



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:05 AM UTC

PDB ID : 9UVR / pdb_00009uvr
Title : Crystal structure of SSA1632 from Streptococcus sanguinis
Authors : Takebe, K.; Suzuki, M.; Nakata, M.
Deposited on : 2025-05-11
Resolution : 3.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

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

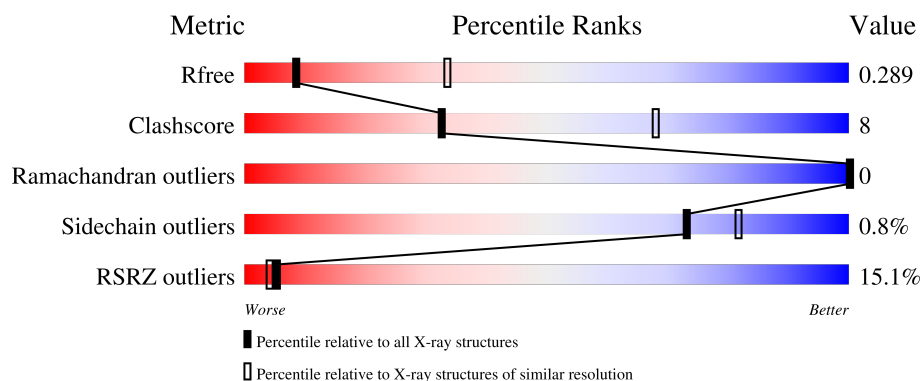
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	479	12% 71% 14% 15%
1	B	479	14% 72% 14% 14%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6227 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.

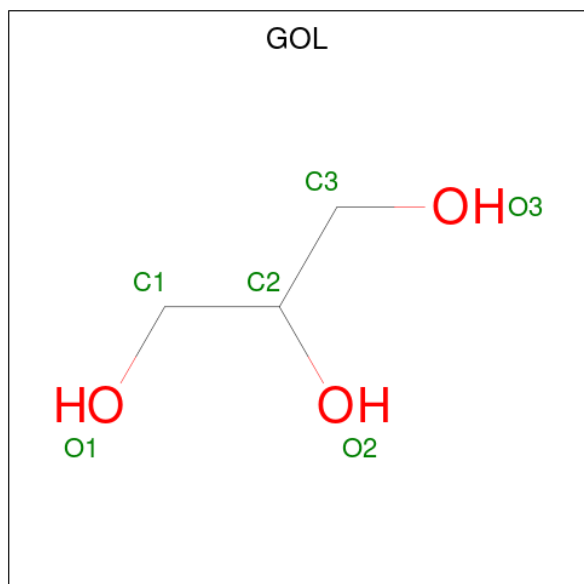
- Molecule 1 is a protein called Surface protein, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	409	Total	C	N	O	S	0	0	0
			3075	1931	506	634	4			
1	B	411	Total	C	N	O	S	0	0	0
			3092	1941	508	639	4			

- Molecule 2 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Cd	0	0
			3	3		
2	B	3	Total	Cd	0	0
			3	3		

