



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:48 AM UTC

PDB ID : 5CQZ / pdb_00005cqz
Title : Human cytosolic 5'-nucleotidase II in complex with 3-(3-Imidazol-1-ylphenyl)-N-(9H-purin-6-yl)benzamide
Authors : Preeti, P.; Aghajari, N.
Deposited on : 2015-07-22
Resolution : 2.90 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

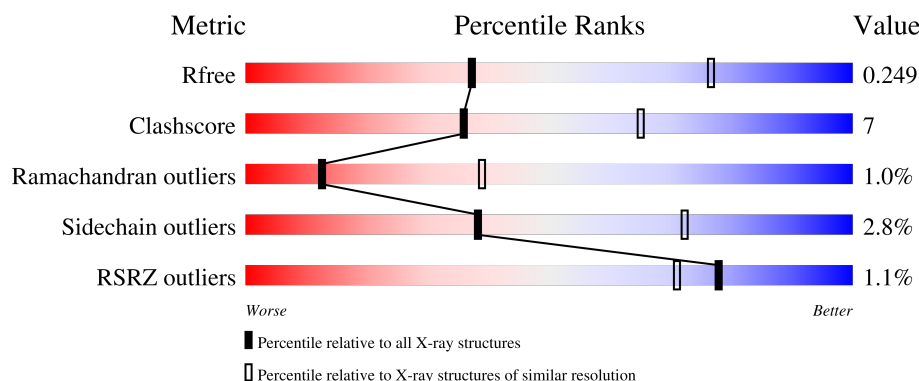
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	555	
1	B	555	

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	SO4	B	604	-	-	X	-
5	53O	B	609	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7702 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 Cytosolic purine 5'-nucleotidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	466	Total	C	N	O	S	0	1	0
			3806	2460	632	695	19			
1	B	453	Total	C	N	O	S	0	1	0
			3699	2393	611	676	19			

There are 38 discrepancies between the modelled and reference sequences:

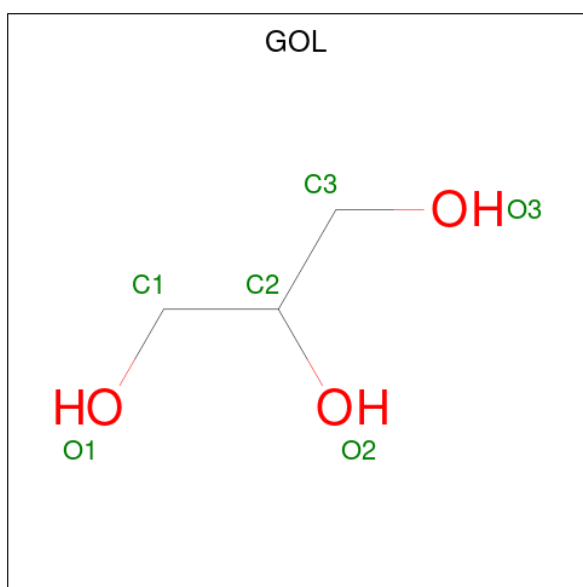
Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	initiating methionine	UNP P49902
A	-17	GLY	-	expression tag	UNP P49902
A	-16	SER	-	expression tag	UNP P49902
A	-15	SER	-	expression tag	UNP P49902
A	-14	HIS	-	expression tag	UNP P49902
A	-13	HIS	-	expression tag	UNP P49902
A	-12	HIS	-	expression tag	UNP P49902
A	-11	HIS	-	expression tag	UNP P49902
A	-10	HIS	-	expression tag	UNP P49902
A	-9	HIS	-	expression tag	UNP P49902
A	-8	SER	-	expression tag	UNP P49902
A	-7	SER	-	expression tag	UNP P49902
A	-6	GLY	-	expression tag	UNP P49902
A	-5	LEU	-	expression tag	UNP P49902
A	-4	VAL	-	expression tag	UNP P49902
A	-3	PRO	-	expression tag	UNP P49902
A	-2	ARG	-	expression tag	UNP P49902
A	-1	GLY	-	expression tag	UNP P49902
A	0	SER	-	expression tag	UNP P49902
B	-18	MET	-	initiating methionine	UNP P49902
B	-17	GLY	-	expression tag	UNP P49902
B	-16	SER	-	expression tag	UNP P49902
B	-15	SER	-	expression tag	UNP P49902
B	-14	HIS	-	expression tag	UNP P49902
B	-13	HIS	-	expression tag	UNP P49902

