



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:25 PM UTC

PDB ID : 3IPK / pdb_00003ipk
Title : Crystal Structure of A3VP1 of AgI/II of Streptococcus mutans
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Deposited on : 2009-08-17
Resolution : 2.04 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

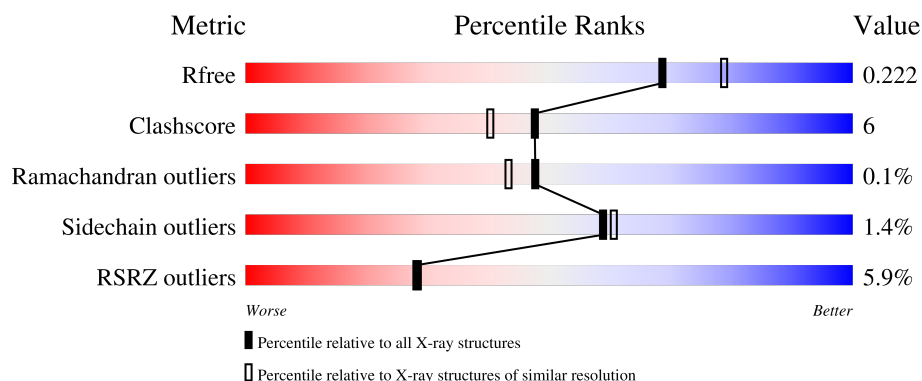
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.04 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	497	 7% 85% 13% ..
1	B	497	 5% 85% 12% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8564 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.

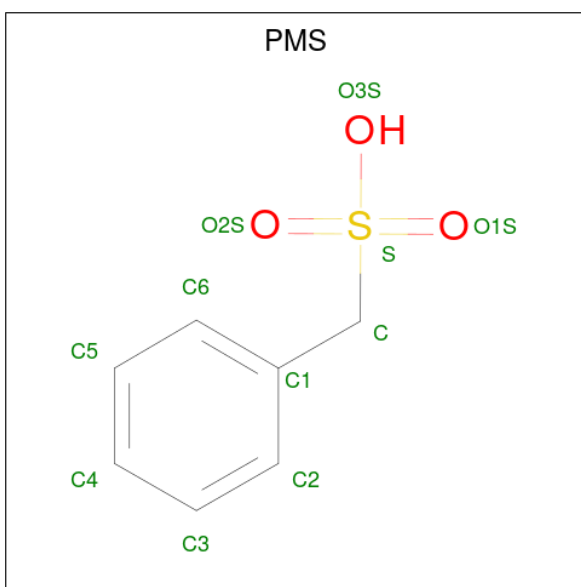
- Molecule 1 is a protein called AgI/II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	489	Total	C	N	O	S	0	0	0
			3760	2364	629	760	7			
1	B	489	Total	C	N	O	S	0	0	0
			3760	2364	629	760	7			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	398	ALA	VAL	SEE REMARK 999	UNP A8R5D9
A	782	TYR	HIS	SEE REMARK 999	UNP A8R5D9
A	875	LEU	-	expression tag	UNP A8R5D9
A	876	GLU	-	expression tag	UNP A8R5D9
A	877	HIS	-	expression tag	UNP A8R5D9
A	878	HIS	-	expression tag	UNP A8R5D9
A	879	HIS	-	expression tag	UNP A8R5D9
A	880	HIS	-	expression tag	UNP A8R5D9
A	881	HIS	-	expression tag	UNP A8R5D9
A	882	HIS	-	expression tag	UNP A8R5D9
B	398	ALA	VAL	SEE REMARK 999	UNP A8R5D9
B	782	TYR	HIS	SEE REMARK 999	UNP A8R5D9
B	875	LEU	-	expression tag	UNP A8R5D9
B	876	GLU	-	expression tag	UNP A8R5D9
B	877	HIS	-	expression tag	UNP A8R5D9
B	878	HIS	-	expression tag	UNP A8R5D9
B	879	HIS	-	expression tag	UNP A8R5D9
B	880	HIS	-	expression tag	UNP A8R5D9
B	881	HIS	-	expression tag	UNP A8R5D9
B	882	HIS	-	expression tag	UNP A8R5D9