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



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

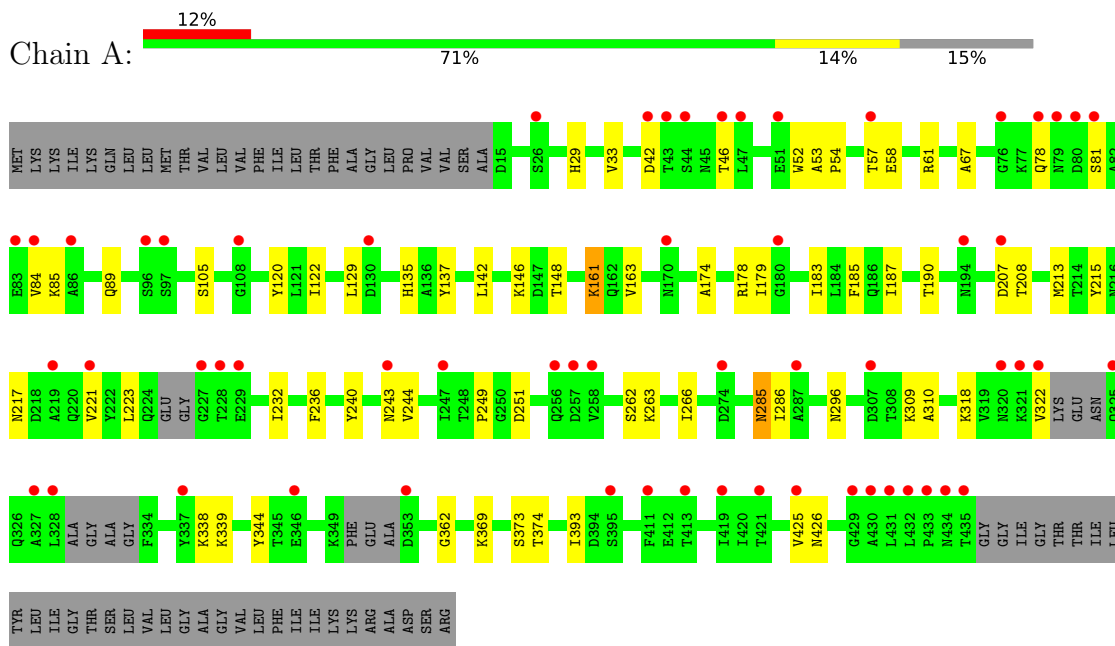
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	28	Total 28	O 28	0	0
4	B	14	Total 14	O 14	0	0

