



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:55 AM UTC

PDB ID : 3SPA / pdb_00003spa
Title : Crystal Structure of Human Mitochondrial RNA Polymerase
Authors : Ringel, R.; Sologub, M.; Morozov, Y.I.; Litonin, D.; Cramer, P.; Temiakov, D.
Deposited on : 2011-07-01
Resolution : 2.50 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

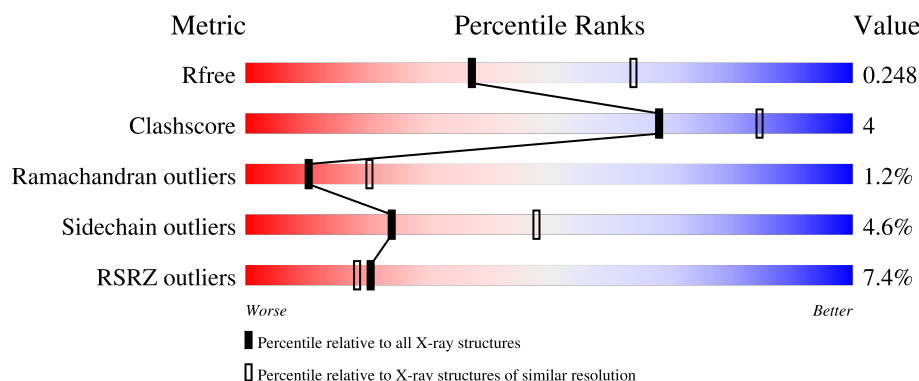
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	1134	6% 71% 11% • 17%
2	B	9	100%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	1	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7666 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 DNA-directed RNA polymerase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	941	Total	C	N	O	S	0	7	0
			7454	4745	1345	1315	49			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	97	MET	-	expression tag	UNP O00411
A	98	GLY	-	expression tag	UNP O00411
A	99	HIS	-	expression tag	UNP O00411
A	100	HIS	-	expression tag	UNP O00411
A	101	HIS	-	expression tag	UNP O00411
A	102	HIS	-	expression tag	UNP O00411
A	103	HIS	-	expression tag	UNP O00411
A	104	HIS	-	expression tag	UNP O00411
A	555	ALA	GLU	engineered mutation	UNP O00411

- Molecule 2 is a protein called Nonamer peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	9	Total	C	N	O	0	0	0
			45	27	9	9			

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

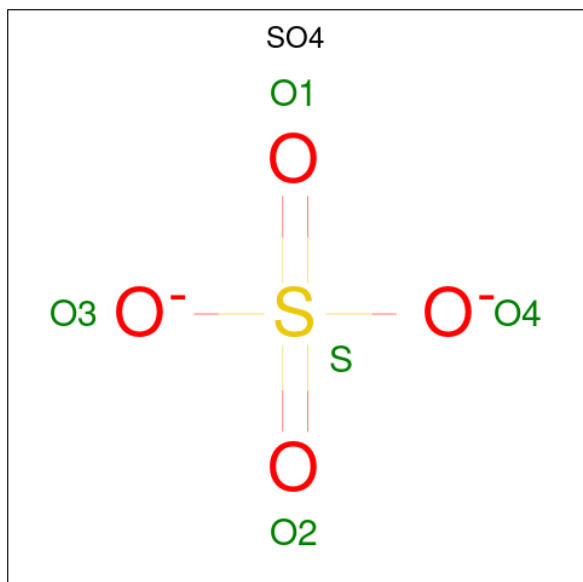


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total	Cl	0	0
			5	5		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		

