



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 07:50 am BST

PDB ID : 5AVN
Title : The 1.03 angstrom structure (P212121) of glucose isomerase crystallized in high-strength agarose hydrogel
Authors : Sugiyama, S.; Shimizu, N.; Maruyama, N.; Sazaki, G.; Adachi, H.; Takano, K.; Murakami, S.; Inoue, T.; Mori, Y.; Matsumura, H.
Deposited on : 2015-06-23
Resolution : 1.03 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

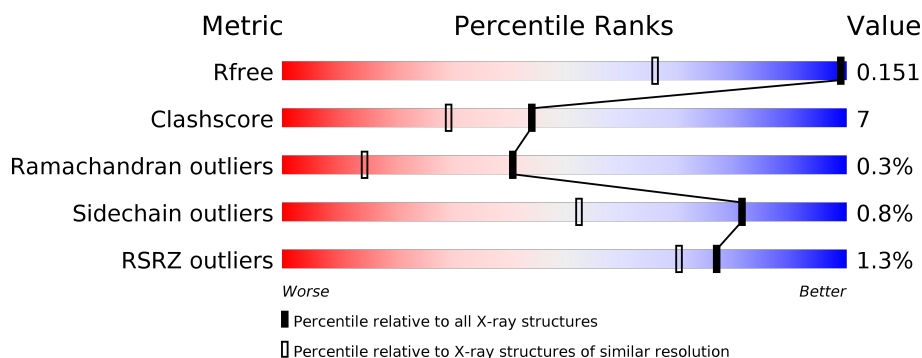
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.03 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}	130704	1596 (1.10-0.98)
Clashscore	141614	1677 (1.10-0.98)
Ramachandran outliers	138981	1591 (1.10-0.98)
Sidechain outliers	138945	1589 (1.10-0.98)
RSRZ outliers	127900	1557 (1.10-0.98)

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 on 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	387	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>11%</div> </div> </div>
1	B	387	<div> <div>2%</div> <div> <div></div> <div>91%</div> <div>9%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7812 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 Xylose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	27	0
			3276	2051	593	621	11			
1	B	387	Total	C	N	O	S	0	21	0
			3250	2029	598	613	10			

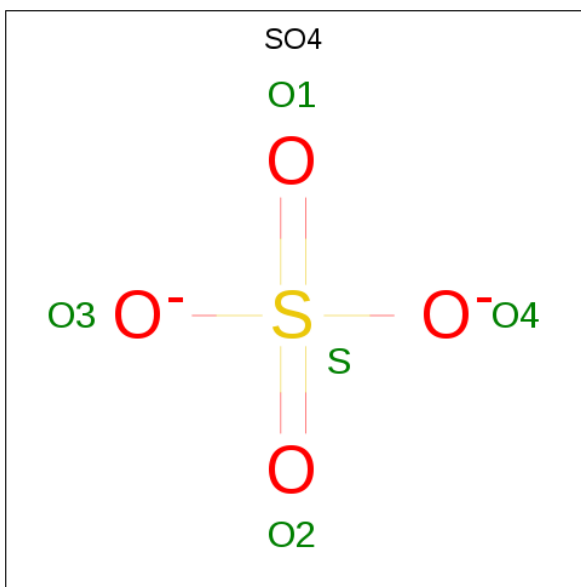
- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mn	0	0
			1	1		
2	A	1	Total	Mn	0	0
			1	1		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).

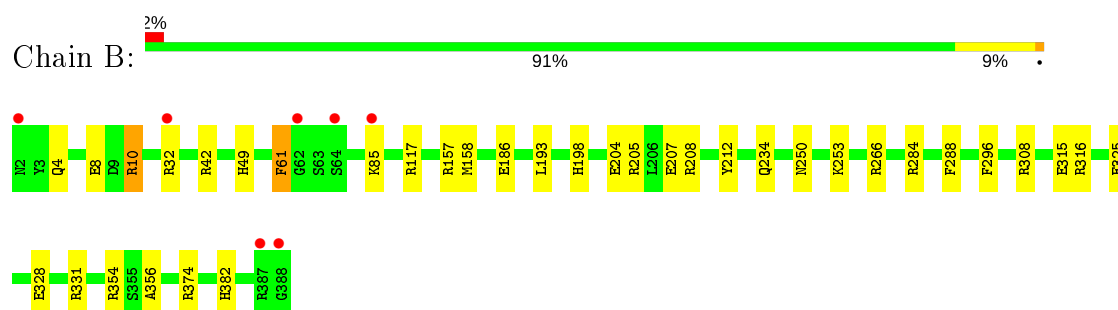
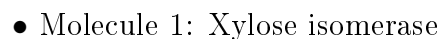


