:LEU:H	2.23	0.47
1:C:214:ILE:HD12	1:C:224:ILE:HD12	1.96	0.47
1:F:169:THR:HG21	1:F:189:PHE:CD1	2.49	0.47
1:D:169:THR:HG21	1:D:189:PHE:CD1	2.49	0.47
1:B:227:TYR:HD1	1:B:228:TRP:CD1	2.33	0.46
1:D:69:MET:HE1	1:D:93:GLN:O	2.15	0.46
1:F:214:ILE:HD12	1:F:224:ILE:HD12	1.97	0.46
1:F:69:MET:HE1	1:F:93:GLN:O	2.15	0.46
1:D:112:ILE:HD12	1:D:143:TYR:HD1	1.80	0.46
1:F:66:TRP:CE2	1:F:67:GLN:HG3	2.51	0.46
1:F:164:ILE:HD11	1:F:220:GLN:HB2	1.97	0.46
1:C:178:GLN:O	1:C:182:VAL:HG13	2.15	0.46
2:I:98:LYS:HD2	2:I:102:TYR:OH	2.16	0.46
1:D:188:ILE:HD13	1:D:209:PRO:HG3	1.98	0.46
1:F:211:ALA:HB2	1:F:265:VAL:HG11	1.97	0.45
1:A:137:GLN:NE2	1:A:141:ASP:OD1	2.49	0.45
1:E:66:TRP:CD1	2:K:81:ASN:HB3	2.51	0.45
1:E:245:GLU:OE1	2:K:123:ARG:NH1	2.46	0.45
2:G:89:ARG:HH22	2:G:112:ASP:CG	2.24	0.45
2:G:113:THR:HG23	2:G:146:THR:HA	1.98	0.45
1:D:159:PHE:CE2	1:D:161:GLN:HB2	2.51	0.45
1:F:188:ILE:HD13	1:F:209:PRO:HG3	1.98	0.45
2:J:49:SER:O	2:J:99:ASN:ND2	2.47	0.45
1:D:178:GLN:O	1:D:182:VAL:HG13	2.17	0.45
1:F:69:MET:HE2	1:F:93:GLN:HB3	1.97	0.45
1:D:161:GLN:OE1	1:D:212:ILE:HD11	2.17	0.45
2:I:54:TYR:O	2:I:94:ARG:NH2	2.50	0.45
1:D:85:TRP:CE2	1:D:88:GLY:HA3	2.52	0.45
2:J:62:ALA:HB3	2:J:65:LYS:HE3	1.99	0.44
2:L:113:THR:HG23	2:L:146:THR:HA	1.98	0.44
2:L:89:ARG:HH22	2:L:112:ASP:CG	2.26	0.44
1:A:161:GLN:OE1	1:A:212:ILE:HD11	2.16	0.44
1:A:192:SER:O	1:A:194:VAL:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:95:ASP:OD2	2:L:98:LYS:HE3	2.16	0.44
1:B:192:SER:O	1:B:194:VAL:N	2.51	0.44
1:D:73:ARG:HA	1:D:73:ARG:HD2	1.72	0.44
1:F:86:ARG:HD3	1:F:91:GLU:OE2	2.18	0.44
1:A:169:THR:HG21	1:A:189:PHE:CD1	2.51	0.44
1:F:73:ARG:HA	1:F:73:ARG:HD2	1.73	0.44
2:J:34:VAL:HG11	2:J:108:LEU:HD12	2.00	0.44
1:C:194:VAL:HG13	1:C:197:PRO:HB3	2.00	0.44
2:L:49:SER:O	2:L:99:ASN:ND2	2.47	0.44
2:K:53:ASP:OD1	2:K:53:ASP:N	2.42	0.43
1:B:188:ILE:HG13	1:B:189:PHE:CD1	2.54	0.43
1:C:158:VAL:HB	1:C:180:LEU:HD21	2.00	0.43
2:H:62:ALA:HB3	2:H:65:LYS:HE3	2.01	0.43
1:E:157:ARG:NE	1:E:208:SER:O	2.42	0.43
1:A:69:MET:HE2	1:A:93:GLN:HB3	1.99	0.43
1:A:188:ILE:HG13	1:A:189:PHE:CD1	2.53	0.43
1:B:188:ILE:HD13	1:B:209:PRO:HG3	2.01	0.43
1:D:215:THR:HA	1:D:216:GLY:HA2	1.84	0.43
2:L:62:ALA:HB3	2:L:65:LYS:HE3	1.99	0.43
1:F:194:VAL:HG13	1:F:197:PRO:HB3	2.01	0.43
1:B:194:VAL:HG13	1:B:197:PRO:HB3	2.01	0.43