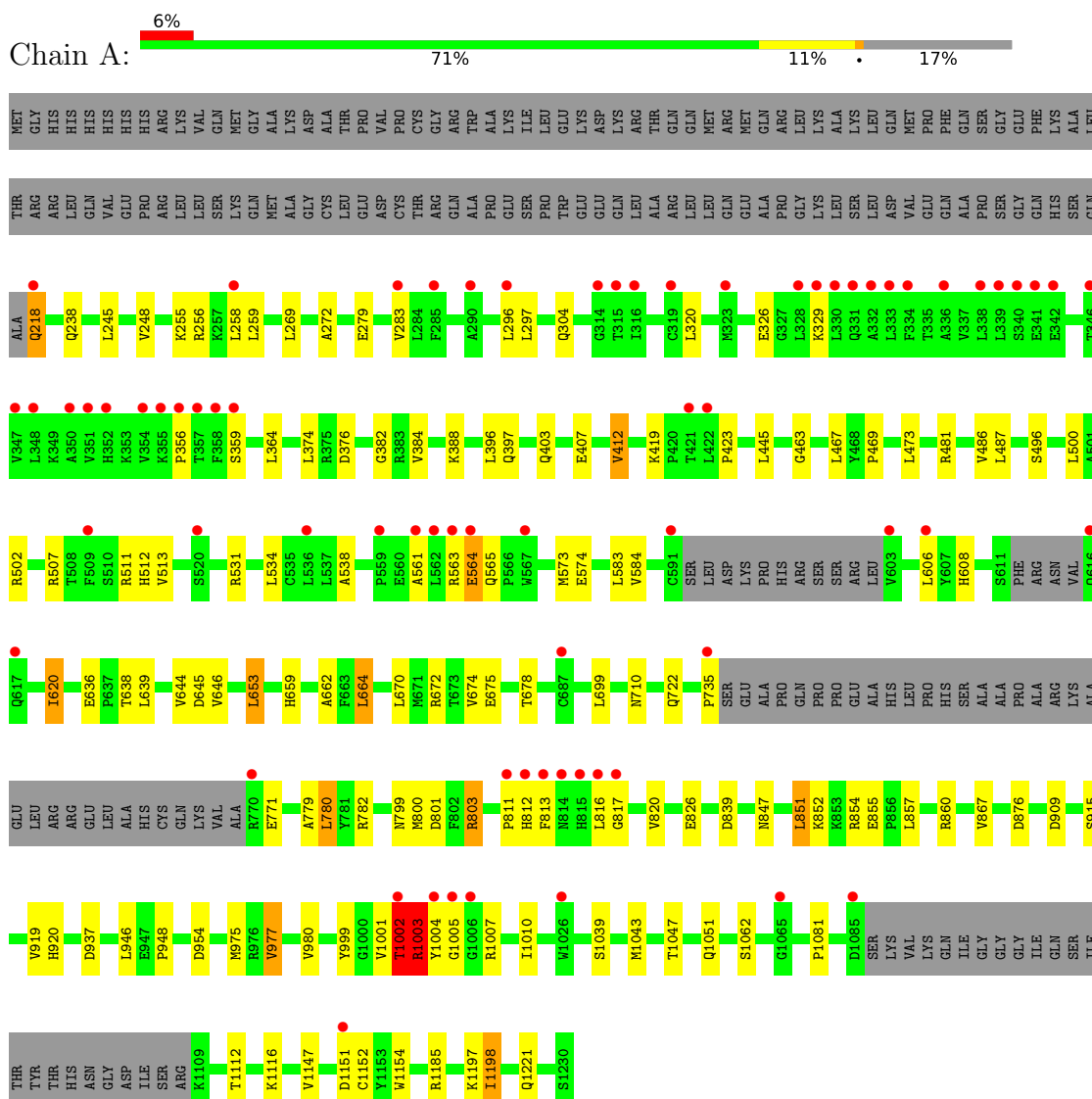
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	145	Total	O	0	0
			145	145		

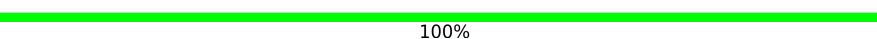
3 Residue-property plots

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: DNA-directed RNA polymerase, mitochondrial



- Molecule 2: Nonamer peptide

Chain B: 

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	211.26Å 211.26Å 60.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	58.13 – 2.50 58.13 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (58.13-2.50) 99.7 (58.13-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 2.51Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, R_{free}	0.185 , 0.231 0.213 , 0.248	Depositor DCC
R_{free} test set	2347 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	61.3	Xtriage
Anisotropy	0.549	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 92.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.013 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7666	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% 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: SO4, CL, 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	0.81	0/7651	1.46	42/10383 (0.4%)