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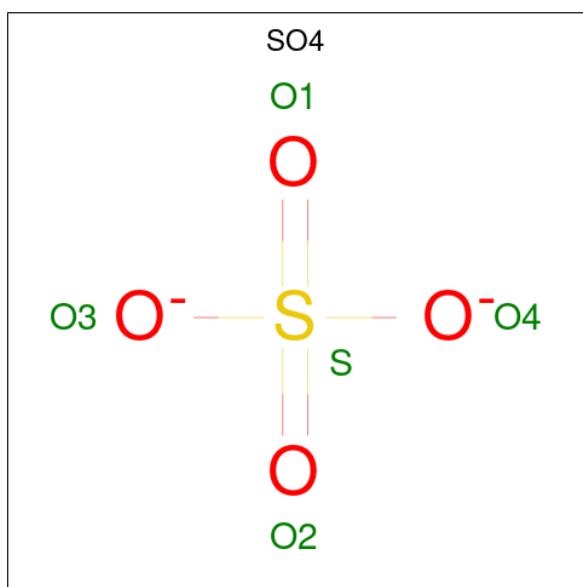
Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	HIS	-	expression tag	UNP P49902
B	-11	HIS	-	expression tag	UNP P49902
B	-10	HIS	-	expression tag	UNP P49902
B	-9	HIS	-	expression tag	UNP P49902
B	-8	SER	-	expression tag	UNP P49902
B	-7	SER	-	expression tag	UNP P49902
B	-6	GLY	-	expression tag	UNP P49902
B	-5	LEU	-	expression tag	UNP P49902
B	-4	VAL	-	expression tag	UNP P49902
B	-3	PRO	-	expression tag	UNP P49902
B	-2	ARG	-	expression tag	UNP P49902
B	-1	GLY	-	expression tag	UNP P49902
B	0	SER	-	expression tag	UNP P49902

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



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

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

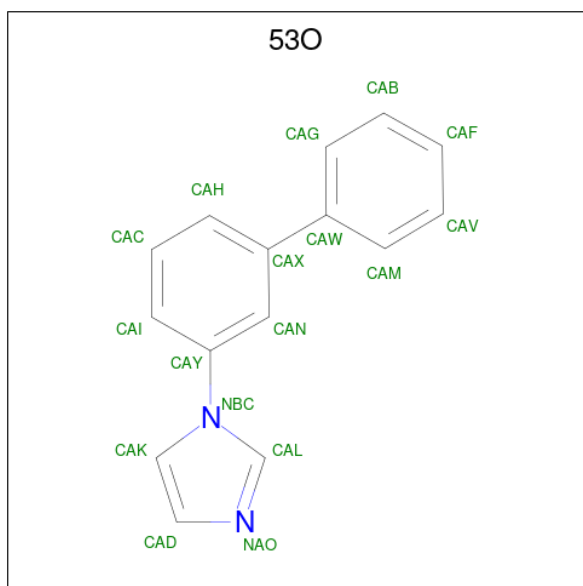
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Mg	0	0
			2	2		

- Molecule 5 is 1-(biphenyl-3-yl)-1H-imidazole (CCD ID: 53O) (formula: C₁₅H₁₂N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	N	0	0
			17	15	2		
5	B	1	Total	C	N	0	0
			17	15	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	44	Total	O	0	0
			44	44		
6	B	36	Total	O	0	0
			36	36		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.89Å 121.98Å 90.15Å 90.00° 114.29° 90.00°	Depositor
Resolution (Å)	19.88 – 2.90 19.88 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.6 (19.88-2.90) 98.3 (19.88-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.88Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.190 , 0.251 0.195 , 0.249	Depositor DCC
R_{free} test set	1550 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	50.5	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7702	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.33% 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, 53O, SO4, MG