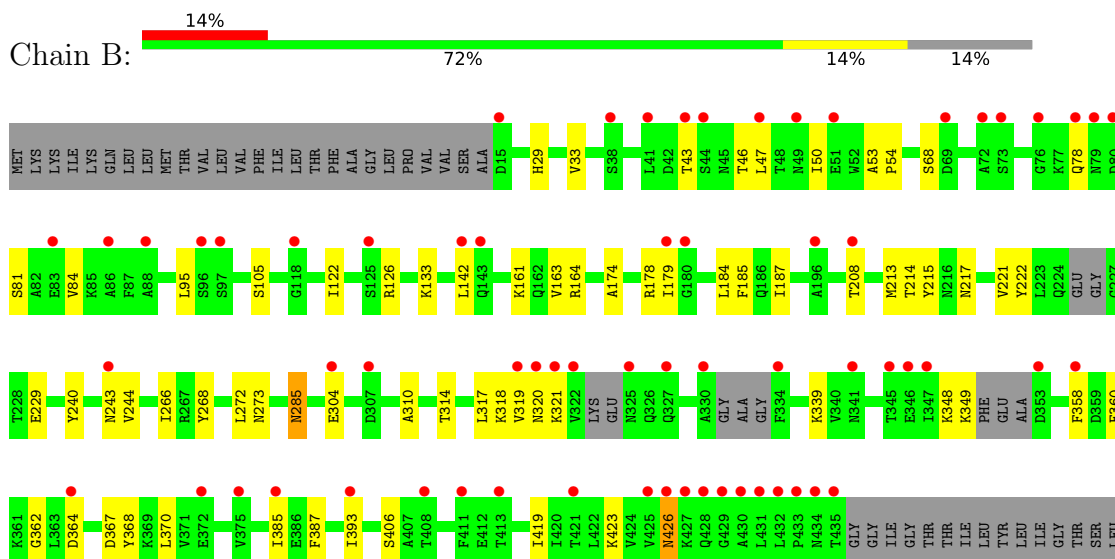
3 Residue-property plots

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: Surface protein, putative



- Molecule 1: Surface protein, putative



VAL
LEU
GLY
ALA
GLY
VAL
LEU
PHE
ILE
ILE
LYS
LYS
ARG
ALA
ASP
SER
ARG

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	42.03Å 64.12Å 92.98Å 78.46° 80.46° 75.77°	Depositor
Resolution (Å)	47.20 – 3.20 47.20 – 3.20	Depositor EDS
% Data completeness (in resolution range)	98.8 (47.20-3.20) 99.1 (47.20-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 3.19Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.278 , 0.289 0.278 , 0.289	Depositor DCC
R_{free} test set	734 reflections (3.23%)	wwPDB-VP
Wilson B-factor (Å ²)	48.8	Xtriage
Anisotropy	0.427	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	6227	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.98% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CD, 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	0.14	0/3126	0.32	0/4258
1	B	0.12	0/3143	0.34	0/4281
All	All	0.13	0/6269	0.33	0/8539

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

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	3075	0	2939	50	0
1	B	3092	0	2960	47	0
2	A	3	0	0	1	0
2	B	3	0	0	0	0
3	A	12	0	16	0	0
4	A	28	0	0	0	0
4	B	14	0	0	0	0
All	All	6227	0	5915	96	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 8.