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	532	Total	O	0	124
			661	661		
5	B	526	Total	O	0	74
			601	601		

- Molecule 1: Xylose isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	94.08Å 98.28Å 85.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.03 26.44 – 1.03	Depositor EDS
% Data completeness (in resolution range)	99.2 (50.00-1.03) 99.2 (26.44-1.03)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 1.03Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.127 , 0.150 0.129 , 0.151	Depositor DCC
R_{free} test set	19438 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	8.5	Xtriage
Anisotropy	0.448	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.013 for k,h,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	7812	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 62.17 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1722e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: CA, MN, SO4

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	1.05	6/3353 (0.2%)	1.08	13/4531 (0.3%)
1	B	1.00	2/3323 (0.1%)	1.06	12/4485 (0.3%)
All	All	1.03	8/6676 (0.1%)	1.07	25/9016 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	250[A]	ASN	C-N	-6.42	1.21	1.33
1	A	250[B]	ASN	C-N	-6.42	1.21	1.33
1	B	207	GLU	CD-OE1	5.95	1.32	1.25
1	B	61	PHE	CB-CG	-5.73	1.41	1.51
1	A	315	GLU	CD-OE2	-5.53	1.19	1.25
1	A	169	VAL	CB-CG2	-5.41	1.41	1.52
1	A	8	GLU	CD-OE2	5.36	1.31	1.25
1	A	386	ALA	CA-C	5.07	1.66	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	250[A]	ASN	CA-C-N	-6.38	103.44	116.20
1	A	250[B]	ASN	CA-C-N	-6.38	103.44	116.20
1	B	212	TYR	CB-CG-CD2	5.97	124.58	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	381	ASP	CB-CG-OD1	5.93	123.64	118.30
1	A	188	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	B	157	ARG	NE-CZ-NH1	-5.79	117.41	120.30
1	B	331	ARG	NE-CZ-NH1	-5.70	117.45	120.30
1	A	212	TYR	CB-CG-CD2	5.69	124.41	121.00
1	A	334	ARG	NE-CZ-NH1	5.69	123.14	120.30
1	A	386	ALA	CB-CA-C	5.66	118.59	110.10
1	B	354	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	B	354	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	B	308	ARG	NE-CZ-NH2	-5.58	117.51	120.30
1	B	208	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	A	361	ASP	CB-CG-OD2	-5.46	113.38	118.30
1	B	32[A]	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	B	32[B]	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	A	109	ARG	NE-CZ-NH2	5.42	123.01	120.30
1	A	152	ARG	NE-CZ-NH2	-5.33	117.64	120.30
1	A	140	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	B	212	TYR	CB-CG-CD1	-5.22	117.87	121.00
1	B	117	ARG	NE-CZ-NH2	-5.15	117.72	120.30
1	B	296	PHE	CB-CG-CD2	-5.10	117.23	120.80
1	A	250[A]	ASN	CA-C-O	5.02	130.64	120.10
1	A	250[B]	ASN	CA-C-O	5.02	130.64	120.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	284	ARG	Sidechain

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	3276	0	3113	49	0
1	B	3250	0	3102	38	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	10	0	0	0	0
4	B	10	0	0	0	0
5	A	661	0	0	29	0
5	B	601	0	0	15	0
All	All	7812	0	6215	85	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 7.