3:B:301:GOL:O3	3:B:301:GOL:O1	2.32	0.42
1:C:69:MET:HE1	1:C:93:GLN:O	2.19	0.42
1:B:36:LEU:HD21	1:B:115:ILE:HG23	2.01	0.42
1:C:192:SER:OG	1:C:197:PRO:HG3	2.20	0.42
1:A:174:SER:O	1:A:178:GLN:NE2	2.52	0.42
1:C:192:SER:O	1:C:194:VAL:N	2.52	0.42
1:D:156:LYS:NZ	1:D:265:VAL:O	2.46	0.42
1:D:188:ILE:HG13	1:D:189:PHE:CD1	2.55	0.42
1:F:201:ARG:NH1	1:F:228:TRP:H	2.16	0.42
2:I:57:GLY:HA2	2:I:72:CYS:HA	2.00	0.42
2:K:49:SER:O	2:K:99:ASN:ND2	2.50	0.42
1:A:202:GLU:O	1:A:206:ALA:N	2.52	0.42
1:C:163:GLY:HA2	2:G:53:ASP:OD2	2.20	0.42
2:I:62:ALA:HB3	2:I:65:LYS:HE3	2.01	0.42
1:A:73:ARG:HA	1:A:73:ARG:HD2	1.77	0.42
2:J:89:ARG:HH22	2:J:112:ASP:CG	2.27	0.42
1:C:69:MET:HE2	1:C:69:MET:HB3	1.94	0.42
1:D:29:LEU:HD22	1:D:64:SER:HB2	2.02	0.42
1:E:156:LYS:HD2	1:E:262:LEU:HG	2.01	0.42
1:F:156:LYS:NZ	1:F:266:ASP:CG	2.57	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:ILE:HD13	1:A:209:PRO:HG3	2.02	0.41
1:E:163:GLY:HA2	2:K:53:ASP:OD2	2.20	0.41
1:E:194:VAL:HG13	1:E:197:PRO:HB3	2.01	0.41
1:B:30:SER:HB3	1:B:85:TRP:CD1	2.55	0.41
1:E:161:GLN:OE1	1:E:212:ILE:HD11	2.21	0.41
1:B:215:THR:HA	1:B:216:GLY:HA2	1.83	0.41
1:C:188:ILE:HG13	1:C:189:PHE:CD1	2.55	0.41
2:K:89:ARG:HH22	2:K:112:ASP:CG	2.27	0.41
1:A:112:ILE:HD12	1:A:143:TYR:CD1	2.54	0.41
1:B:228:TRP:CB	1:B:232:LEU:HG	2.51	0.41
2:H:113:THR:HG23	2:H:146:THR:HA	2.02	0.41
1:D:30:SER:HB3	1:D:85:TRP:CD1	2.55	0.41
1:F:188:ILE:HG13	1:F:189:PHE:CD1	2.56	0.41
1:B:71:LEU:H	1:B:71:LEU:HG	1.73	0.41
2:H:89:ARG:HH22	2:H:112:ASP:CG	2.28	0.41
1:A:194:VAL:HG13	1:A:197:PRO:HB3	2.02	0.41
1:C:112:ILE:HD12	1:C:143:TYR:HD1	1.86	0.41
2:G:57:GLY:HA2	2:G:72:CYS:HA	2.03	0.41
2:H:148:SER:O	2:H:149:SER:OG	2.37	0.41
1:B:213:VAL:HG22	1:B:237:ILE:HB	2.03	0.41
1:B:232:LEU:HD23	1:B:232:LEU:HA	1.84	0.41
1:B:245:GLU:OE1	2:I:123:ARG:NH1	2.48	0.41
1:C:188:ILE:HD13	1:C:209:PRO:HG3	2.01	0.41
1:C:110:THR:O	1:C:178:GLN:NE2	2.54	0.40
1:C:215:THR:HA	1:C:216:GLY:HA2	1.84	0.40
2:L:34:VAL:HG12	2:L:35:GLN:O	2.20	0.40
1:A:217:GLY:HA2	1:A:218:PRO:HD3	1.93	0.40
1:D:194:VAL:HG13	1:D:197:PRO:HB3	2.03	0.40
1:E:159:PHE:CE2	1:E:161:GLN:HB2	2.55	0.40
1:E:215:THR:HA	1:E:216:GLY:HA2	1.84	0.40
2:G:36:PRO:HD2	2:G:149:SER:HB2	2.03	0.40
2:I:49:SER:O	2:I:99:ASN:ND2	2.53	0.40
2:I:53:ASP:OD1	2:I:53:ASP:N	2.41	0.40
2:K:54:TYR:O	2:K:94:ARG:NH2	2.53	0.40
1:A:74:ILE:O	1:A:79:PRO:HD3	2.22	0.40
