



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:45 AM UTC

PDB ID : 5C9S / pdb_00005c9s
Title : Crystal Structure of the C-terminal domain of Rrp5
Authors : Khoshnevis, S.; Karbstein, K.
Deposited on : 2015-06-28
Resolution : 2.70 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

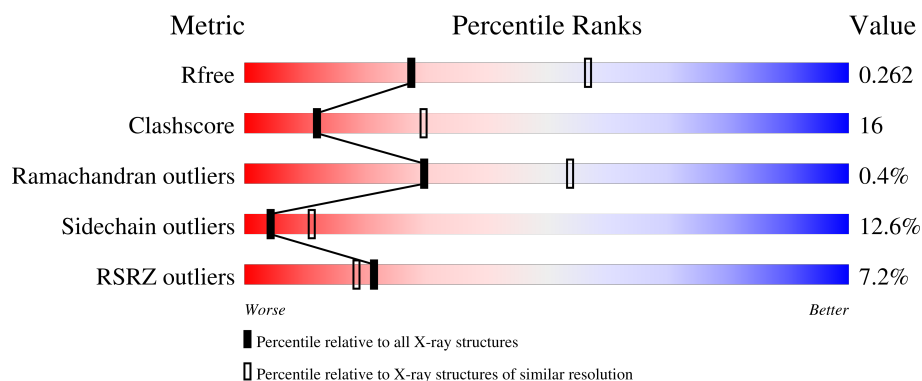
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	B	572	 3% 28% 15% . 52%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2317 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.

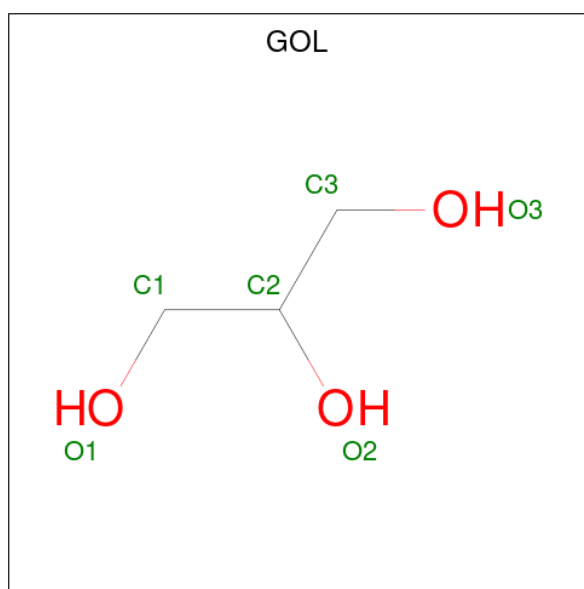
- Molecule 1 is a protein called rRNA biogenesis protein RRP5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	276	2258	1449	378	426	5	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1158	GLY	-	expression tag	UNP Q05022
B	1159	PRO	-	expression tag	UNP Q05022
B	1160	LEU	-	expression tag	UNP Q05022
B	1161	GLY	-	expression tag	UNP Q05022
B	1162	SER	-	expression tag	UNP Q05022