All (96) 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:161:LYS:NZ	1:A:285:ASN:CG	1.67	1.50
1:B:321:LYS:NZ	1:B:426:ASN:CG	1.70	1.48
1:B:321:LYS:HZ1	1:B:426:ASN:CG	1.45	1.02
1:A:161:LYS:HZ3	1:A:285:ASN:CG	1.80	0.85
1:A:135:HIS:ND1	2:A:501:CD:CD	1.68	0.78
1:A:161:LYS:CE	1:A:285:ASN:CG	2.60	0.75
1:B:163:VAL:HG23	1:B:310:ALA:HB1	1.70	0.74
1:A:163:VAL:HG23	1:A:310:ALA:HB1	1.71	0.73
1:A:322:VAL:HG23	1:A:425:VAL:HA	1.70	0.72
1:B:84:VAL:HG13	1:B:142:LEU:HD12	1.72	0.71
1:A:84:VAL:HG13	1:A:142:LEU:HD12	1.75	0.68
1:A:208:THR:HG22	1:A:244:VAL:HG22	1.77	0.66
1:B:29:HIS:HB2	1:B:105:SER:HB2	1.77	0.66
1:A:251:ASP:OD2	1:B:126:ARG:NH2	2.31	0.63
1:A:161:LYS:HD3	1:A:187:ILE:HG12	1.81	0.62
1:B:43:THR:O	1:B:46:THR:OG1	2.18	0.61
1:A:81:SER:O	1:A:85:LYS:HG3	2.00	0.60
1:B:178:ARG:HH12	1:B:348:LYS:HE2	1.67	0.59
1:B:208:THR:HG22	1:B:244:VAL:HG22	1.85	0.59
1:B:370:LEU:HB3	1:B:385:ILE:HD11	1.85	0.59
1:B:240:TYR:CZ	1:B:243:ASN:HA	2.40	0.56
1:B:214:THR:HG23	1:B:273:ASN:HD22	1.70	0.56
1:A:185:PHE:HZ	1:A:213:MET:HE1	1.71	0.56
1:A:178:ARG:HG3	1:A:362:GLY:O	2.06	0.56
1:B:321:LYS:CE	1:B:426:ASN:CG	2.75	0.55
1:B:185:PHE:HZ	1:B:213:MET:HE1	1.72	0.54
1:B:78:GLN:O	1:B:81:SER:HB3	2.07	0.54
1:A:29:HIS:HB2	1:A:105:SER:HB3	1.90	0.54
1:A:161:LYS:NZ	1:A:285:ASN:CB	2.64	0.53
1:B:318:LYS:HE2	1:B:419:ILE:HG21	1.91	0.53
1:B:314:THR:CG2	1:B:362:GLY:HA2	2.38	0.53
1:B:178:ARG:HG3	1:B:362:GLY:O	2.09	0.52
1:A:53:ALA:HB1	1:A:54:PRO:HD2	1.92	0.52
1:A:207:ASP:OD1	1:A:285:ASN:OD1	2.28	0.52
1:A:33:VAL:HG22	1:A:122:ILE:HG12	1.93	0.51
1:B:164:ARG:HD2	1:B:184:LEU:HD23	1.93	0.51
1:B:321:LYS:HD2	1:B:426:ASN:CG	2.36	0.51
1:A:161:LYS:HZ3	1:A:285:ASN:CB	2.22	0.51
1:A:78:GLN:O	1:A:81:SER:HB3	2.11	0.50
1:B:53:ALA:HB1	1:B:54:PRO:HD2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:TYR:CZ	1:A:243:ASN:HA	2.46	0.50
1:A:221:VAL:HG22	1:A:266:ILE:HG12	1.93	0.50
1:B:349:LYS:H	1:B:349:LYS:HD2	1.77	0.49
1:B:319:VAL:HG22	1:B:358:PHE:HB2	1.95	0.49
1:B:164:ARG:HD3	1:B:184:LEU:HB3	1.94	0.49
1:B:268:TYR:OH	1:B:285:ASN:OD1	2.30	0.49
1:A:42:ASP:N	1:A:46:THR:O	2.47	0.48
1:A:286:ILE:HD11	1:A:309:LYS:HD3	1.96	0.48
1:A:338:LYS:HD3	1:A:339:LYS:N	2.28	0.47
1:B:387:PHE:HB3	1:B:406:SER:HB2	1.97	0.47
1:B:221:VAL:HG22	1:B:266:ILE:HG12	1.97	0.47
1:A:78:GLN:HG2	1:A:81:SER:HB3	1.97	0.47