1:A:93:GLN:HA	1:A:96:GLN:HE21	1.86	0.40
1:D:69:MET:HE2	1:D:69:MET:HB3	1.82	0.40
1:D:207:ARG:C	1:D:209:PRO:HD3	2.47	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	243/289 (84%)	228 (94%)	12 (5%)	3 (1%)	10	17
1	B	243/289 (84%)	228 (94%)	13 (5%)	2 (1%)	16	27
1	C	243/289 (84%)	229 (94%)	12 (5%)	2 (1%)	16	27
1	D	243/289 (84%)	230 (95%)	11 (4%)	2 (1%)	16	27
1	E	243/289 (84%)	230 (95%)	11 (4%)	2 (1%)	16	27
1	F	243/289 (84%)	228 (94%)	14 (6%)	1 (0%)	30	45
2	G	124/159 (78%)	120 (97%)	4 (3%)	0	100	100
2	H	124/159 (78%)	120 (97%)	4 (3%)	0	100	100
2	I	124/159 (78%)	120 (97%)	4 (3%)	0	100	100
2	J	124/159 (78%)	120 (97%)	4 (3%)	0	100	100
2	K	124/159 (78%)	120 (97%)	4 (3%)	0	100	100
2	L	124/159 (78%)	120 (97%)	4 (3%)	0	100	100
All	All	2202/2688 (82%)	2093 (95%)	97 (4%)	12 (0%)	24	39

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	99	SER
1	A	193	ARG
1	B	193	ARG
1	C	193	ARG
1	E	193	ARG
1	F	193	ARG
1	A	88	GLY
1	D	193	ARG
1	E	88	GLY
1	C	88	GLY
1	B	88	GLY
1	D	88	GLY

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	205/235 (87%)	200 (98%)	5 (2%)	43	65
1	B	205/235 (87%)	201 (98%)	4 (2%)	48	69
1	C	205/235 (87%)	204 (100%)	1 (0%)	81	91
1	D	205/235 (87%)	202 (98%)	3 (2%)	57	75
1	E	205/235 (87%)	203 (99%)	2 (1%)	68	81
1	F	205/235 (87%)	203 (99%)	2 (1%)	68	81
2	G	100/124 (81%)	100 (100%)	0	100	100
2	H	100/124 (81%)	100 (100%)	0	100	100
2	I	100/124 (81%)	99 (99%)	1 (1%)	68	81
2	J	100/124 (81%)	100 (100%)	0	100	100
2	K	100/124 (81%)	99 (99%)	1 (1%)	68	81
2	L	100/124 (81%)	100 (100%)	0	100	100
All	All	1830/2154 (85%)	1811 (99%)	19 (1%)	68	81

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	LEU
1	A	182	VAL
1	A	194	VAL
1	A	233	LYS
1	A	240	THR
1	B	71	LEU
1	B	145	GLN
1	B	182	VAL
1	B	194	VAL
1	C	194	VAL
1	D	137	GLN
1	D	194	VAL
1	D	240	THR
1	E	100	LEU

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Mol	Chain	Res	Type
1	E	194	VAL
1	F	89	ASN
1	F	194	VAL
2	I	34	VAL
2	K	34	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	GLN
1	A	96	GLN
1	A	149	GLN
1	A	256	GLN
1	A	260	ASN
1	B	67	GLN
1	B	145	GLN
1	B	178	GLN
1	B	220	GLN
1	C	67	GLN
1	C	128	GLN
1	D	67	GLN
1	D	137	GLN
1	D	149	GLN
1	D	178	GLN
1	E	220	GLN
1	E	256	GLN
1	E	260	ASN
1	F	198	GLN
2	G	104	GLN
2	G	106	ASN
2	J	104	GLN
2	J	106	ASN
2	L	35	GLN
2	L	104	GLN