- Molecule 2 is phenylmethanesulfonic acid (CCD ID: PMS) (formula: C₇H₈O₃S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	S	0	0
			11	7	3	1		
2	B	1	Total	C	O	S	0	0
			11	7	3	1		

- Molecule 3 is SULFATE ION (CCD ID: SO₄) (formula: O₄S).



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

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Ca 1	0	0
4	B	1	Total 1	Ca 1	0	0

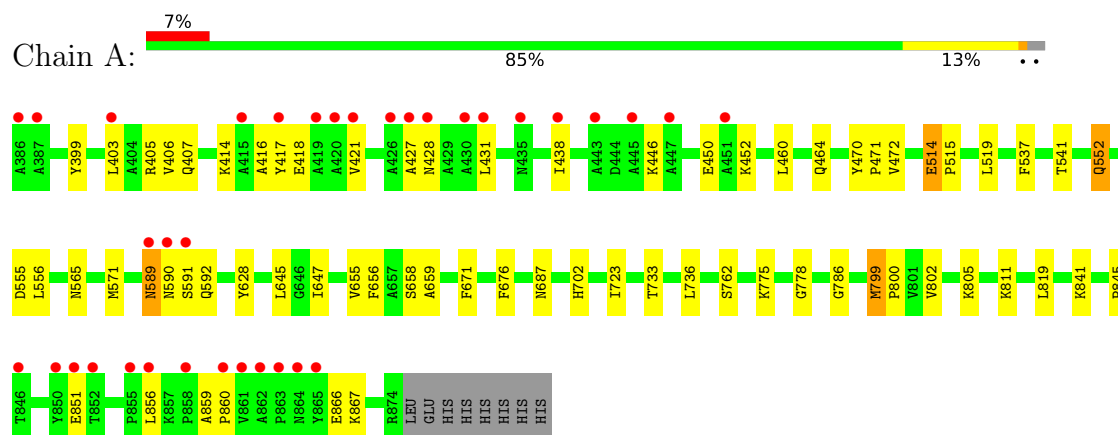
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	450	Total 450	O 450	0	0
5	B	560	Total 560	O 560	0	0

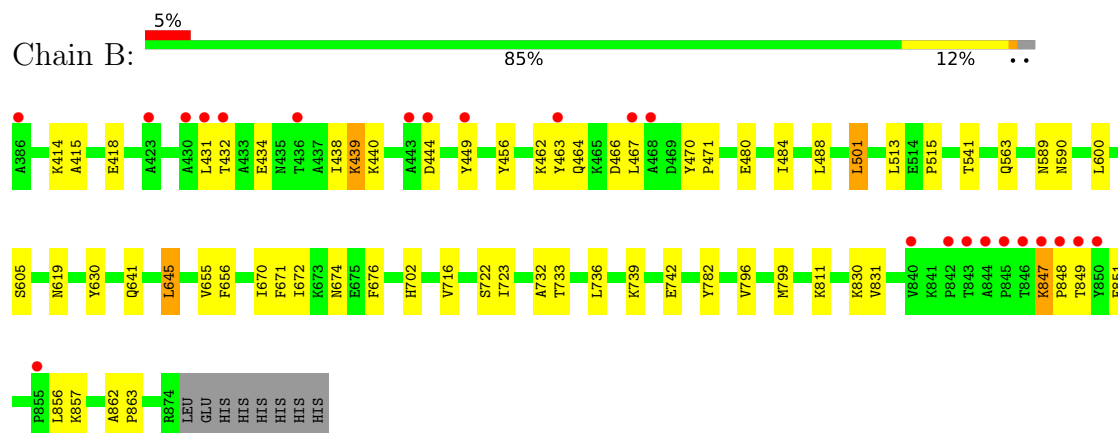
3 Residue-property plots [i](#)