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	851	LEU	N-CA-C	-9.41	98.39	110.53
1	A	1151	ASP	CA-CB-CG	7.86	120.46	112.60
1	A	326	GLU	N-CA-C	7.62	119.22	111.07
1	A	1152	CYS	N-CA-C	7.25	121.65	109.76
1	A	672	ARG	N-CA-C	7.13	121.21	112.23
1	A	852	LYS	N-CA-C	6.96	121.35	112.86
1	A	356	PRO	CA-C-N	6.47	133.35	121.70
1	A	356	PRO	C-N-CA	6.47	133.35	121.70
1	A	735	PRO	N-CA-CB	6.35	109.98	103.00
1	A	710	ASN	CA-C-N	6.13	126.78	119.98
1	A	710	ASN	C-N-CA	6.13	126.78	119.98
1	A	909	ASP	CA-CB-CG	6.13	118.73	112.60
1	A	1002	THR	CB-CA-C	5.96	122.27	110.42
1	A	937	ASP	CA-CB-CG	5.86	118.46	112.60
1	A	1198	ILE	N-CA-C	5.82	117.25	110.62
1	A	1001	VAL	CA-C-N	5.75	132.53	121.54
1	A	1001	VAL	C-N-CA	5.75	132.53	121.54
1	A	356	PRO	N-CA-CB	5.70	110.50	103.15
1	A	487	LEU	CA-C-N	5.66	127.87	120.28
1	A	487	LEU	C-N-CA	5.66	127.87	120.28
1	A	801	ASP	CA-CB-CG	5.62	118.22	112.60
1	A	678	THR	CA-C-N	5.44	127.57	120.28
1	A	678	THR	C-N-CA	5.44	127.57	120.28
1	A	839	ASP	CA-CB-CG	5.43	118.03	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1197	LYS	CA-C-N	5.40	128.15	120.53
1	A	1197	LYS	C-N-CA	5.40	128.15	120.53
1	A	256	ARG	CA-C-N	5.38	128.24	120.38
1	A	256	ARG	C-N-CA	5.38	128.24	120.38
1	A	376	ASP	CA-CB-CG	5.37	117.97	112.60
1	A	329	LYS	CA-C-N	5.35	127.39	120.44
1	A	329	LYS	C-N-CA	5.35	127.39	120.44
1	A	259	LEU	N-CA-C	5.27	116.84	108.67
1	A	999	TYR	N-CA-C	5.26	120.70	113.97
1	A	486	VAL	N-CA-C	-5.24	105.74	110.82
1	A	771	GLU	CA-C-N	5.14	127.17	120.28
1	A	771	GLU	C-N-CA	5.14	127.17	120.28
1	A	876	ASP	N-CA-C	5.11	116.94	111.36
1	A	855	GLU	N-CA-C	5.07	116.21	109.93
1	A	296	LEU	CA-C-N	5.04	127.04	120.28
1	A	296	LEU	C-N-CA	5.04	127.04	120.28
1	A	374	LEU	CA-C-N	5.01	127.30	120.54
1	A	374	LEU	C-N-CA	5.01	127.30	120.54