2	L	106	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	$\# Z > 2$
3	GOL	E	301	-	5,5,5	0.35	0	5,5,5	0.32	0
3	GOL	B	301	-	5,5,5	0.38	0	5,5,5	0.35	0
3	GOL	H	201	-	5,5,5	0.37	0	5,5,5	0.32	0
3	GOL	D	301	-	5,5,5	0.35	0	5,5,5	0.34	0
3	GOL	F	301	-	5,5,5	0.38	0	5,5,5	0.39	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
3	GOL	E	301	-	-	2/4/4/4	-
3	GOL	B	301	-	-	2/4/4/4	-
3	GOL	H	201	-	-	2/4/4/4	-
3	GOL	D	301	-	-	2/4/4/4	-
3	GOL	F	301	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	301	GOL	O1-C1-C2-O2
3	B	301	GOL	O1-C1-C2-C3
3	D	301	GOL	O1-C1-C2-C3
3	H	201	GOL	O1-C1-C2-O2
3	H	201	GOL	O1-C1-C2-C3
3	E	301	GOL	O1-C1-C2-C3
3	D	301	GOL	O1-C1-C2-O2
3	E	301	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	301	GOL	1	0
3	F	301	GOL	1	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, 95th 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 ²)	Q < 0.9
1	A	245/289 (84%)	0.46	16 (6%) 25 19	55, 77, 118, 156	0
1	B	245/289 (84%)	0.46	14 (5%) 29 23	50, 83, 139, 185	0
1	C	245/289 (84%)	0.79	18 (7%) 21 15	65, 102, 168, 201	0
1	D	245/289 (84%)	0.34	10 (4%) 41 33	49, 74, 116, 166	0
1	E	245/289 (84%)	0.63	16 (6%) 25 19	56, 88, 136, 170	0
1	F	245/289 (84%)	0.97	35 (14%) 6 5	59, 101, 163, 212	0
2	G	126/159 (79%)	1.42	23 (18%) 3 2	94, 146, 180, 216	0
2	H	126/159 (79%)	0.90	15 (11%) 9 7	61, 95, 138, 159	0
2	I	126/159 (79%)	2.04	62 (49%) 0 0	78, 158, 221, 252	0
2	J	126/159 (79%)	0.76	10 (7%) 18 14	53, 86, 132, 165	0
2	K	126/159 (79%)	0.91	14 (11%) 10 8	69, 103, 148, 182	0
2	L	126/159 (79%)	1.67	41 (32%) 1 0	79, 114, 163, 186	0
All	All	2226/2688 (82%)	0.84	274 (12%) 8 7	49, 95, 167, 252	0

All (274) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	145	VAL	5.1
2	I	116	TYR	4.5
2	H	31	GLY	4.2
2	H	77	GLY	4.2
2	H	33	LEU	4.2
2	I	40	LEU	4.2
2	I	144	PRO	4.0
2	I	108	LEU	4.0
2	G	33	LEU	4.0
2	I	31	GLY	3.9
2	I	143	THR	3.9

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Mol	Chain	Res	Type	RSRZ
2	I	38	GLY	3.8
2	I	42	LEU	3.8
2	I	37	GLY	3.8
2	I	146	THR	3.7
1	A	217	GLY	3.6
2	I	88	GLY	3.6
2	L	32	GLY	3.6
2	I	34	VAL	3.6
1	F	167	PRO	3.6
2	K	31	GLY	3.5
2	I	90	PHE	3.5
1	F	215	THR	3.5
1	A	208	SER	3.4
2	I	118	CYS	3.4
1	C	218	PRO	3.3
2	L	33	LEU	3.3
2	I	30	GLY	3.2
2	L	118	CYS	3.2
2	I	41	ARG	3.2
2	K	87	LYS	3.2
2	I	112	ASP	3.2