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

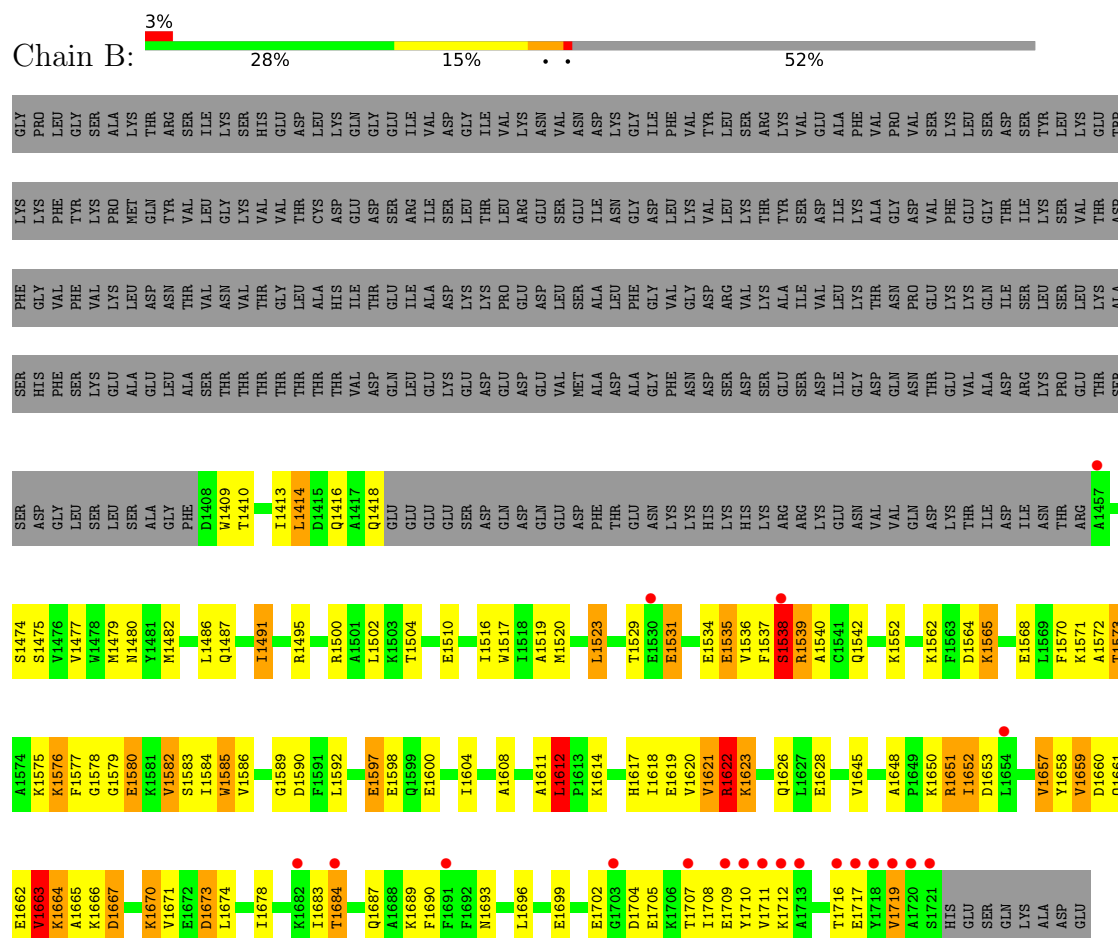
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	53	Total 53	O 53	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: rRNA biogenesis protein RRP5



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.14Å 114.14Å 67.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.67 – 2.70 43.67 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (43.67-2.70) 99.9 (43.67-2.70)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.7.1_743	Depositor
R, R_{free}	0.221 , 0.283 0.225 , 0.262	Depositor DCC
R_{free} test set	719 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.529	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.047 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2317	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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:
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	B	1.39	42/2302 (1.8%)	1.04	5/3097 (0.2%)