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: AgI/II



• Molecule 1: AgI/II



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.03Å 164.15Å 67.73Å 90.00° 91.03° 90.00°	Depositor
Resolution (Å)	42.72 – 2.04 42.72 – 2.04	Depositor EDS
% Data completeness (in resolution range)	94.8 (42.72-2.04) 94.9 (42.72-2.04)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.73 (at 2.05Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.181 , 0.222 0.181 , 0.222	Depositor DCC
R_{free} test set	6734 reflections (10.16%)	wwPDB-VP
Wilson B-factor (Å ²)	15.2	Xtriage
Anisotropy	0.504	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.039 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8564	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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: PMS, SO4, CA

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	0.34	0/3839	0.85	10/5211 (0.2%)
1	B	0.34	0/3839	0.88	6/5211 (0.1%)
All	All	0.34	0/7678	0.86	16/10422 (0.2%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	733	THR	N-CA-C	7.69	119.74	111.36
1	B	799	MET	N-CA-C	6.75	118.57	110.07
1	A	541	THR	N-CA-C	6.53	120.34	112.38
1	B	733	THR	N-CA-C	6.19	119.00	111.82
1	B	736	LEU	N-CA-C	-5.99	101.91	110.59
1	A	799	MET	N-CA-C	5.88	118.40	110.29
1	A	736	LEU	N-CA-C	-5.86	102.10	110.59
1	B	541	THR	N-CA-C	5.66	119.28	112.38
1	A	514	GLU	CA-C-N	5.59	125.43	119.28
1	A	514	GLU	C-N-CA	5.59	125.43	119.28
1	B	723	ILE	N-CA-C	5.52	116.53	108.58
1	A	723	ILE	N-CA-C	5.24	116.13	108.58
1	A	676	PHE	N-CA-C	5.17	117.24	109.23
1	A	659	ALA	N-CA-C	-5.06	105.42	112.30
1	A	658	SER	N-CA-C	5.05	117.39	108.75
1	B	676	PHE	N-CA-C	5.04	117.04	109.23