2	I	46	ALA	3.2
2	J	133	SER	3.2
2	I	140	GLY	3.2
2	I	29	SER	3.2
2	L	87	LYS	3.1
1	F	88	GLY	3.1
2	L	31	GLY	3.1
1	E	90	ALA	3.1
1	F	151	ALA	3.1
1	A	153	LYS	3.1
2	L	77	GLY	3.1
2	G	121	ALA	3.1
2	I	36	PRO	3.0
2	I	70	VAL	3.0
1	E	228	TRP	3.0
2	G	32	GLY	3.0
2	H	87	LYS	3.0
2	I	45	ALA	3.0
2	I	115	VAL	3.0
2	L	58	TRP	2.9
2	H	78	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	F	164	ILE	2.9
2	L	34	VAL	2.9
1	E	153	LYS	2.9
1	A	232	LEU	2.9
2	J	33	LEU	2.9
1	F	220	GLN	2.8
1	F	212	ILE	2.8
2	G	142	GLY	2.8
1	F	224	ILE	2.8
1	A	216	GLY	2.8
1	B	232	LEU	2.8
2	I	27	VAL	2.8
2	L	133	SER	2.8
2	I	43	SER	2.8
2	K	49	SER	2.8
1	B	165	ASN	2.8
2	K	33	LEU	2.8
2	L	52	ASP	2.8
2	G	146	THR	2.8
2	L	41	ARG	2.8
2	I	63	PRO	2.7
1	F	149	GLN	2.7
2	K	142	GLY	2.7
1	C	164	ILE	2.7
2	I	57	GLY	2.7
1	A	90	ALA	2.7
1	B	215	THR	2.7
1	F	194	VAL	2.7
2	L	36	PRO	2.7
2	I	39	SER	2.7
1	E	194	VAL	2.7
2	G	145	VAL	2.7
2	I	62	ALA	2.7
2	L	115	VAL	2.7
1	D	191	ASP	2.7
1	C	178	GLN	2.7
2	I	117	TYR	2.7
1	A	89	ASN	2.7
1	E	66	TRP	2.7
2	I	58	TRP	2.7
1	C	154	PRO	2.7
1	F	241	SER	2.6

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Mol	Chain	Res	Type	RSRZ
2	I	99	ASN	2.6
1	B	152	ASP	2.6
1	E	191	ASP	2.6
2	I	105	MET	2.6
2	L	40	LEU	2.6
2	I	65	LYS	2.6
1	B	89	ASN	2.6
1	D	89	ASN	2.6
2	I	149	SER	2.6
1	F	152	ASP	2.6
2	L	51	LEU	2.6
2	I	87	LYS	2.6
2	H	142	GLY	2.6
2	J	74	GLY	2.6
1	F	196	TRP	2.6
1	F	228	TRP	2.6
2	I	113	THR	2.6
1	C	266	ASP	2.6
1	D	141	ASP	2.6
1	F	141	ASP	2.6
1	F	50	TYR	2.6
2	H	132	HIS	2.6
2	J	31	GLY	2.6
2	I	106	ASN	2.6
2	I	119	ALA	2.6
2	L	83	ALA	2.6
2	L	96	ASN	2.6
1	C	237	ILE	2.5
1	E	71	LEU	2.5
2	H	32	GLY	2.5
2	J	142	GLY	2.5
1	E	154	PRO	2.5
1	F	189	PHE	2.5
1	C	89	ASN	2.5
1	A	154	PRO	2.5
2	I	114	ALA	2.5
1	C	215	THR	2.5
1	F	169	THR	2.5
1	F	155	LYS	2.5
2	I	26	LEU	2.5
2	I	103	LEU	2.5
2	L	109	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
2	J	48	GLU	2.5
1	D	53	TYR	2.5
1	D	217	GLY	2.5
2	I	69	GLY	2.5
1	D	215	THR	2.4
2	L	103	LEU	2.4
2	I	148	SER	2.4
1	A	59	LYS	2.4
1	C	151	ALA	2.4
2	G	134	SER	2.4
2	G	52	ASP	2.4
2	K	118	CYS	2.4
2	L	44	CYS	2.4
2	I	28	GLU	2.4
1	F	153	LYS	2.4
2	L	90	PHE	2.4
1	E	215	THR	2.4
1	E	89	ASN	2.4
2	I	49	SER	2.4
2	I	25	GLN	2.4
2	I	147	VAL	2.4
2	K	133	SER	2.4
2	G	87	LYS	2.4
2	G	31	GLY	2.4
2	K	141	GLN	2.4
1	F	148	ALA	2.4
2	K	62	ALA	2.4
1	D	150	TYR	2.3
1	F	59	LYS	2.3