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1586	VAL	C-O	-7.38	1.15	1.24
1	B	1589	GLY	C-O	-7.34	1.15	1.23
1	B	1622	ARG	C-O	-7.02	1.16	1.24
1	B	1582	VAL	C-O	-6.97	1.15	1.24
1	B	1537	PHE	C-O	-6.83	1.16	1.24
1	B	1592	LEU	C-O	-6.76	1.16	1.24
1	B	1565	LYS	C-O	-6.74	1.16	1.24
1	B	1536	VAL	C-O	-6.58	1.16	1.24
1	B	1572	ALA	C-O	-6.38	1.16	1.24
1	B	1540	ALA	C-O	-6.15	1.16	1.24
1	B	1584	ILE	C-O	-6.15	1.16	1.24
1	B	1537	PHE	CA-C	-6.13	1.45	1.52
1	B	1619	GLU	C-O	-5.96	1.17	1.24
1	B	1611	ALA	CA-C	-5.93	1.44	1.52
1	B	1539	ARG	CA-C	-5.84	1.45	1.52
1	B	1664	LYS	C-O	-5.83	1.17	1.24
1	B	1539	ARG	C-O	-5.82	1.17	1.24
1	B	1621	VAL	C-O	-5.80	1.17	1.24
1	B	1570	PHE	C-O	-5.73	1.17	1.24
1	B	1620	VAL	C-O	-5.70	1.16	1.24
1	B	1658	TYR	C-O	-5.69	1.17	1.24
1	B	1612	LEU	C-N	5.66	1.40	1.33
1	B	1585	TRP	C-O	-5.66	1.17	1.24
1	B	1580	GLU	C-O	-5.54	1.17	1.24
1	B	1618	ILE	C-O	-5.48	1.17	1.24
1	B	1573	THR	C-O	-5.48	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1622	ARG	CA-C	-5.44	1.45	1.52
1	B	1623	LYS	C-O	-5.39	1.17	1.24
1	B	1568	GLU	C-O	-5.39	1.17	1.24
1	B	1652	ILE	C-O	-5.39	1.17	1.24
1	B	1660	ASP	C-O	-5.26	1.17	1.24
1	B	1590	ASP	C-O	-5.26	1.18	1.24
1	B	1661	GLN	CA-C	-5.25	1.46	1.52
1	B	1585	TRP	N-CA	-5.25	1.40	1.46
1	B	1583	SER	C-O	-5.23	1.17	1.24
1	B	1652	ILE	CA-C	-5.19	1.46	1.52
1	B	1661	GLN	C-O	-5.19	1.18	1.24
1	B	1659	VAL	C-O	-5.17	1.17	1.24
1	B	1621	VAL	N-CA	-5.11	1.40	1.46
1	B	1657	VAL	C-O	-5.11	1.18	1.24
1	B	1651	ARG	CA-C	-5.10	1.46	1.52
1	B	1564	ASP	C-O	-5.07	1.18	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1538	SER	N-CA-C	6.08	117.57	111.07
1	B	1664	LYS	N-CA-C	-5.39	107.96	114.75
1	B	1612	LEU	CA-C-N	-5.13	114.36	120.11
1	B	1612	LEU	C-N-CA	-5.13	114.36	120.11
1	B	1540	ALA	N-CA-C	5.03	119.67	111.37

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	B	2258	0	2248	71	0
2	B	6	0	8	2	0
3	B	53	0	0	0	0
All	All	2317	0	2256	71	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 16.