All (85) 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:374[A]:ARG:NH2	5:B:502:HOH:O	1.63	1.32
1:A:266[B]:ARG:HD2	5:A:821:HOH:O	1.30	1.26
1:A:7[B]:PRO:HD2	1:A:8:GLU:OE2	1.36	1.25
1:B:325[A]:GLU:OE2	5:B:502:HOH:O	1.54	1.20
1:B:10[A]:ARG:CG	1:B:10[A]:ARG:HH11	1.55	1.19
1:B:266[B]:ARG:HD2	5:B:800:HOH:O	1.43	1.17
1:A:6[A]:THR:HG22	5:A:935:HOH:O	1.47	1.14
1:A:380[B]:MET:HA	1:A:380[B]:MET:HE2	1.18	1.11
1:B:315:GLU:OE1	1:B:316[B]:ARG:NH1	1.81	1.10
1:B:328[B]:GLU:OE1	5:B:503:HOH:O	1.67	1.10
1:A:380[B]:MET:HA	1:A:380[B]:MET:CE	1.84	1.08
1:B:10[A]:ARG:HH11	1:B:10[A]:ARG:HG2	1.09	1.06
1:B:266[B]:ARG:CD	5:B:800:HOH:O	1.99	0.98
1:A:132:GLU:OE2	5:A:503[A]:HOH:O	1.81	0.96
1:A:167:GLU:HG3	5:A:717:HOH:O	1.68	0.93
1:B:10[A]:ARG:NH1	1:B:10[A]:ARG:HG2	1.78	0.92
1:B:10[A]:ARG:CB	1:B:10[A]:ARG:HH11	1.81	0.92
1:A:167:GLU:HG3	5:A:585:HOH:O	1.70	0.90
1:A:266[B]:ARG:CD	5:A:821:HOH:O	1.95	0.87
1:A:7[B]:PRO:CD	1:A:8:GLU:OE2	2.21	0.86
1:A:6[A]:THR:HA	5:A:509:HOH:O	1.74	0.85
1:B:10[A]:ARG:CG	1:B:10[A]:ARG:NH1	2.31	0.85
1:A:207[B]:GLU:OE1	5:A:504:HOH:O	1.93	0.85
1:B:204[B]:GLU:O	1:B:205[B]:ARG:HD3	1.79	0.83
1:A:7[A]:PRO:HD3	5:A:509:HOH:O	1.78	0.83
1:B:10[A]:ARG:HD2	5:B:869:HOH:O	1.81	0.80
1:A:69[B]:GLU:HG2	5:A:527:HOH:O	1.81	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6[A]:THR:CG2	5:A:843:HOH:O	2.31	0.77
1:A:380[B]:MET:CA	1:A:380[B]:MET:CE	2.55	0.77
1:A:6[A]:THR:HG22	5:A:843:HOH:O	1.86	0.75
1:A:42[B]:ARG:NH2	5:A:505:HOH:O	2.19	0.72
1:B:8[B]:GLU:OE2	5:B:504:HOH:O	2.06	0.72
1:B:10[A]:ARG:NH1	1:B:10[A]:ARG:CB	2.51	0.71
1:A:387:ARG:NH1	5:A:507[A]:HOH:O	2.24	0.71
1:B:61:PHE:CD2	5:B:573:HOH:O	2.46	0.69
1:A:380[B]:MET:CA	1:A:380[B]:MET:HE2	2.07	0.69
1:B:204[B]:GLU:C	1:B:205[B]:ARG:HD3	2.13	0.68
1:A:7[A]:PRO:CD	5:A:509:HOH:O	2.41	0.67
1:A:77[B]:GLN:NE2	5:A:540[B]:HOH:O	2.27	0.67
1:B:8[B]:GLU:HA	1:B:10[B]:ARG:NH2	2.14	0.63
1:A:6[A]:THR:CA	5:A:509:HOH:O	2.38	0.63
1:B:4:GLN:OE1	5:B:505:HOH:O	2.15	0.62
1:A:42[B]:ARG:NE	5:A:505:HOH:O	2.25	0.62
1:B:10[A]:ARG:HB3	1:B:10[A]:ARG:NH1	2.15	0.62
1:B:205[B]:ARG:HH11	1:B:205[B]:ARG:HG3	1.64	0.61
1:A:6[A]:THR:CG2	5:A:935:HOH:O	2.23	0.60
1:A:380[B]:MET:HE3	1:A:380[B]:MET:CA	2.32	0.59
1:A:167:GLU:CG	5:A:717:HOH:O	2.40	0.59
1:B:8[B]:GLU:N	1:B:8[B]:GLU:OE1	2.31	0.59
1:B:205[B]:ARG:NH1	5:B:507:HOH:O	2.26	0.58
1:A:42[B]:ARG:CZ	5:A:505:HOH:O	2.51	0.57
1:A:7[A]:PRO:N	5:A:509:HOH:O	2.38	0.56
1:B:325[A]:GLU:CG	5:B:502:HOH:O	2.52	0.56
1:B:8[B]:GLU:O	1:B:8[B]:GLU:HG2	2.06	0.56
1:B:250:ASN:ND2	1:B:253[A]:LYS:NZ	2.54	0.55
1:B:205[B]:ARG:NH1	1:B:205[B]:ARG:HG3	2.23	0.54
1:B:316[B]:ARG:HD2	1:B:382:HIS:O	2.08	0.53
1:B:10[A]:ARG:HH12	1:B:49:HIS:CD2	2.26	0.53
1:A:6[A]:THR:C	5:A:509:HOH:O	2.46	0.52
1:A:32[A]:ARG:HG2	1:A:296[A]:PHE:CE2	2.45	0.52
1:A:370[B]:MET:CE	5:B:763:HOH:O	2.58	0.51
1:B:158[B]:MET:SD	1:B:193:LEU:HD11	2.52	0.50
1:A:2:ASN:N	5:A:511:HOH:O	2.44	0.50
1:A:380[B]:MET:HE3	1:A:380[B]:MET:HA	1.85	0.50
1:A:10:ARG:NH1	1:A:281:SER:O	2.45	0.49
1:A:380[B]:MET:HE3	1:A:380[B]:MET:HB2	1.75	0.47
1:B:316[A]:ARG:HD3	5:B:854[A]:HOH:O	2.15	0.46
1:A:167:GLU:CG	5:A:585:HOH:O	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158[B]:MET:SD	1:A:193:LEU:HD11	2.57	0.44
1:B:85[B]:LYS:HA	1:B:85[B]:LYS:HD2	1.82	0.44
1:B:250:ASN:ND2	1:B:253[A]:LYS:HZ3	2.14	0.44
1:A:234:GLN:HE22	1:B:198:HIS:HB3	1.83	0.44
1:A:370[B]:MET:HE1	5:B:763:HOH:O	2.18	0.44
1:B:316[A]:ARG:HA	1:B:316[A]:ARG:HD3	1.80	0.43
1:A:32[A]:ARG:HG2	1:A:296[A]:PHE:CZ	2.54	0.42
5:A:543[A]:HOH:O	1:B:356:ALA:HA	2.19	0.42
1:A:198:HIS:HB3	1:B:234:GLN:HE22	1.84	0.42
1:A:25[A]:PRO:HB2	1:A:26:PHE:CE2	2.55	0.42
1:A:324[A]:PRO:HG3	5:A:685:HOH:O	2.20	0.42
1:A:325[A]:GLU:OE2	5:A:506:HOH:O	2.20	0.42
1:B:42[A]:ARG:NH2	5:B:516:HOH:O	2.51	0.41
1:A:266[B]:ARG:NE	1:A:380[B]:MET:HE1	2.35	0.41
1:A:316:ARG:HA	1:A:316:ARG:HD3	1.81	0.41
1:A:25[A]:PRO:HB2	1:A:26:PHE:CD2	2.56	0.41
1:A:325[A]:GLU:CG	5:A:506:HOH:O	2.69	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	412/387 (106%)	400 (97%)	10 (2%)	2 (0%)	29	7
1	B	407/387 (105%)	396 (97%)	9 (2%)	2 (0%)	29	7
All	All	819/774 (106%)	796 (97%)	19 (2%)	4 (0%)	41	7