1	A	195	PRO	2.3
2	L	101	VAL	2.3
1	E	205	LEU	2.3
1	F	247	ALA	2.3
2	G	73	ILE	2.3
2	L	146	THR	2.3
2	H	133	SER	2.3
2	I	74	GLY	2.3
2	L	37	GLY	2.3
1	F	199	VAL	2.3
1	D	206	ALA	2.3
1	E	150	TYR	2.3
2	L	136	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	193	ARG	2.3
1	F	222	PRO	2.3
1	C	87	GLY	2.3
2	G	74	GLY	2.3
1	C	204	VAL	2.3
1	E	206	ALA	2.3
2	G	54	TYR	2.3
1	B	187	ASN	2.3
2	L	49	SER	2.3
1	C	262	LEU	2.3
2	I	33	LEU	2.3
2	K	115	VAL	2.3
1	F	150	TYR	2.2
2	L	99	ASN	2.2
2	I	109	ARG	2.2
1	B	154	PRO	2.2
1	B	191	ASP	2.2
1	B	88	GLY	2.2
2	G	34	VAL	2.2
1	F	233	LYS	2.2
2	H	96	ASN	2.2
2	H	35	GLN	2.2
2	I	126	GLY	2.2
2	G	108	LEU	2.2
1	E	59	LYS	2.2
1	F	182	VAL	2.2
2	I	86	VAL	2.2
1	A	215	THR	2.2
2	G	144	PRO	2.2
1	A	152	ASP	2.2
1	F	232	LEU	2.2
2	L	148	SER	2.2
1	B	228	TRP	2.2
1	E	50	TYR	2.2
2	H	117	TYR	2.2
2	K	65	LYS	2.2
1	C	232	LEU	2.2
2	G	26	LEU	2.2
1	F	188	ILE	2.2
2	L	139	TRP	2.1
2	H	48	GLU	2.1
1	B	193	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
2	I	104	GLN	2.1
2	L	54	TYR	2.1
1	B	217	GLY	2.1
2	I	142	GLY	2.1
2	L	88	GLY	2.1
1	F	265	VAL	2.1
2	K	132	HIS	2.1
1	C	201	ARG	2.1
1	B	59	LYS	2.1
1	C	155	LYS	2.1
1	D	190	LYS	2.1
1	F	154	PRO	2.1
2	J	47	PRO	2.1
2	L	108	LEU	2.1
1	E	68	GLY	2.1
2	J	77	GLY	2.1
2	G	107	ASP	2.1
2	L	39	SER	2.1
2	L	143	THR	2.1
1	A	233	LYS	2.1
2	L	63	PRO	2.1
2	G	117	TYR	2.1
2	L	117	TYR	2.1
2	L	126	GLY	2.1
2	G	27	VAL	2.1
1	F	208	SER	2.1
2	G	148	SER	2.1
1	C	173	GLU	2.1
2	K	36	PRO	2.1
1	A	266	ASP	2.0
1	C	152	ASP	2.0
2	I	60	ARG	2.0
2	H	134	SER	2.0
2	J	148	SER	2.0
2	L	100	THR	2.0
1	B	195	PRO	2.0
2	K	63	PRO	2.0
2	L	47	PRO	2.0
1	C	58	GLN	2.0
1	F	165	ASN	2.0
2	I	132	HIS	2.0
2	I	77	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
2	J	32	GLY	2.0
1	A	204	VAL	2.0
1	A	92	ARG	2.0
2	G	83	ALA	2.0
2	I	98	LYS	2.0
1	F	66	TRP	2.0
2	L	25	GLN	2.0
2	H	106	ASN	2.0
2	I	32	GLY	2.0
2	L	140	GLY	2.0
2	G	82	TYR	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 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, 95th 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 ²)	Q<0.9
3	GOL	B	301	6/6	0.84	0.18	93,100,103,108	0
3	GOL	F	301	6/6	0.84	0.19	124,126,129,130	0
3	GOL	E	301	6/6	0.85	0.16	73,76,82,85	0
3	GOL	H	201	6/6	0.87	0.13	70,77,82,92	0
3	GOL	D	301	6/6	0.91	0.12	66,68,70,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.