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	0.61	0/3908	0.97	8/5278 (0.2%)
1	B	0.62	0/3797	0.96	10/5126 (0.2%)
All	All	0.62	0/7705	0.97	18/10404 (0.2%)

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	A	0	1

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	130	GLY	N-CA-C	-7.17	97.72	112.34
1	B	128	ILE	N-CA-C	6.99	117.78	110.72
1	B	129	ARG	N-CA-C	6.86	119.91	110.24
1	A	325	ILE	N-CA-C	6.57	118.14	109.21
1	B	435	GLY	N-CA-C	-6.53	102.01	110.69
1	A	435	GLY	N-CA-C	-6.27	102.36	110.69
1	A	211	LYS	N-CA-C	-5.80	106.83	114.31
1	B	189	ASP	N-CA-C	-5.71	103.57	110.88
1	A	323	HIS	N-CA-C	5.48	117.66	108.73
1	A	130	GLY	CA-C-N	5.37	125.18	119.28
1	A	130	GLY	C-N-CA	5.37	125.18	119.28
1	B	95	PHE	CA-C-N	-5.16	114.67	120.14
1	B	95	PHE	C-N-CA	-5.16	114.67	120.14
1	B	175	CYS	CA-C-N	5.14	124.75	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	175	CYS	C-N-CA	5.14	124.75	119.56
1	A	362	LYS	N-CA-C	5.10	121.65	110.80
1	B	148	GLU	N-CA-C	-5.08	105.83	112.23
1	A	81	GLY	N-CA-C	5.07	122.09	114.95

