



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 20, 2026 – 01:19 AM UTC

PDB ID : 7DQC / pdb_00007dqc
Title : Crystal structure of nucleotide-free mutant A(S23C)3B(N64C)3 complex from *Enterococcus hirae* V-ATPase
Authors : Maruyama, S.; Suzuki, K.; Mizutani, K.; Imai, F.L.; Ishizuka-Katsura, Y.; Shirouzu, M.; Murata, M.
Deposited on : 2020-12-23
Resolution : 2.71 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

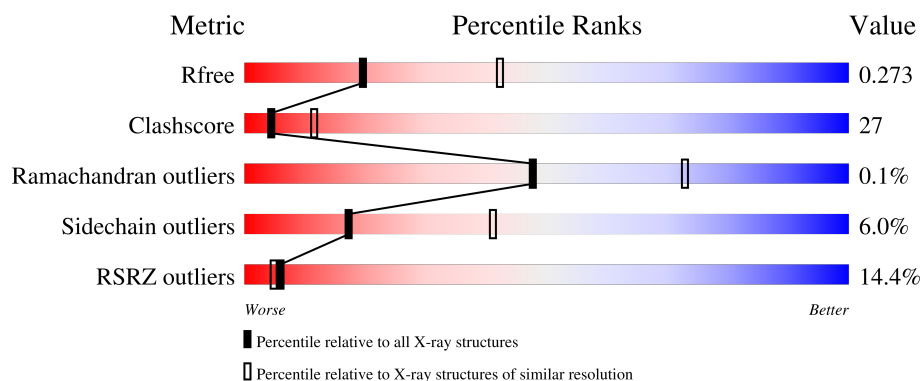
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	600	
1	B	600	
1	C	600	
2	D	465	
2	E	465	

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Mol	Chain	Length	Quality of chain
2	F	465	29% 47% 45% 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23860 atoms, of which 48 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 V-type sodium ATPase catalytic subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	586	Total	C	N	O	S	0	0	0
			4529	2847	761	895	26			
1	B	586	Total	C	N	O	S	0	0	0
			4501	2827	758	889	27			
1	C	584	Total	C	N	O	S	0	0	0
			4456	2802	746	881	27			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	GLY	-	expression tag	UNP Q08636
A	-5	SER	-	expression tag	UNP Q08636
A	-4	SER	-	expression tag	UNP Q08636
A	-3	GLY	-	expression tag	UNP Q08636
A	-2	SER	-	expression tag	UNP Q08636
A	-1	SER	-	expression tag	UNP Q08636
A	0	GLY	-	expression tag	UNP Q08636
A	23	CYS	SER	engineered mutation	UNP Q08636
B	-6	GLY	-	expression tag	UNP Q08636
B	-5	SER	-	expression tag	UNP Q08636
B	-4	SER	-	expression tag	UNP Q08636
B	-3	GLY	-	expression tag	UNP Q08636
B	-2	SER	-	expression tag	UNP Q08636
B	-1	SER	-	expression tag	UNP Q08636
B	0	GLY	-	expression tag	UNP Q08636
B	23	CYS	SER	engineered mutation	UNP Q08636
C	-6	GLY	-	expression tag	UNP Q08636
C	-5	SER	-	expression tag	UNP Q08636
C	-4	SER	-	expression tag	UNP Q08636
C	-3	GLY	-	expression tag	UNP Q08636
C	-2	SER	-	expression tag	UNP Q08636
C	-1	SER	-	expression tag	UNP Q08636
C	0	GLY	-	expression tag	UNP Q08636

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Chain	Residue	Modelled	Actual	Comment	Reference
C	23	CYS	SER	engineered mutation	UNP Q08636