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	3760	0	3684	44	0
1	B	3760	0	3684	40	0
2	A	11	0	8	1	0
2	B	11	0	8	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	450	0	0	9	0
5	B	560	0	0	8	0
All	All	8564	0	7384	83	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 (83) 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:438:ILE:HG23	1:A:851:GLU:HG3	1.27	1.15
1:B:470:TYR:HB3	1:B:471:PRO:HD3	1.71	0.73
1:A:431:LEU:HD23	1:A:856:LEU:HA	1.71	0.73
1:A:775:LYS:HD2	5:A:1259:HOH:O	1.93	0.68
1:A:647:ILE:HD13	1:A:655:VAL:HG23	1.75	0.67
1:A:470:TYR:HB3	1:A:471:PRO:HD3	1.79	0.63
1:B:672:ILE:HD12	1:B:672:ILE:N	2.18	0.58
1:A:414:LYS:O	1:A:418:GLU:HG3	2.05	0.57
1:A:417:TYR:O	1:A:421:VAL:HG23	2.04	0.57
1:B:484:ILE:O	1:B:488:LEU:HD13	2.04	0.57
1:A:472:VAL:HG23	5:A:1105:HOH:O	2.05	0.56
1:A:406:VAL:HG13	1:A:866:GLU:HG3	1.88	0.56
1:A:403:LEU:O	1:A:407:GLN:HG3	2.07	0.55
1:B:671:PHE:C	1:B:672:ILE:HD12	2.31	0.55
1:B:849:THR:HG23	1:B:849:THR:O	2.07	0.54
1:B:439:LYS:HB2	1:B:439:LYS:NZ	2.22	0.54
1:A:399:TYR:CZ	1:A:403:LEU:HD12	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:589:ASN:HD22	1:A:589:ASN:H	1.56	0.54
1:A:841:LYS:HE2	5:A:1237:HOH:O	2.10	0.51
1:B:438:ILE:HG23	1:B:851:GLU:HB2	1.93	0.51
1:A:565:ASN:ND2	5:A:1230:HOH:O	2.44	0.51
1:A:860:PRO:HD2	5:A:1819:HOH:O	2.10	0.51
1:B:432:THR:HG23	1:B:856:LEU:HD21	1.92	0.51
1:A:556:LEU:C	1:A:556:LEU:HD23	2.36	0.51
1:A:859:ALA:HB1	5:A:1819:HOH:O	2.11	0.50
1:A:552:GLN:HG2	1:A:555:ASP:CG	2.36	0.50
1:A:460:LEU:HG	1:A:464:GLN:NE2	2.27	0.50
1:A:537:PHE:CE2	1:A:556:LEU:HD12	2.47	0.49
1:B:456:TYR:HA	5:B:1498:HOH:O	2.13	0.49
1:B:513:LEU:HG	1:B:515:PRO:HD3	1.94	0.49
1:A:647:ILE:HD13	1:A:655:VAL:CG2	2.39	0.49
1:B:414:LYS:O	1:B:418:GLU:HG3	2.12	0.48
1:B:462:LYS:HE3	1:B:466:ASP:OD2	2.14	0.48
1:B:415:ALA:HB3	5:B:1899:HOH:O	2.12	0.48
1:B:847:LYS:HD3	1:B:848:PRO:N	2.29	0.48
1:A:537:PHE:CZ	1:A:556:LEU:HD12	2.49	0.48
1:A:446:LYS:O	1:A:450:GLU:HG3	2.14	0.47
1:A:702:HIS:HD2	5:A:922:HOH:O	1.97	0.47
1:A:762:SER:HB3	2:A:900:PMS:O3S	2.13	0.47
1:B:857:LYS:HB3	5:B:1762:HOH:O	2.13	0.47
1:A:591:SER:HB2	5:A:1805:HOH:O	2.15	0.47