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	7454	0	7444	54	0
2	B	45	0	11	0	0
3	A	12	0	16	5	0
4	A	5	0	0	0	0
5	A	5	0	0	0	0
6	A	145	0	0	2	0
All	All	7666	0	7471	54	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (54) 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:218:GLN:O	1:A:218:GLN:HG3	1.77	0.78
1:A:1002:THR:HA	1:A:1003:ARG:O	1.93	0.67
1:A:1002:THR:H	1:A:1003:ARG:HG2	1.60	0.67
1:A:1039:SER:O	1:A:1043:MET:HG2	1.98	0.63
1:A:779:ALA:HB1	3:A:1:GOL:H32	1.82	0.60
1:A:507:ARG:HB3	1:A:511:ARG:HH21	1.68	0.59
1:A:1147:VAL:HG23	1:A:1154:TRP:HB2	1.84	0.58
1:A:502:ARG:HG3	1:A:573:MET:HE1	1.86	0.56
1:A:397:GLN:HG2	1:A:531:ARG:HG2	1.87	0.56
1:A:248:VAL:HG11	1:A:463:GLY:HA3	1.87	0.56
1:A:218:GLN:O	1:A:218:GLN:CG	2.49	0.56
1:A:564:GLU:HG3	1:A:659:HIS:CD2	2.41	0.55
1:A:854[A]:ARG:HD3	1:A:954:ASP:HB2	1.88	0.54
1:A:1051[B]:GLN:NE2	6:A:1295:HOH:O	2.33	0.53
1:A:645:ASP:HB3	1:A:812:HIS:HD2	1.74	0.53
1:A:467:LEU:HD11	1:A:574:GLU:HB3	1.90	0.52
1:A:653:LEU:HD13	1:A:664:LEU:HD12	1.91	0.52
1:A:699:LEU:HD11	1:A:800:MET:HG3	1.92	0.52
1:A:238:GLN:HE22	1:A:469:PRO:HA	1.75	0.52
1:A:1002:THR:HA	1:A:1003:ARG:C	2.36	0.51
1:A:255:LYS:HA	1:A:258:LEU:HD13	1.93	0.50
1:A:803:ARG:HH22	1:A:1003:ARG:HH12	1.59	0.50
1:A:584:VAL:HG22	1:A:606:LEU:HB2	1.93	0.50
1:A:407:GLU:OE2	1:A:664:LEU:HB2	2.13	0.49
1:A:513:VAL:HG22	1:A:563:ARG:HG3	1.94	0.49
1:A:218:GLN:HA	6:A:1271:HOH:O	2.13	0.48
1:A:1007:ARG:HA	1:A:1010:ILE:HD12	1.95	0.48
1:A:255:LYS:O	1:A:258:LEU:HB2	2.13	0.48
1:A:388:LYS:HD2	1:A:538:ALA:O	2.13	0.48
1:A:847:ASN:HD21	1:A:860:ARG:HH11	1.60	0.47
1:A:675:GLU:HG2	1:A:1081:PRO:HB2	1.95	0.47
1:A:412:VAL:HG21	1:A:646:VAL:HG21	1.96	0.47
1:A:272:ALA:HB1	1:A:304:GLN:HB3	1.97	0.47
1:A:396:LEU:HB3	1:A:534:LEU:HD22	1.97	0.47
1:A:419:LYS:HE3	1:A:722:GLN:HB3	1.97	0.46
1:A:645:ASP:HA	1:A:811:PRO:HD2	1.96	0.46
1:A:403:GLN:HG3	1:A:664:LEU:HD22	1.97	0.46
1:A:473:LEU:HD22	1:A:512:HIS:CE1	2.51	0.46
1:A:820:VAL:HG13	3:A:1:GOL:H2	1.98	0.45
1:A:496:SER:HA	1:A:620:ILE:HA	1.99	0.45
1:A:279:GLU:O	1:A:283:VAL:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:782:ARG:HG3	1:A:813:PHE:HD1	1.82	0.45
1:A:639:LEU:HD11	1:A:780:LEU:HD13	1.98	0.45
1:A:645:ASP:HB3	1:A:812:HIS:CD2	2.52	0.44
1:A:670:LEU:HD11	1:A:699:LEU:HD12	1.99	0.44
1:A:826:GLU:OE2	3:A:2:GOL:H12	2.18	0.43
1:A:445:LEU:HD21	1:A:583:LEU:HA	2.02	0.42
1:A:779:ALA:CB	3:A:1:GOL:H32	2.49	0.42
1:A:563:ARG:HA	1:A:563:ARG:HD2	1.91	0.41
1:A:1062:SER:OG	1:A:1112:THR:HG22	2.20	0.41
1:A:948:PRO:HD3	1:A:1221:GLN:HB3	2.02	0.41
1:A:975:MET:HG2	1:A:977:VAL:HG23	2.01	0.41
1:A:1047:THR:O	1:A:1051[A]:GLN:HB2	2.20	0.41
1:A:820:VAL:CG1	3:A:1:GOL:H2	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	A	938/1134 (83%)	891 (95%)	36 (4%)	11 (1%)	10	20

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	359	SER
1	A	561	ALA
1	A	816	LEU
1	A	817	GLY
1	A	1002	THR
1	A	1003	ARG
1	A	382	GLY
1	A	662	ALA

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Mol	Chain	Res	Type
1	A	565	GLN
1	A	1005	GLY
1	A	423	PRO

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	795/982 (81%)	759 (96%)	36 (4%)	24	49

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	218	GLN
1	A	245	LEU
1	A	269	LEU
1	A	297	LEU
1	A	320	LEU
1	A	364	LEU
1	A	384	VAL
1	A	412	VAL
1	A	481	ARG
1	A	500	LEU
1	A	564	GLU
1	A	608	HIS
1	A	620	ILE
1	A	636	GLU
1	A	638	THR
1	A	644	VAL
1	A	653	LEU
1	A	664	LEU
1	A	674	VAL
1	A	780	LEU
1	A	799	ASN
1	A	803	ARG
1	A	851	LEU