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	323	HIS	Peptide

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	3806	0	3740	68	0
1	B	3699	0	3626	47	0
2	A	12	0	16	1	0
2	B	12	0	16	1	0
3	A	35	0	0	1	0
3	B	20	0	0	3	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	17	0	12	5	0
5	B	17	0	12	11	0
6	A	44	0	0	1	0
6	B	36	0	0	0	0
All	All	7702	0	7422	109	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 (109) 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:446:ARG:HH21	5:B:609:53O:H11	1.27	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:612:53O:H9	1:B:445:SER:H	1.38	0.89
1:A:344:LYS:NZ	3:B:604:SO4:O1	2.08	0.86
1:B:456:ARG:HH21	1:B:457:TYR:HB2	1.47	0.79
1:A:344:LYS:NZ	3:B:604:SO4:S	2.61	0.73
1:A:446:ARG:HE	5:B:609:53O:H10	1.55	0.71
1:A:343:GLY:HA2	1:A:366:TRP:HD1	1.56	0.70
1:B:478:ARG:HH21	5:B:609:53O:H5	1.56	0.68
1:A:144:ARG:NH2	6:A:701:HOH:O	2.22	0.68
1:A:442:ARG:NH1	1:A:445:SER:O	2.24	0.67
1:A:445:SER:OG	5:B:609:53O:H9	1.94	0.67
1:A:360:SER:OG	1:A:361:LYS:N	2.25	0.67
1:B:478:ARG:NH2	5:B:609:53O:H5	2.10	0.65
1:B:29[B]:ARG:HB2	1:B:34:ARG:HG2	1.79	0.64
1:A:115:TYR:CD1	1:B:456:ARG:HG3	2.34	0.63
1:B:134:ARG:NH1	3:B:604:SO4:O4	2.32	0.63
1:B:478:ARG:HH21	5:B:609:53O:H3	1.64	0.62
1:A:343:GLY:HA2	1:A:366:TRP:CD1	2.34	0.62
1:B:386:SER:O	1:B:390:GLU:HG2	2.00	0.61
1:A:386:SER:O	1:A:390:GLU:HG3	1.99	0.61
5:A:612:53O:H11	1:B:446:ARG:CZ	2.31	0.61
5:A:612:53O:H10	1:B:446:ARG:HG3	1.83	0.60
1:A:85:GLU:OE2	1:A:177:ARG:NH2	2.35	0.59
1:A:247:LEU:HD23	1:A:286:ILE:HG23	1.84	0.59
1:A:342:LYS:HG2	1:A:364:GLN:HB3	1.85	0.59
1:A:356:ASP:N	1:A:356:ASP:OD1	2.36	0.59
1:B:97:THR:HG21	1:B:438:GLY:O	2.03	0.58
1:B:217:LYS:HA	1:B:220:GLU:HG2	1.84	0.58
1:A:446:ARG:HE	5:B:609:53O:CAB	2.15	0.58
1:A:29[B]:ARG:HH11	1:A:34:ARG:HE	1.51	0.58
1:B:305:VAL:HB	1:B:326:VAL:HG12	1.86	0.57
1:B:393:GLN:O	1:B:397:ILE:HG12	2.05	0.56
1:A:229:ASP:OD2	1:A:231:LYS:HG3	2.07	0.55
1:A:143:GLN:OE1	1:A:145:ASP:HB2	2.06	0.55
1:A:361:LYS:HD3	1:A:363:ARG:HH12	1.71	0.55
1:B:49:PHE:CD1	1:B:239:MET:HE2	2.41	0.54
1:A:39:ARG:HD2	1:B:446:ARG:HH21	1.72	0.54
1:A:140:LYS:NZ	3:A:604:SO4:O1	2.41	0.54
1:A:393:GLN:O	1:A:397:ILE:HG12	2.08	0.54
1:A:195:ARG:NH1	1:A:199:GLN:OE1	2.38	0.54
1:A:57:LEU:HD23	1:A:465:PHE:CZ	2.43	0.53
1:A:49:PHE:CD2	1:A:239:MET:HE2	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:GLN:NE2	1:B:482:VAL:O	2.36	0.53
1:B:143:GLN:HG3	1:B:145:ASP:HB2	1.90	0.52
1:A:304:GLN:NE2	1:A:325:ILE:HG23	2.24	0.52
1:B:354:PHE:O	1:B:456:ARG:NH2	2.41	0.52
1:A:378:GLU:HB3	2:A:601:GOL:H12	1.91	0.51
1:A:456:ARG:HG3	1:B:115:TYR:CD1	2.45	0.51
1:B:423:ILE:O	1:B:427:THR:HG23	2.10	0.51
1:A:395:LEU:HD21	1:A:422:ARG:HG3	1.93	0.51
1:A:146:ASP:C	1:A:148:GLU:H	2.19	0.50