- Molecule 2 is a protein called V-type sodium ATPase subunit B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	448	Total	C	N	O	S	0	0	0
			3380	2137	590	638	15			
2	E	450	Total	C	N	O	S	0	0	0
			3347	2112	582	640	13			
2	F	450	Total	C	N	O	S	0	0	0
			3277	2068	568	627	14			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-6	GLY	-	expression tag	UNP Q08637
D	-5	SER	-	expression tag	UNP Q08637
D	-4	SER	-	expression tag	UNP Q08637
D	-3	GLY	-	expression tag	UNP Q08637
D	-2	SER	-	expression tag	UNP Q08637
D	-1	SER	-	expression tag	UNP Q08637
D	0	GLY	-	expression tag	UNP Q08637
D	64	CYS	ASN	engineered mutation	UNP Q08637
E	-6	GLY	-	expression tag	UNP Q08637
E	-5	SER	-	expression tag	UNP Q08637
E	-4	SER	-	expression tag	UNP Q08637
E	-3	GLY	-	expression tag	UNP Q08637
E	-2	SER	-	expression tag	UNP Q08637
E	-1	SER	-	expression tag	UNP Q08637
E	0	GLY	-	expression tag	UNP Q08637
E	64	CYS	ASN	engineered mutation	UNP Q08637
F	-6	GLY	-	expression tag	UNP Q08637
F	-5	SER	-	expression tag	UNP Q08637
F	-4	SER	-	expression tag	UNP Q08637
F	-3	GLY	-	expression tag	UNP Q08637
F	-2	SER	-	expression tag	UNP Q08637
F	-1	SER	-	expression tag	UNP Q08637
F	0	GLY	-	expression tag	UNP Q08637
F	64	CYS	ASN	engineered mutation	UNP Q08637

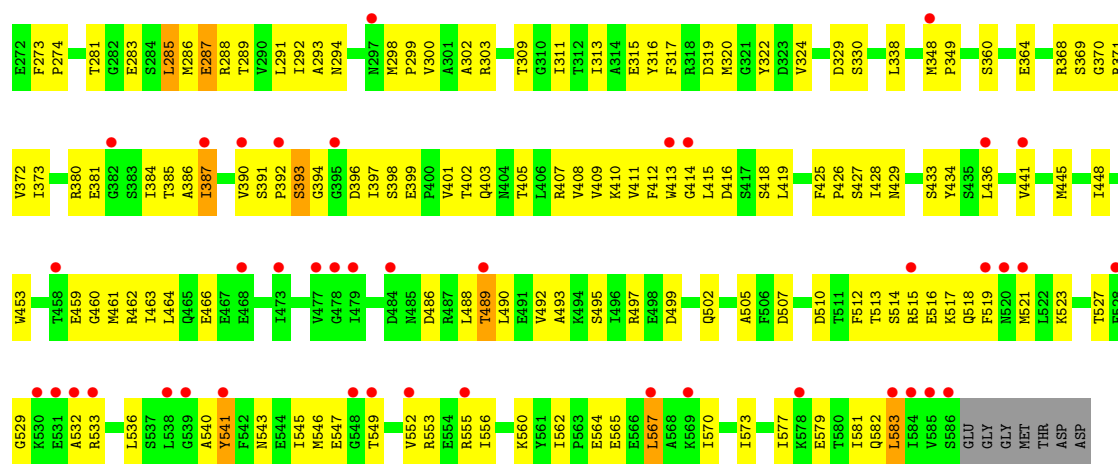
- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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
			14	3	8	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	103	Total	O	0	0
			103	103		
4	B	45	Total	O	0	0
			45	45		
4	D	34	Total	O	0	0
			34	34		
4	E	42	Total	O	0	0
			42	42		
4	C	43	Total	O	0	0
			43	43		
4	F	19	Total	O	0	0
			19	19		