All (71) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1517:TRP:HA	1:B:1520:MET:HE3	1.32	1.08
1:B:1516:ILE:HG22	1:B:1520:MET:HE2	1.30	1.06
1:B:1659:VAL:O	1:B:1663:VAL:HG13	1.78	0.84
1:B:1653:ASP:O	1:B:1657:VAL:HG23	1.80	0.80
1:B:1689:LYS:O	1:B:1693:ASN:ND2	2.17	0.78
1:B:1622:ARG:NH1	1:B:1626:GLN:OE1	2.20	0.74
1:B:1502:LEU:HD11	1:B:1520:MET:HE1	1.69	0.72
1:B:1410:THR:HG21	1:B:1480:ASN:OD1	1.89	0.72
1:B:1651:ARG:HH22	2:B:1801:GOL:H32	1.57	0.69
1:B:1608:ALA:O	1:B:1612:LEU:HB2	1.93	0.69
1:B:1622:ARG:HG2	1:B:1622:ARG:HH11	1.59	0.68
1:B:1500:ARG:O	1:B:1504:THR:HG23	1.93	0.68
1:B:1617:HIS:O	1:B:1621:VAL:HG23	1.95	0.65
1:B:1578:GLY:O	1:B:1612:LEU:CD1	2.47	0.62
1:B:1517:TRP:CA	1:B:1520:MET:HE3	2.20	0.61
1:B:1667:ASP:OD1	1:B:1670:LYS:HB2	2.02	0.59
1:B:1705:GLU:HA	1:B:1708:ILE:HB	1.86	0.57
1:B:1579:GLY:O	1:B:1580:GLU:HG3	2.05	0.57
1:B:1409:TRP:CH2	1:B:1519:ALA:HB1	2.39	0.57
1:B:1662:GLU:C	1:B:1664:LYS:H	2.11	0.57
1:B:1716:THR:O	1:B:1719:VAL:HG12	2.05	0.56
1:B:1475:SER:O	1:B:1479:MET:HG3	2.06	0.55
1:B:1705:GLU:O	1:B:1709:GLU:HG2	2.06	0.55
1:B:1663:VAL:HG23	1:B:1663:VAL:O	2.06	0.54
1:B:1529:THR:OG1	1:B:1531:GLU:HG2	2.10	0.52
1:B:1622:ARG:NH1	1:B:1622:ARG:HG2	2.22	0.52
1:B:1707:THR:HA	1:B:1710:TYR:HB3	1.91	0.51
1:B:1683:ILE:HG23	1:B:1687:GLN:HB2	1.92	0.51
1:B:1539:ARG:O	1:B:1542:GLN:HB2	2.11	0.51
1:B:1534:GLU:O	1:B:1538:SER:HB2	2.12	0.50
1:B:1579:GLY:C	1:B:1580:GLU:HG3	2.37	0.50
1:B:1693:ASN:ND2	1:B:1693:ASN:H	2.09	0.50
1:B:1648:ALA:C	1:B:1650:LYS:H	2.19	0.50
1:B:1665:ALA:O	1:B:1666:LYS:HG2	2.11	0.50
1:B:1582:VAL:HG22	1:B:1612:LEU:HD21	1.95	0.49
1:B:1416:GLN:O	1:B:1552:LYS:NZ	2.38	0.49
1:B:1693:ASN:H	1:B:1693:ASN:HD22	1.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1598:GLU:OE2	1:B:1628:GLU:OE2	2.31	0.49
1:B:1652:ILE:O	1:B:1652:ILE:HG22	2.11	0.49
1:B:1673:ASP:N	1:B:1673:ASP:OD1	2.47	0.47
1:B:1707:THR:O	1:B:1711:VAL:HG23	2.14	0.47
1:B:1662:GLU:C	1:B:1664:LYS:N	2.73	0.47
1:B:1573:THR:O	1:B:1573:THR:HG22	2.15	0.47
1:B:1684:THR:HG22	1:B:1687:GLN:H	1.80	0.47
1:B:1674:LEU:O	1:B:1678:ILE:HG13	2.14	0.46
1:B:1659:VAL:HG12	1:B:1674:LEU:HD23	1.99	0.45
1:B:1585:TRP:CD2	1:B:1608:ALA:HB2	2.51	0.45
1:B:1659:VAL:HG12	1:B:1674:LEU:HG	1.98	0.45
1:B:1409:TRP:HH2	1:B:1519:ALA:HB1	1.78	0.45
1:B:1576:LYS:HD3	1:B:1577:PHE:CZ	2.51	0.45
1:B:1535:GLU:O	1:B:1539:ARG:HG3	2.17	0.44
1:B:1474:SER:OG	1:B:1477:VAL:HG23	2.18	0.43
1:B:1597:GLU:HB2	1:B:1600:GLU:HB2	2.01	0.43
1:B:1414:LEU:HD12	1:B:1414:LEU:HA	1.89	0.43
1:B:1479:MET:HA	1:B:1482:MET:HG2	2.01	0.42
1:B:1578:GLY:O	1:B:1612:LEU:HD13	2.19	0.42
1:B:1684:THR:O	1:B:1687:GLN:HB2	2.18	0.42
1:B:1699:GLU:HA	1:B:1702:GLU:HB2	2.00	0.42
1:B:1648:ALA:O	1:B:1650:LYS:N	2.51	0.42
1:B:1414:LEU:HD13	1:B:1479:MET:HE1	2.01	0.42
1:B:1651:ARG:HH22	2:B:1801:GOL:C3	2.28	0.42
1:B:1486:LEU:HD21	1:B:1523:LEU:CD2	2.50	0.42
1:B:1704:ASP:O	1:B:1708:ILE:HG13	2.20	0.42
1:B:1622:ARG:NH1	1:B:1626:GLN:CD	2.76	0.42
1:B:1665:ALA:C	1:B:1666:LYS:HG2	2.45	0.41
1:B:1491:ILE:HD11	1:B:1495:ARG:NH2	2.35	0.41
1:B:1667:ASP:O	1:B:1671:VAL:HG23	2.21	0.41
1:B:1696:LEU:HD12	1:B:1696:LEU:HA	1.88	0.41
1:B:1687:GLN:O	1:B:1690:PHE:HB3	2.21	0.40
1:B:1659:VAL:HG12	1:B:1674:LEU:CG	2.51	0.40
1:B:1659:VAL:HG12	1:B:1674:LEU:CD2	2.51	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	B	272/572 (48%)	261 (96%)	10 (4%)	1 (0%)	30	54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1663	VAL