1:A:416:ALA:HA	5:A:1804:HOH:O	2.15	0.46
1:B:480:GLU:O	1:B:484:ILE:HG12	2.16	0.46
1:B:431:LEU:HD13	5:B:1435:HOH:O	2.15	0.46
1:A:802:VAL:HG11	1:A:805:LYS:HD2	1.98	0.46
1:A:589:ASN:HD22	1:A:589:ASN:N	2.14	0.46
1:B:501:LEU:HD22	1:B:716:VAL:HG12	1.98	0.45
1:A:405:ARG:HH11	1:A:405:ARG:HG2	1.82	0.45
1:A:655:VAL:HG22	1:A:656:PHE:N	2.30	0.45
1:A:403:LEU:O	1:A:403:LEU:HD23	2.17	0.45
1:A:427:ALA:O	1:A:431:LEU:HB2	2.17	0.45
1:A:552:GLN:HE21	1:A:552:GLN:HB2	1.55	0.45
1:B:655:VAL:HG22	1:B:656:PHE:N	2.32	0.44
1:B:619:ASN:HA	1:B:831:VAL:HG11	1.99	0.44
1:B:722:SER:O	1:B:732:ALA:HA	2.17	0.44
1:A:452:LYS:HB3	1:A:845:PRO:HB3	2.00	0.44
1:B:830:LYS:HE2	5:B:1729:HOH:O	2.17	0.44
1:B:463:TYR:CZ	1:B:467:LEU:HD22	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:739:LYS:HD3	1:B:742:GLU:CD	2.43	0.43
1:A:514:GLU:N	1:A:515:PRO:HD3	2.33	0.43
1:A:628:TYR:OH	1:A:819:LEU:HD12	2.18	0.43
1:B:670:ILE:O	1:B:670:ILE:HG23	2.17	0.43
1:A:428:ASN:OD1	1:A:856:LEU:HG	2.19	0.43
1:B:600:LEU:HD11	1:B:645:LEU:HD22	1.99	0.43
1:B:862:ALA:HA	1:B:863:PRO:HD3	1.92	0.42
1:A:592:GLN:NE2	1:B:641:GLN:HE22	2.17	0.42
1:A:671:PHE:CD1	1:A:786:GLY:HA3	2.54	0.42
1:B:796:VAL:HG13	5:B:1465:HOH:O	2.19	0.42
1:B:434:GLU:O	1:B:438:ILE:HD13	2.20	0.42
1:A:687:ASN:OD1	1:A:778:GLY:HA2	2.20	0.42
1:B:440:LYS:O	1:B:444:ASP:HB2	2.20	0.42
1:A:519:LEU:O	1:A:571:MET:HE1	2.20	0.41
1:B:449:TYR:CD1	1:B:848:PRO:HD3	2.56	0.41
1:B:589:ASN:OD1	1:B:590:ASN:N	2.53	0.41
1:A:799:MET:HA	1:A:800:PRO:HD3	1.96	0.41
1:A:589:ASN:HD22	1:A:589:ASN:C	2.27	0.41
1:B:674:ASN:O	1:B:782:TYR:HA	2.20	0.41
1:B:702:HIS:HD2	5:B:1284:HOH:O	2.02	0.41
1:B:605:SER:HB2	1:B:630:TYR:O	2.20	0.41
1:B:600:LEU:CD1	1:B:645:LEU:HD22	2.51	0.41
1:B:439:LYS:HB2	1:B:439:LYS:HZ1	1.85	0.40
1:B:456:TYR:HB2	5:B:1897:HOH:O	2.22	0.40
1:B:563:GLN:N	1:B:563:GLN:CD	2.80	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	487/497 (98%)	475 (98%)	11 (2%)	1 (0%)	43	37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	487/497 (98%)	480 (99%)	7 (1%)	0	100	100
All	All	974/994 (98%)	955 (98%)	18 (2%)	1 (0%)	48	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	590	ASN