1:A:232:LEU:HB3	1:A:233:PRO:HD3	1.93	0.50
1:B:104:THR:HB	1:B:195:ARG:HA	1.92	0.50
1:A:175:CYS:SG	1:A:177:ARG:HG3	2.52	0.50
1:B:302:LEU:HD13	1:B:334:THR:HG21	1.94	0.49
1:A:445:SER:HB3	5:B:609:53O:H7	1.93	0.49
1:A:456:ARG:NH2	1:A:457:TYR:HB2	2.28	0.49
1:B:304:GLN:HG2	1:B:313:LYS:HD2	1.94	0.48
1:A:333:ASP:OD2	1:A:363:ARG:HD3	2.14	0.48
1:B:390:GLU:HA	1:B:393:GLN:HB2	1.95	0.48
1:A:385:LYS:HE3	1:A:388:LEU:HD13	1.97	0.47
1:A:382:TRP:CD2	1:B:492:HIS:HE1	2.33	0.46
1:B:475:TYR:CZ	2:B:602:GOL:H32	2.50	0.46
1:A:385:LYS:HE2	1:A:433:CYS:SG	2.55	0.46
1:B:129:ARG:HA	1:B:133:THR:OG1	2.16	0.46
1:B:40:SER:HB2	5:B:609:53O:CAM	2.45	0.46
1:A:52:ASP:O	1:A:56:THR:HB	2.15	0.46
1:A:134:ARG:HA	1:A:137:TYR:O	2.16	0.46
1:B:306:ASP:OD2	1:B:309:THR:HG23	2.16	0.45
1:A:98:ARG:NE	1:A:374:GLU:OE2	2.33	0.45
1:A:104:THR:HB	1:A:195:ARG:HA	1.97	0.45
1:A:442:ARG:NH2	1:B:482:VAL:O	2.41	0.44
1:A:40:SER:HB2	5:A:612:53O:H8	1.99	0.44
1:B:229:ASP:OD2	1:B:231:LYS:HE2	2.18	0.44
1:B:30:GLU:O	1:B:34:ARG:HG3	2.17	0.44
1:A:489:THR:CB	1:B:34:ARG:HH12	2.31	0.43
1:B:222:LEU:HD13	1:B:262:TYR:HB2	1.99	0.43
1:A:29[B]:ARG:HB2	1:A:34:ARG:HG2	2.01	0.43
5:B:609:53O:H8	5:B:609:53O:H4	1.66	0.43
1:A:342:LYS:CG	1:A:343:GLY:H	2.31	0.43
1:A:109:LEU:HA	1:A:109:LEU:HD23	1.80	0.43
1:A:288:VAL:O	1:A:328:SER:HA	2.19	0.43
1:B:146:ASP:C	1:B:148:GLU:H	2.26	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:MET:HB2	1:A:475:TYR:O	2.19	0.42
1:B:186:LYS:HA	1:B:190:LEU:O	2.19	0.42
1:B:286:ILE:HG13	1:B:325:ILE:HG21	2.01	0.42
1:B:325:ILE:HG22	1:B:326:VAL:N	2.34	0.42
1:A:423:ILE:O	1:A:427:THR:HG23	2.19	0.42
5:A:612:53O:H8	5:A:612:53O:H4	1.62	0.42
1:A:421:ARG:HA	1:A:424:LYS:HD2	2.00	0.42
1:B:57:LEU:HD23	1:B:465:PHE:CZ	2.54	0.42
1:A:29[B]:ARG:NH1	1:A:34:ARG:HH21	2.17	0.42
1:A:50:GLY:HA3	1:A:246:PHE:CZ	2.55	0.42
1:A:323:HIS:CE1	1:A:325:ILE:O	2.72	0.42
1:A:342:LYS:HG3	1:A:343:GLY:H	1.85	0.42
1:B:247:LEU:HD23	1:B:286:ILE:HG23	2.02	0.42
1:A:197:MET:O	1:A:201:VAL:HG23	2.20	0.41
1:B:28:ARG:HG2	1:B:29[B]:ARG:HH11	1.85	0.41
1:A:197:MET:HE2	1:A:197:MET:HB3	1.88	0.41
1:A:361:LYS:O	1:A:363:ARG:N	2.53	0.41
5:B:609:53O:H4	5:B:609:53O:H1	1.65	0.41
1:B:57:LEU:HD23	1:B:57:LEU:HA	1.84	0.41
1:A:29[B]:ARG:H	1:A:29[B]:ARG:HG3	1.53	0.40
1:A:224:LYS:HD3	1:A:225:TYR:CZ	2.57	0.40
1:A:362:LYS:HA	1:A:362:LYS:HD2	1.83	0.40
1:A:392:LEU:HD21	1:A:427:THR:HG22	2.03	0.40
1:B:466:ILE:HD12	1:B:466:ILE:HG23	1.77	0.40
1:A:27:TYR:CG	1:B:404:LYS:HG3	2.56	0.40
1:A:489:THR:HB	1:B:34:ARG:HH12	1.85	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	463/555 (83%)	438 (95%)	18 (4%)	7 (2%)	8	28
1	B	448/555 (81%)	430 (96%)	14 (3%)	4 (1%)	14	41
All	All	911/1110 (82%)	868 (95%)	32 (4%)	11 (1%)	12	34