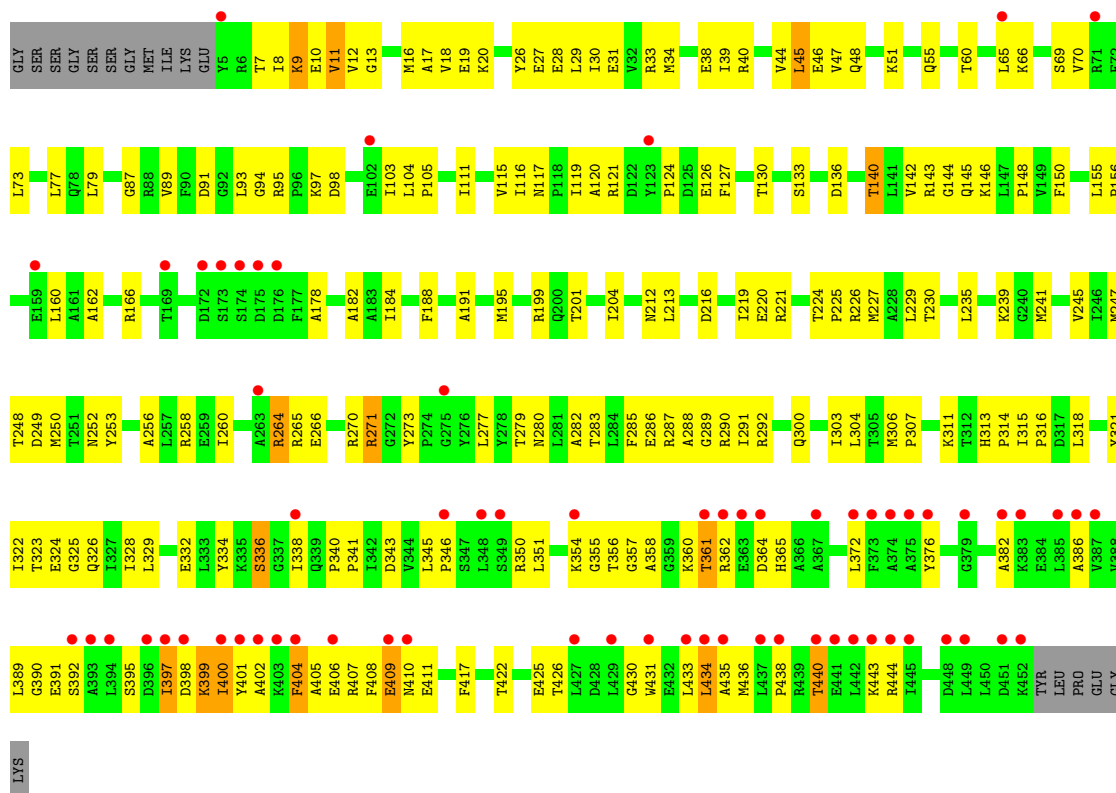


• Molecule 1: V-type sodium ATPase catalytic subunit A

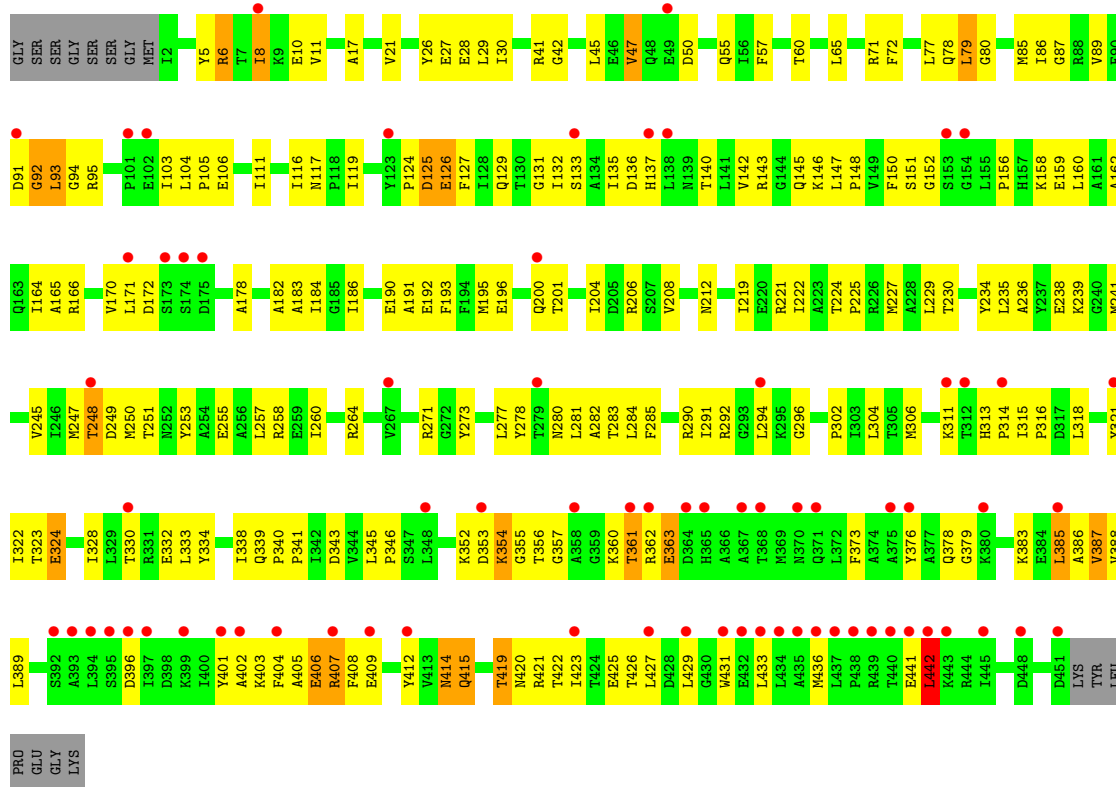


• Molecule 2: V-type sodium ATPase subunit B

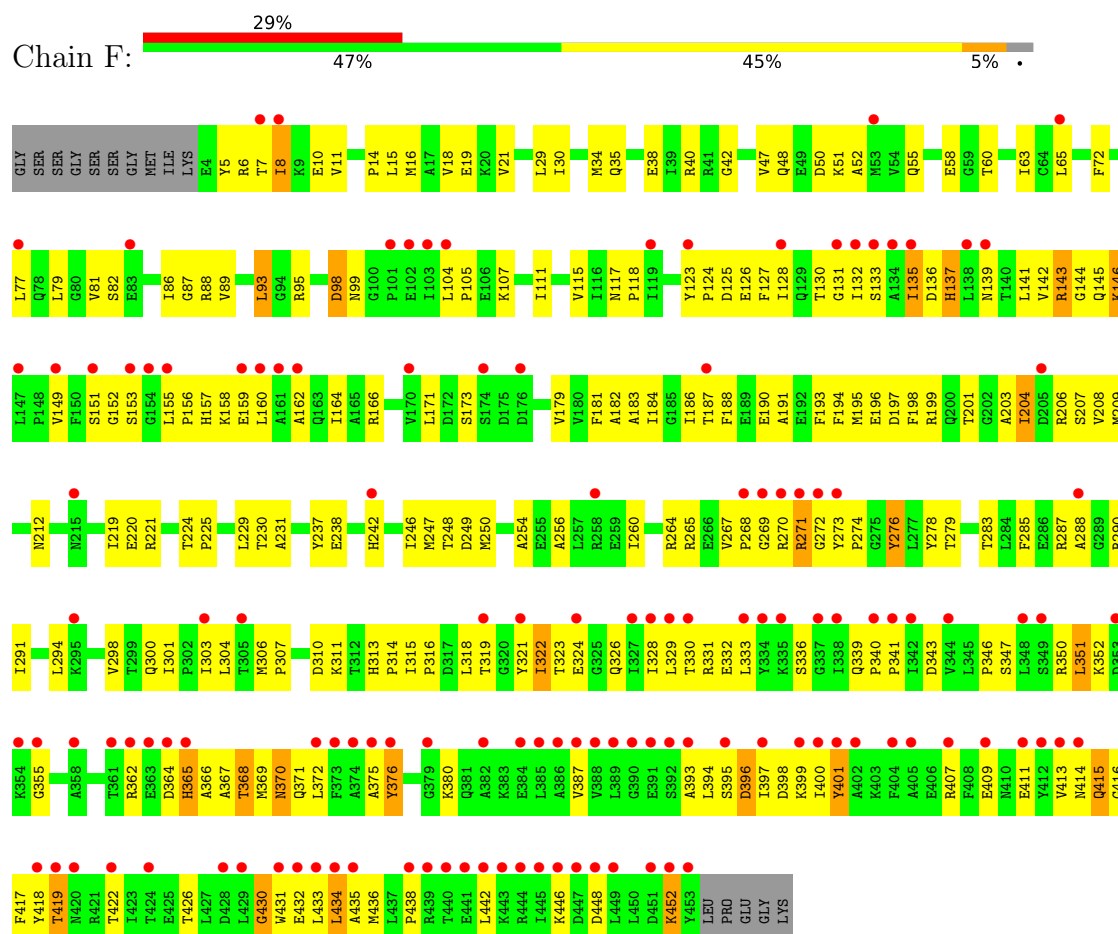




• Molecule 2: V-type sodium ATPase subunit B



● Molecule 2: V-type sodium ATPase subunit B



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	122.08Å 121.83Å 129.11Å 90.00° 90.22° 90.00°	Depositor
Resolution (Å)	44.30 – 2.71 44.30 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.30-2.71) 99.8 (44.30-2.71)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.234 , 0.266 0.245 , 0.273	Depositor DCC
R_{free} test set	5266 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	61.6	Xtriage