1:B:222:TYR:HB3	1:B:229:GLU:HG3	1.97	0.47
1:A:236:PHE:CE2	1:A:249:PRO:HG3	2.50	0.46
1:B:179:ILE:HG13	1:B:362:GLY:HA3	1.96	0.46
1:B:317:LEU:HB3	1:B:360:PHE:HB2	1.97	0.46
1:A:179:ILE:HG13	1:A:362:GLY:HA3	1.99	0.45
1:B:321:LYS:CD	1:B:426:ASN:CG	2.88	0.45
1:A:146:LYS:HG3	1:A:148:THR:HB	1.98	0.45
1:A:33:VAL:HG12	1:A:120:TYR:CD2	2.51	0.45
1:B:179:ILE:HA	1:B:272:LEU:HB2	1.98	0.45
1:B:33:VAL:HG22	1:B:122:ILE:HG12	1.99	0.45
1:B:163:VAL:HG22	1:B:185:PHE:CD2	2.52	0.45
1:A:57:THR:O	1:A:61:ARG:HG3	2.18	0.44
1:A:344:TYR:CE2	1:A:369:LYS:HB2	2.52	0.44
1:A:58:GLU:HA	1:A:61:ARG:HD2	1.99	0.44
1:A:215:TYR:CE2	1:A:217:ASN:HA	2.52	0.44
1:A:89:GLN:HE22	1:A:296:ASN:HA	1.82	0.44
1:A:174:ALA:HB3	1:A:393:ILE:HG21	1.99	0.43
1:B:47:LEU:HD23	1:B:50:ILE:HD12	2.01	0.43
1:B:215:TYR:CE2	1:B:217:ASN:HA	2.53	0.43
1:B:320:ASN:O	1:B:423:LYS:HA	2.19	0.43
1:A:161:LYS:HG3	1:A:310:ALA:HB2	2.01	0.42
1:A:190:THR:HG22	1:A:263:LYS:HG2	2.01	0.42
1:B:161:LYS:HD2	1:B:187:ILE:HG12	2.00	0.42
1:A:163:VAL:CG2	1:A:310:ALA:HB1	2.47	0.42
1:B:178:ARG:HG2	1:B:179:ILE:N	2.33	0.42
1:A:78:GLN:HG2	1:A:81:SER:CB	2.48	0.42
1:A:163:VAL:HG22	1:A:185:PHE:CD2	2.55	0.42
1:A:318:LYS:HB2	1:A:318:LYS:HE3	1.80	0.42
1:A:373:SER:OG	1:A:374:THR:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:LEU:HB2	1:A:137:TYR:CZ	2.55	0.42
1:A:161:LYS:HZ3	1:A:285:ASN:HB3	1.85	0.41
1:A:183:ILE:HD12	1:A:213:MET:HE3	2.02	0.41
1:B:364:ASP:OD1	1:B:368:TYR:OH	2.38	0.41
1:A:53:ALA:HB1	1:A:54:PRO:CD	2.51	0.41
1:A:52:TRP:CD1	1:A:67:ALA:HB2	2.55	0.41
1:B:174:ALA:HB3	1:B:393:ILE:HG21	2.02	0.41
1:B:314:THR:HG23	1:B:362:GLY:HA2	2.03	0.41
1:A:232:ILE:HD12	1:A:236:PHE:HE1	1.86	0.41
1:A:223:LEU:HD11	1:A:262:SER:HB3	2.03	0.40
1:B:133:LYS:HG2	1:B:304:GLU:HG3	2.03	0.40
1:B:179:ILE:CD1	1:B:362:GLY:HA3	2.51	0.40
1:B:339:LYS:HB2	1:B:367:ASP:HB2	2.03	0.40
1:B:50:ILE:HB	1:B:68:SER:HA	2.04	0.40
1:B:95:LEU:HD23	1:B:95:LEU:HA	1.93	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	399/479 (83%)	387 (97%)	12 (3%)	0	100	100
1	B	401/479 (84%)	389 (97%)	12 (3%)	0	100	100
All	All	800/958 (84%)	776 (97%)	24 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	330/401 (82%)	327 (99%)	3 (1%)	70	81
1	B	333/401 (83%)	331 (99%)	2 (1%)	78	84
All	All	663/802 (83%)	658 (99%)	5 (1%)	73	82