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	317	TYR
1	A	362	LYS
1	B	29[A]	ARG
1	B	29[B]	ARG
1	B	325	ILE
1	B	326	VAL
1	A	29[A]	ARG
1	A	29[B]	ARG
1	A	129	ARG
1	A	147	THR
1	A	361	LYS

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	415/497 (84%)	402 (97%)	13 (3%)	35	69
1	B	403/497 (81%)	392 (97%)	11 (3%)	39	73
All	All	818/994 (82%)	794 (97%)	24 (3%)	38	71

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLU
1	A	143	GLN
1	A	173	THR
1	A	177	ARG
1	A	231	LYS

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Mol	Chain	Res	Type
1	A	254	LYS
1	A	356	ASP
1	A	363	ARG
1	A	392	LEU
1	A	413	ARG
1	A	425	LYS
1	A	436	MET
1	A	466	ILE
1	B	29[A]	ARG
1	B	29[B]	ARG
1	B	129	ARG
1	B	173	THR
1	B	254	LYS
1	B	298	GLU
1	B	308	LYS
1	B	349	ILE
1	B	395	LEU
1	B	424	LYS
1	B	466	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	209	HIS
1	A	304	GLN
1	A	323	HIS
1	A	405	HIS
1	A	420	GLN
1	B	123	HIS
1	B	174	ASN
1	B	420	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 21 ligands modelled in this entry, 4 are monoatomic - leaving 17 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	605	-	4,4,4	0.30	0	6,6,6	0.29	0
5	53O	B	609	-	19,19,19	1.83	3 (15%)	25,25,25	1.90	5 (20%)
3	SO4	B	604	-	4,4,4	0.25	0	6,6,6	0.30	0
3	SO4	A	605	-	4,4,4	0.30	0	6,6,6	0.23	0
3	SO4	A	603	4	4,4,4	0.27	0	6,6,6	0.16	0
3	SO4	B	603	4	4,4,4	0.32	0	6,6,6	0.53	0
5	53O	A	612	-	19,19,19	1.86	3 (15%)	25,25,25	1.62	5 (20%)
2	GOL	B	601	-	5,5,5	0.44	0	5,5,5	0.94	0
3	SO4	A	609	-	4,4,4	0.26	0	6,6,6	0.30	0
3	SO4	A	608	-	4,4,4	0.29	0	6,6,6	0.22	0
2	GOL	A	601	-	5,5,5	0.41	0	5,5,5	0.27	0
2	GOL	A	602	-	5,5,5	0.44	0	5,5,5	0.37	0
3	SO4	B	606	-	4,4,4	0.25	0	6,6,6	0.15	0
2	GOL	B	602	-	5,5,5	0.43	0	5,5,5	0.35	0
3	SO4	A	604	-	4,4,4	0.24	0	6,6,6	0.36	0
3	SO4	A	607	-	4,4,4	0.39	0	6,6,6	0.43	0
3	SO4	A	606	-	4,4,4	0.37	0	6,6,6	0.18	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
5	53O	A	612	-	-	0/8/8/8	0/3/3/3
2	GOL	B	601	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	601	-	-	2/4/4/4	-
2	GOL	A	602	-	-	2/4/4/4	-
2	GOL	B	602	-	-	4/4/4/4	-
5	53O	B	609	-	-	0/8/8/8	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	609	53O	CAY-NBC	-5.82	1.33	1.43
5	A	612	53O	CAY-NBC	-5.63	1.34	1.43
5	A	612	53O	CAW-CAX	-4.41	1.38	1.49
5	B	609	53O	CAW-CAX	-4.11	1.39	1.49
5	A	612	53O	CAK-NBC	-2.64	1.33	1.38
5	B	609	53O	CAK-NBC	-2.33	1.33	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	609	53O	CAY-NBC-CAK	-4.37	120.44	126.84
5	B	609	53O	CAI-CAY-NBC	4.24	125.68	119.80
5	A	612	53O	CAI-CAY-NBC	3.45	124.58	119.80
5	B	609	53O	CAY-NBC-CAL	3.06	132.21	127.20
5	B	609	53O	CAN-CAX-CAW	-3.04	115.74	120.84
5	A	612	53O	CAY-NBC-CAK	-2.95	122.52	126.84
5	A	612	53O	CAN-CAX-CAW	-2.79	116.16	120.84
5	B	609	53O	CAN-CAY-NBC	-2.71	115.68	119.27
5	A	612	53O	CAN-CAY-NBC	-2.25	116.29	119.27
5	A	612	53O	CAB-CAG-CAW	-2.15	118.09	120.54