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	186[A]	GLU

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Mol	Chain	Res	Type
1	A	186[B]	GLU
1	B	186[A]	GLU
1	B	186[B]	GLU

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	330/303 (109%)	327 (99%)	3 (1%)	78	50	
1	B	325/303 (107%)	322 (99%)	3 (1%)	78	50	
All	All	655/606 (108%)	649 (99%)	6 (1%)	81	50	

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLU
1	A	91	THR
1	A	288	PHE
1	B	10[A]	ARG
1	B	10[B]	ARG
1	B	288	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	185	ASN
1	A	234	GLN
1	A	348	GLN
1	B	2	ASN
1	B	234	GLN
1	B	250	ASN
1	B	348	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 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	$\# Z > 2$
4	SO4	B	403	-	4,4,4	1.91	1 (25%)	6,6,6	0.70	0
4	SO4	B	404	-	4,4,4	0.39	0	6,6,6	0.99	0
4	SO4	A	404	-	4,4,4	0.47	0	6,6,6	0.68	0
4	SO4	A	403	-	4,4,4	1.18	0	6,6,6	0.59	0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	403	SO4	O1-S	3.58	1.65	1.46

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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	387/387 (100%)	-0.12	3 (0%) 86 81	6, 8, 16, 78	15 (3%)
1	B	387/387 (100%)	-0.10	7 (1%) 68 61	6, 9, 15, 95	17 (4%)
All	All	774/774 (100%)	-0.11	10 (1%) 77 70	6, 8, 16, 95	32 (4%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	388	GLY	13.4
1	A	388	GLY	8.1
1	B	387	ARG	3.4
1	B	62	GLY	3.4
1	B	85[A]	LYS	3.2
1	B	2	ASN	3.0
1	A	2	ASN	2.5
1	A	387	ARG	2.4
1	B	64	SER	2.1
1	B	32[A]	ARG	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 ⓘ

There are no carbohydrates 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(\AA^2)	Q<0.9
4	SO4	B	404	5/5	0.85	0.23	30,30,30,30	0
4	SO4	B	403	5/5	0.92	0.13	10,12,17,19	5
4	SO4	A	403	5/5	0.93	0.13	10,11,17,17	5
3	CA	B	402	1/1	0.94	0.24	17,17,17,17	1
4	SO4	A	404	5/5	0.96	0.08	15,21,25,32	5
3	CA	A	402	1/1	0.99	0.25	10,10,10,10	1
2	MN	B	401	1/1	1.00	0.03	7,7,7,7	0
2	MN	A	401	1/1	1.00	0.03	7,7,7,7	0

6.5 Other polymers [i](#)

There are no such residues in this entry.