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	401/409 (98%)	396 (99%)	5 (1%)	63	65
1	B	401/409 (98%)	395 (98%)	6 (2%)	57	58
All	All	802/818 (98%)	791 (99%)	11 (1%)	59	60

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	552	GLN
1	A	589	ASN
1	A	645	LEU
1	A	811	LYS
1	A	867	LYS
1	B	439	LYS
1	B	464	GLN
1	B	501	LEU
1	B	645	LEU
1	B	811	LYS
1	B	847	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	400	GLN
1	A	407	GLN
1	A	410	ASN
1	A	425	ASN
1	A	464	GLN
1	A	494	HIS
1	A	516	ASN
1	A	548	GLN
1	A	552	GLN
1	A	565	ASN
1	A	589	ASN
1	A	590	ASN
1	A	592	GLN
1	A	663	GLN
1	A	702	HIS
1	B	393	GLN
1	B	424	ASN
1	B	464	GLN
1	B	560	ASN
1	B	592	GLN
1	B	690	ASN
1	B	702	HIS
1	B	740	GLN

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](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	SO4	B	903	-	4,4,4	0.38	0	6,6,6	0.10	0
2	PMS	B	901	-	10,11,11	3.24	6 (60%)	15,15,15	4.77	3 (20%)
2	PMS	A	900	-	10,11,11	3.29	6 (60%)	15,15,15	4.75	3 (20%)
3	SO4	A	902	-	4,4,4	0.34	0	6,6,6	0.07	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
2	PMS	B	901	-	-	0/5/5/5	0/1/1/1
2	PMS	A	900	-	-	0/5/5/5	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	901	PMS	C3-C2	4.78	1.47	1.38
2	A	900	PMS	C3-C2	4.75	1.47	1.38
2	A	900	PMS	C5-C6	4.52	1.46	1.38
2	B	901	PMS	C5-C6	4.49	1.46	1.38
2	A	900	PMS	C6-C1	4.14	1.47	1.38
2	B	901	PMS	C6-C1	4.03	1.46	1.38
2	A	900	PMS	C4-C3	4.02	1.47	1.38
2	A	900	PMS	C5-C4	3.88	1.46	1.38
2	B	901	PMS	C4-C3	3.87	1.46	1.38
2	B	901	PMS	C5-C4	3.82	1.46	1.38
2	A	900	PMS	C2-C1	3.78	1.46	1.38
2	B	901	PMS	C2-C1	3.64	1.46	1.38

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	PMS	C1-C-S	17.11	155.96	112.46
2	A	900	PMS	C1-C-S	17.02	155.73	112.46
2	B	901	PMS	O1S-S-C	5.63	119.29	105.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	900	PMS	O1S-S-C	5.54	119.08	105.82
2	A	900	PMS	O3S-S-O2S	-3.39	102.91	111.40
2	B	901	PMS	O3S-S-O2S	-3.36	103.00	111.40

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	900	PMS	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	489/497 (98%)	0.03	35 (7%) 21 21	7, 16, 49, 59	0
1	B	489/497 (98%)	-0.06	23 (4%) 36 36	7, 15, 46, 57	0
All	All	978/994 (98%)	-0.01	58 (5%) 28 28	7, 16, 48, 59	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	590	ASN	3.6
1	B	463	TYR	3.5
1	A	861	VAL	3.5
1	A	865	TYR	3.4
1	A	387	ALA	3.2
1	A	856	LEU	3.2
1	B	444	ASP	3.2
1	B	443	ALA	3.1
1	B	431	LEU	3.0
1	B	432	THR	3.0
1	A	438	ILE	2.9
1	B	840	VAL	2.9
1	A	420	ALA	2.8
1	A	589	ASN	2.8
1	B	849	THR	2.8
1	A	445	ALA	2.7
1	A	862	ALA	2.6
1	A	443	ALA	2.6
1	B	386	ALA	2.6
1	A	863	PRO	2.6
1	B	843	THR	2.6
1	A	435	ASN	2.5
1	A	421	VAL	2.5
1	A	415	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	860	PRO	2.5
1	A	403	LEU	2.5
1	B	467	LEU	2.5
1	B	847	LYS	2.5
1	A	591	SER	2.4
1	A	386	ALA	2.4
1	A	427	ALA	2.4
1	A	430	ALA	2.4
1	A	428	ASN	2.4
1	A	864	ASN	2.4
1	A	858	PRO	2.4
1	A	447	ALA	2.2
1	A	431	LEU	2.2
1	B	449	TYR	2.2
1	B	846	THR	2.2
1	A	855	PRO	2.2
1	B	468	ALA	2.2
1	A	417	TYR	2.2
1	B	423	ALA	2.1
1	B	430	ALA	2.1
1	A	846	THR	2.1
1	A	852	THR	2.1
1	B	845	PRO	2.1
1	B	436	THR	2.1
1	B	848	PRO	2.1
1	B	855	PRO	2.1
1	A	426	ALA	2.1
1	B	844	ALA	2.1
1	A	419	ALA	2.1
1	A	851	GLU	2.0
1	B	850	TYR	2.0
1	A	451	ALA	2.0
1	B	842	PRO	2.0
1	A	850	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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	PMS	B	901	11/11	0.85	0.16	38,39,43,43	0
2	PMS	A	900	11/11	0.89	0.15	37,39,39,40	0
3	SO4	B	903	5/5	0.91	0.13	48,48,49,50	0
3	SO4	A	902	5/5	0.93	0.12	39,41,42,43	0
4	CA	A	904	1/1	0.98	0.18	26,26,26,26	0
4	CA	B	905	1/1	0.98	0.15	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.