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	GOL	O1-C1-C2-C3
2	B	602	GOL	O1-C1-C2-C3
2	A	601	GOL	O1-C1-C2-O2
2	A	602	GOL	O1-C1-C2-C3
2	B	601	GOL	O1-C1-C2-C3
2	B	602	GOL	C1-C2-C3-O3
2	A	602	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	B	602	GOL	O1-C1-C2-O2
2	B	602	GOL	O2-C2-C3-O3
2	B	601	GOL	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	609	53O	11	0
3	B	604	SO4	3	0
5	A	612	53O	5	0
2	A	601	GOL	1	0
2	B	602	GOL	1	0
3	A	604	SO4	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 [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	A	466/555 (83%)	-0.27	4 (0%) 81 75	28, 40, 66, 90	1 (0%)
1	B	453/555 (81%)	-0.23	6 (1%) 75 67	29, 43, 66, 80	1 (0%)
All	All	919/1110 (82%)	-0.25	10 (1%) 78 71	28, 42, 67, 90	2 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	29[A]	ARG	3.7
1	B	29[A]	ARG	3.5
1	A	363	ARG	2.9
1	A	189	ASP	2.6
1	B	323	HIS	2.4
1	A	360	SER	2.4
1	B	32	TYR	2.2
1	B	363	ARG	2.1
1	B	492	HIS	2.1
1	B	325	ILE	2.1

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

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
5	53O	A	612	17/17	0.73	0.17	59,65,69,69	0
3	SO4	A	607	5/5	0.83	0.17	61,68,70,92	0
5	53O	B	609	17/17	0.83	0.14	50,64,69,69	0
3	SO4	A	606	5/5	0.88	0.21	51,52,65,77	0
3	SO4	A	604	5/5	0.89	0.10	61,70,81,84	0
3	SO4	A	609	5/5	0.89	0.21	70,72,88,97	0
3	SO4	B	604	5/5	0.90	0.09	63,65,75,79	0
3	SO4	B	606	5/5	0.90	0.18	81,82,100,102	0
3	SO4	B	605	5/5	0.92	0.20	58,63,72,85	0
3	SO4	A	605	5/5	0.92	0.17	61,63,77,90	0
2	GOL	A	602	6/6	0.93	0.14	44,46,51,53	0
3	SO4	A	608	5/5	0.94	0.21	51,54,76,77	0
2	GOL	B	602	6/6	0.96	0.08	41,45,49,53	0
2	GOL	A	601	6/6	0.96	0.07	42,44,48,48	0
4	MG	A	610	1/1	0.97	0.11	28,28,28,28	0
2	GOL	B	601	6/6	0.98	0.07	36,38,39,41	0
3	SO4	A	603	5/5	0.98	0.05	30,31,36,38	0
4	MG	B	607	1/1	0.99	0.05	28,28,28,28	0
4	MG	B	608	1/1	0.99	0.12	41,41,41,41	0
3	SO4	B	603	5/5	0.99	0.04	36,36,38,43	0
4	MG	A	611	1/1	0.99	0.18	44,44,44,44	0

6.5 Other polymers

There are no such residues in this entry.