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	161	LYS
1	A	285	ASN
1	A	426	ASN
1	B	285	ASN
1	B	426	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	ASN
1	A	216	ASN
1	A	320	ASN
1	B	66	ASN
1	B	220	GLN
1	B	342	ASN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 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	$\# Z > 2$
3	GOL	A	505	-	5,5,5	0.95	0	5,5,5	1.06	0
3	GOL	A	504	-	5,5,5	0.94	0	5,5,5	1.07	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	505	-	-	2/4/4/4	-
3	GOL	A	504	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	505	GOL	C1-C2-C3-O3
3	A	505	GOL	O2-C2-C3-O3

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	409/479 (85%)	1.16	59 (14%) 6 5	29, 44, 76, 138	0
1	B	411/479 (85%)	1.33	65 (15%) 5 4	29, 48, 98, 167	0
All	All	820/958 (85%)	1.25	124 (15%) 5 4	29, 45, 89, 167	0

All (124) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	435	THR	8.4
1	B	435	THR	8.4
1	B	434	ASN	5.8
1	B	432	LEU	5.6
1	B	431	LEU	5.6
1	A	434	ASN	5.3
1	B	322	VAL	4.7
1	A	431	LEU	4.5
1	B	433	PRO	4.4
1	B	83	GLU	4.4
1	B	41	LEU	4.3
1	A	320	ASN	4.1
1	B	78	GLN	4.0
1	B	430	ALA	4.0
1	B	80	ASP	3.9
1	A	432	LEU	3.9
1	B	320	ASN	3.8
1	A	325	GLN	3.7
1	A	227	GLY	3.6
1	B	411	PHE	3.4
1	B	325	ASN	3.4
1	B	429	GLY	3.4
1	A	256	GLN	3.4
1	B	79	ASN	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	96	SER	3.3
1	B	421	THR	3.3
1	A	411	PHE	3.2
1	B	353	ASP	3.2
1	A	430	ALA	3.2
1	A	79	ASN	3.2
1	B	428	GLN	3.2
1	A	97	SER	3.1
1	B	44	SER	3.1
1	B	334	PHE	3.1
1	B	425	VAL	3.1
1	A	78	GLN	3.1
1	B	346	GLU	3.1
1	A	433	PRO	3.1
1	A	429	GLY	3.0
1	B	180	GLY	3.0
1	A	80	ASP	2.9
1	A	257	ASP	2.9
1	A	307	ASP	2.9
1	A	108	GLY	2.9
1	A	42	ASP	2.8
1	B	330	ALA	2.8
1	B	413	THR	2.8
1	B	208	THR	2.8
1	A	180	GLY	2.8
1	A	46	THR	2.7
1	A	243	ASN	2.7
1	A	26	SER	2.7
1	A	83	GLU	2.6
1	B	97	SER	2.6
1	B	72	ALA	2.6
1	A	353	ASP	2.6
1	B	73	SER	2.6
1	B	76	GLY	2.6
1	A	57	THR	2.6
1	B	43	THR	2.6
1	B	408	THR	2.5
1	A	44	SER	2.5
1	A	421	THR	2.5
1	B	47	LEU	2.5
1	B	142	LEU	2.5
1	B	347	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	69	ASP	2.4
1	B	243	ASN	2.4
1	B	321	LYS	2.4
1	A	337	TYR	2.4
1	B	86	ALA	2.4
1	A	207	ASP	2.3
1	B	51	GLU	2.3
1	B	125	SER	2.3
1	B	385	ILE	2.3
1	B	327	GLN	2.3
1	B	118	GLY	2.3
1	B	375	VAL	2.3
1	A	395	SER	2.3
1	A	321	LYS	2.3
1	A	322	VAL	2.3
1	A	346	GLU	2.3
1	A	86	ALA	2.3
1	B	49	ASN	2.3
1	B	427	LYS	2.3
1	A	51	GLU	2.2
1	A	287	ALA	2.2
1	B	358	PHE	2.2
1	B	319	VAL	2.2
1	B	393	ILE	2.2
1	B	345	THR	2.2
1	B	426	ASN	2.2
1	B	372	GLU	2.2
1	A	130	ASP	2.2
1	A	328	LEU	2.2
1	A	43	THR	2.2
1	B	143	GLN	2.2
1	B	179	ILE	2.2
1	A	47	LEU	2.2
1	A	84	VAL	2.1
1	A	221	VAL	2.1
1	A	247	ILE	2.1
1	B	38	SER	2.1
1	A	229	GLU	2.1
1	B	304	GLU	2.1
1	A	170	ASN	2.1
1	A	194	ASN	2.1
1	B	96	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	274	ASP	2.1
1	A	258	VAL	2.1
1	B	364	ASP	2.1
1	B	88	ALA	2.1
1	B	196	ALA	2.1
1	B	341	ASN	2.1
1	A	228	THR	2.1
1	A	413	THR	2.1
1	A	327	ALA	2.1
1	A	425	VAL	2.0
1	A	81	SER	2.0
1	A	419	ILE	2.0
1	A	219	ALA	2.0
1	B	15	ASP	2.0
1	B	307	ASP	2.0
1	A	76	GLY	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	A	505	6/6	0.57	0.15	38,42,43,45	0
3	GOL	A	504	6/6	0.80	0.12	39,42,53,63	0
2	CD	A	503	1/1	0.80	0.12	51,51,51,51	0
2	CD	A	502	1/1	0.88	0.11	48,48,48,48	0
2	CD	A	501	1/1	0.88	0.09	36,36,36,36	0
2	CD	B	502	1/1	0.97	0.11	47,47,47,47	0
2	CD	B	503	1/1	0.98	0.07	36,36,36,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CD	B	501	1/1	0.99	0.06	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.