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Mol	Chain	Res	Type
1	A	857	LEU
1	A	867	VAL
1	A	915	SER
1	A	919	VAL
1	A	920	HIS
1	A	946	LEU
1	A	977	VAL
1	A	980	VAL
1	A	1003	ARG
1	A	1004	TYR
1	A	1116	LYS
1	A	1185	ARG
1	A	1198	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	238	GLN
1	A	244	HIS
1	A	252	GLN
1	A	390	HIS
1	A	485	GLN
1	A	608	HIS
1	A	616	GLN
1	A	799	ASN
1	A	812	HIS
1	A	829	GLN
1	A	847	ASN
1	A	903	ASN
1	A	920	HIS
1	A	962	GLN
1	A	1041	GLN
1	A	1170	GLN
1	A	1209	GLN

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](#)

Of 8 ligands modelled in this entry, 5 are monoatomic - leaving 3 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	2	-	5,5,5	0.53	0	5,5,5	0.39	0
3	GOL	A	1	-	5,5,5	0.75	0	5,5,5	1.09	0
5	SO4	A	1233	-	4,4,4	0.46	0	6,6,6	0.43	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	2	-	-	0/4/4/4	-
3	GOL	A	1	-	-	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
3	A	1	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	A	1	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2	GOL	1	0
3	A	1	GOL	4	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	941/1134 (82%)	0.45	70 (7%) 20 18	32, 82, 144, 194	7 (0%)
2	B	0/9	-	-	-	-
All	All	941/1143 (82%)	0.45	70 (7%) 20 18	32, 82, 144, 194	7 (0%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	815	HIS	9.3
1	A	1085	ASP	6.4
1	A	562	LEU	5.4
1	A	563	ARG	5.4
1	A	340	SER	5.1
1	A	1151	ASP	4.7
1	A	218	GLN	4.5
1	A	816	LEU	4.4
1	A	350	ALA	4.1
1	A	561	ALA	4.0
1	A	332	ALA	4.0
1	A	520	SER	3.9
1	A	735	PRO	3.9
1	A	567	TRP	3.8
1	A	351	VAL	3.6
1	A	559	PRO	3.6
1	A	358	PHE	3.6
1	A	813	PHE	3.5
1	A	812	HIS	3.5
1	A	1005	GLY	3.4
1	A	591	CYS	3.3
1	A	603	VAL	3.2
1	A	356	PRO	3.2
1	A	814	ASN	3.1
1	A	1006	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	285	PHE	3.0
1	A	357	THR	2.9
1	A	811	PRO	2.9
1	A	359	SER	2.8
1	A	354	VAL	2.8
1	A	1065	GLY	2.7
1	A	348	LEU	2.7
1	A	346	THR	2.7
1	A	341	GLU	2.6
1	A	331	GLN	2.6
1	A	323	MET	2.6
1	A	339	LEU	2.6
1	A	421	THR	2.6
1	A	333	LEU	2.6
1	A	347	VAL	2.5
1	A	329	LYS	2.4
1	A	1026	TRP	2.4
1	A	330	LEU	2.4
1	A	564	GLU	2.4
1	A	352	HIS	2.3
1	A	509	PHE	2.3
1	A	617	GLN	2.3
1	A	422	LEU	2.3
1	A	342	GLU	2.3
1	A	314	GLY	2.3
1	A	1002	THR	2.3
1	A	328	LEU	2.3
1	A	283	VAL	2.3
1	A	315	THR	2.2
1	A	616	GLN	2.2
1	A	338	LEU	2.2
1	A	536	LEU	2.2
1	A	258	LEU	2.2
1	A	296	LEU	2.2
1	A	290	ALA	2.2
1	A	336	ALA	2.2
1	A	687	CYS	2.2
1	A	817	GLY	2.2
1	A	316	ILE	2.1
1	A	1004	TYR	2.1
1	A	319	CYS	2.1
1	A	606	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	334	PHE	2.0
1	A	770	ARG	2.0
1	A	355	LYS	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
4	CL	A	3	1/1	0.77	0.16	122,122,122,122	0
3	GOL	A	2	6/6	0.84	0.12	107,108,109,109	0
4	CL	A	1232	1/1	0.87	0.11	102,102,102,102	0
3	GOL	A	1	6/6	0.88	0.16	85,86,86,87	0
4	CL	A	5	1/1	0.88	0.08	97,97,97,97	0
4	CL	A	4	1/1	0.90	0.11	121,121,121,121	0
4	CL	A	1231	1/1	0.93	0.13	113,113,113,113	0
5	SO4	A	1233	5/5	0.97	0.08	73,77,79,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.