Anisotropy	0.155	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 72.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l 0.000 for -k,-h,-l 0.015 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	23860	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% 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	0.25	0/4605	0.71	4/6234 (0.1%)
1	B	0.26	0/4577	0.74	4/6199 (0.1%)
1	C	0.25	0/4532	0.72	1/6148 (0.0%)
2	D	0.26	0/3441	0.71	0/4658
2	E	0.28	0/3406	0.76	3/4616 (0.1%)
2	F	0.30	0/3336	0.77	2/4536 (0.0%)
All	All	0.27	0/23897	0.73	14/32391 (0.0%)

There are no bond length outliers.

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	399	GLU	CA-C-N	6.83	126.34	119.24
1	A	399	GLU	C-N-CA	6.83	126.34	119.24
1	B	230	VAL	CA-C-N	6.72	126.35	119.56
1	B	230	VAL	C-N-CA	6.72	126.35	119.56
2	E	92	GLY	N-CA-C	-5.95	107.53	115.32

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	4529	0	4462	160	0
1	B	4501	0	4407	237	0
1	C	4456	0	4329	230	0
2	D	3380	0	3279	195	0
2	E	3347	0	3194	198	0
2	F	3277	0	3037	254	0
3	A	36	48	48	3	0
4	A	103	0	0	5	0
4	B	45	0	0	2	0
4	C	43	0	0	0	0
4	D	34	0	0	4	0
4	E	42	0	0	6	0
4	F	19	0	0	1	0
All	All	23812	48	22756	1230	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 27.

The worst 5 of 1230 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)
2:F:267:VAL:HG22	2:F:268:PRO:HD2	1.31	1.10
2:F:155:LEU:HD22	2:F:156:PRO:HD2	1.25	1.09
1:A:27:MET:HE3	1:A:71:LEU:HB2	1.28	1.08
1:B:8:LYS:HD3	1:B:15:MET:HE2	1.40	1.00
1:B:260:GLY:HA3	1:B:261:GLU:HG2	1.39	1.00