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	B	238/502 (47%)	208 (87%)	30 (13%)	4	11

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1413	ILE
1	B	1414	LEU
1	B	1418	GLN
1	B	1487	GLN
1	B	1491	ILE
1	B	1510	GLU
1	B	1523	LEU
1	B	1531	GLU
1	B	1535	GLU
1	B	1538	SER

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Mol	Chain	Res	Type
1	B	1562	LYS
1	B	1565	LYS
1	B	1571	LYS
1	B	1575	LYS
1	B	1576	LYS
1	B	1597	GLU
1	B	1604	ILE
1	B	1612	LEU
1	B	1614	LYS
1	B	1622	ARG
1	B	1623	LYS
1	B	1645	VAL
1	B	1663	VAL
1	B	1667	ASP
1	B	1670	LYS
1	B	1673	ASP
1	B	1684	THR
1	B	1712	LYS
1	B	1717	GLU
1	B	1719	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	1471	ASN
1	B	1522	ASN
1	B	1656	ASN
1	B	1661	GLN
1	B	1693	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

1 ligand is 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	$\# Z > 2$
2	GOL	B	1801	-	5,5,5	0.31	0	5,5,5	0.40	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
2	GOL	B	1801	-	-	2/4/4/4	-

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
2	B	1801	GOL	O1-C1-C2-C3
2	B	1801	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1801	GOL	2	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	B	276/572 (48%)	0.34	20 (7%) 21 18	19, 32, 85, 96	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1711	VAL	4.0
1	B	1719	VAL	3.8
1	B	1713	ALA	2.9
1	B	1721	SER	2.8
1	B	1720	ALA	2.8
1	B	1710	TYR	2.8
1	B	1716	THR	2.7
1	B	1682	LYS	2.7
1	B	1530	GLU	2.6
1	B	1457	ALA	2.6
1	B	1712	LYS	2.6
1	B	1717	GLU	2.5
1	B	1709	GLU	2.5
1	B	1707	THR	2.5
1	B	1691	PHE	2.5
1	B	1684	THR	2.4
1	B	1718	TYR	2.3
1	B	1654	LEU	2.1
1	B	1703	GLY	2.1
1	B	1538	SER	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($\text{\AA}^2$)	Q<0.9
2	GOL	B	1801	6/6	0.85	0.13	42,44,51,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.