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	584/600 (97%)	576 (99%)	8 (1%)	0	100	100
1	B	584/600 (97%)	573 (98%)	10 (2%)	1 (0%)	43	68
1	C	582/600 (97%)	569 (98%)	13 (2%)	0	100	100
2	D	446/465 (96%)	431 (97%)	15 (3%)	0	100	100
2	E	448/465 (96%)	427 (95%)	20 (4%)	1 (0%)	43	68
2	F	448/465 (96%)	431 (96%)	16 (4%)	1 (0%)	43	68
All	All	3092/3195 (97%)	3007 (97%)	82 (3%)	3 (0%)	48	73

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	387	VAL
1	B	287	GLU
2	E	387	VAL

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	491/511 (96%)	475 (97%)	16 (3%)	33	63
1	B	484/511 (95%)	463 (96%)	21 (4%)	26	54
1	C	474/511 (93%)	449 (95%)	25 (5%)	20	46
2	D	335/387 (87%)	313 (93%)	22 (7%)	15	36
2	E	325/387 (84%)	298 (92%)	27 (8%)	10	26
2	F	304/387 (79%)	270 (89%)	34 (11%)	6	15
All	All	2413/2694 (90%)	2268 (94%)	145 (6%)	17	41

5 of 145 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	F	93	LEU
2	F	434	LEU

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Mol	Chain	Res	Type
2	F	135	ILE
2	F	310	ASP
2	D	326	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
2	F	167	GLN
2	F	370	ASN
2	D	326	GLN
2	D	414	ASN
2	E	112	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	A	603	-	5,5,5	0.38	0	5,5,5	0.16	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	A	602	-	5,5,5	0.36	0	5,5,5	0.28	0
3	GOL	A	601	-	5,5,5	0.40	0	5,5,5	0.21	0
3	GOL	A	604	-	5,5,5	0.38	0	5,5,5	0.20	0
3	GOL	A	605	-	5,5,5	0.38	0	5,5,5	0.22	0
3	GOL	A	606	-	5,5,5	0.38	0	5,5,5	0.24	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	A	603	-	-	2/4/4/4	-
3	GOL	A	602	-	-	2/4/4/4	-
3	GOL	A	601	-	-	1/4/4/4	-
3	GOL	A	604	-	-	1/4/4/4	-
3	GOL	A	605	-	-	2/4/4/4	-
3	GOL	A	606	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	603	GOL	O1-C1-C2-C3
3	A	606	GOL	O1-C1-C2-C3
3	A	602	GOL	O1-C1-C2-C3
3	A	605	GOL	O1-C1-C2-C3
3	A	602	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	GOL	2	0
3	A	605	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 [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, 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	586/600 (97%)	0.28	21 (3%) 46 42	32, 52, 104, 139	0
1	B	586/600 (97%)	0.83	56 (9%) 13 11	42, 82, 137, 164	0
1	C	584/600 (97%)	1.18	99 (16%) 4 3	54, 87, 139, 158	0
2	D	448/465 (96%)	1.00	67 (14%) 5 4	48, 80, 152, 202	0
2	E	450/465 (96%)	0.91	72 (16%) 5 4	37, 72, 158, 209	0
2	F	450/465 (96%)	1.44	133 (29%) 1 1	51, 95, 172, 201	0
All	All	3104/3195 (97%)	0.92	448 (14%) 6 5	32, 77, 149, 209	0

The worst 5 of 448 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	396	ASP	5.8
1	C	545	ILE	5.2
1	C	479	ILE	4.9
2	E	137	HIS	4.8
2	E	393	ALA	4.7

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	A	605	6/6	0.65	0.23	113,136,138,138	0
3	GOL	A	601	6/6	0.71	0.27	114,136,139,139	0
3	GOL	A	603	6/6	0.76	0.26	101,121,125,125	0
3	GOL	A	604	6/6	0.77	0.19	77,93,95,95	0
3	GOL	A	602	6/6	0.86	0.20	88,106,108,108	0
3	GOL	A	606	6/6	0.90	0.21	91,109,112,112	0

6.5 Other polymers [i](#)

There are no such residues in this entry.