



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 02:49 PM EDT

PDB ID : 1RHY / pdb_00001rhy
Title : Crystal structure of Imidazole Glycerol Phosphate Dehydratase
Authors : Sinha, S.C.; Chaudhuri, B.N.; Burgner, J.W.; Yakovleva, G.; Davisson, V.J.;
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Deposited on : 2003-11-14
Resolution : 2.30 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 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

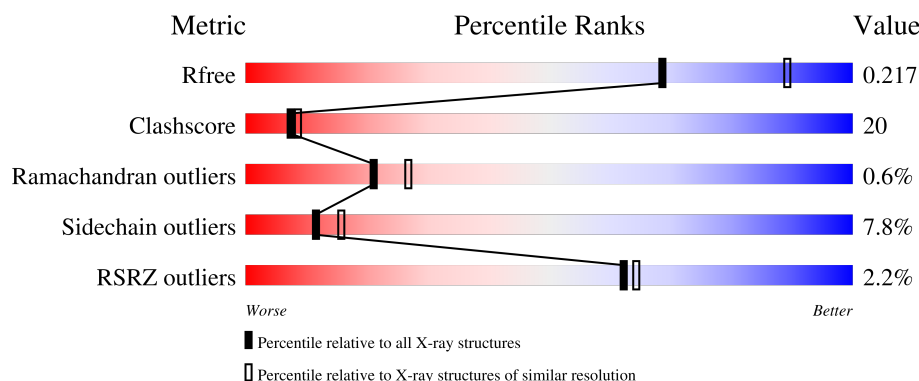
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	202	 63% 23% • 11%
1	B	202	 3% 61% 20% 6% • 11%

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
5	GOL	A	326	-	X	-	-
5	GOL	B	321	-	X	-	-
5	GOL	B	322	-	X	-	-
5	GOL	B	323	-	X	-	-
5	GOL	B	324	-	X	-	-
5	GOL	B	325	-	X	-	-
5	GOL	B	327	-	X	-	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 3141 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.

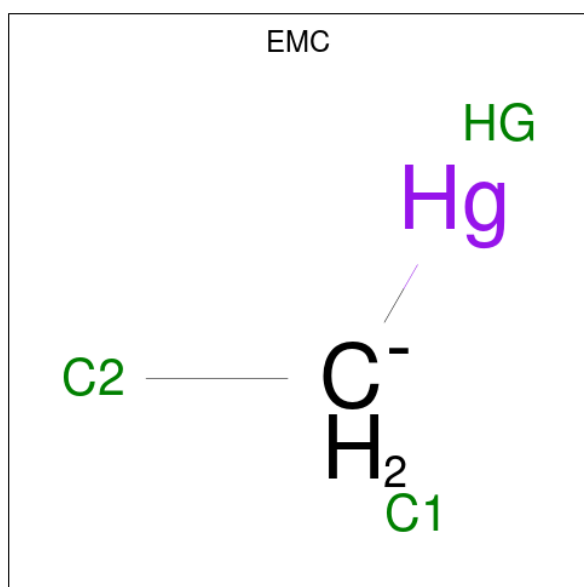
- Molecule 1 is a protein called Imidazole glycerol phosphate dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	180	Total	C	N	O	S	0	5	0
			1395	876	246	262	11			
1	B	180	Total	C	N	O	S	0	5	0
			1390	872	244	264	10			

- Molecule 2 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

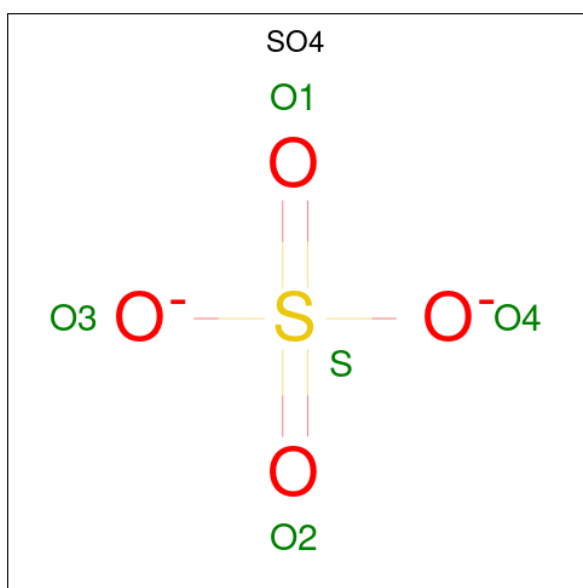
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Hg	0	0
			1	1		
2	B	1	Total	Hg	0	0
			1	1		

- Molecule 3 is ETHYL MERCURY ION (CCD ID: EMC) (formula: C₂H₅Hg).



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

- Molecule 4 is SULFATE ION (CCD ID: 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	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 GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 6 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		

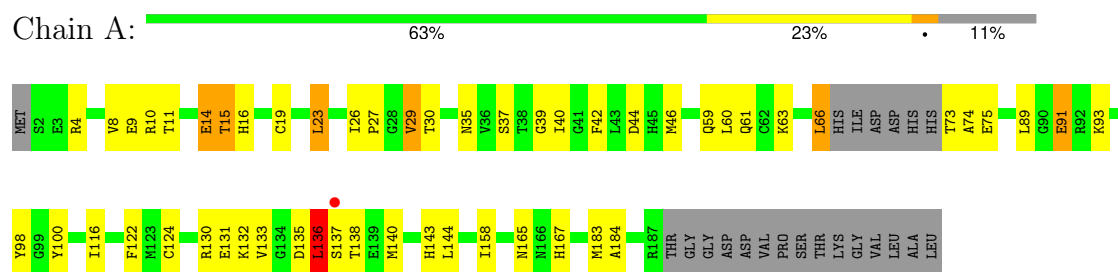
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	118	Total	O	0	0
			118	118		
7	B	129	Total	O	0	0
			129	129		

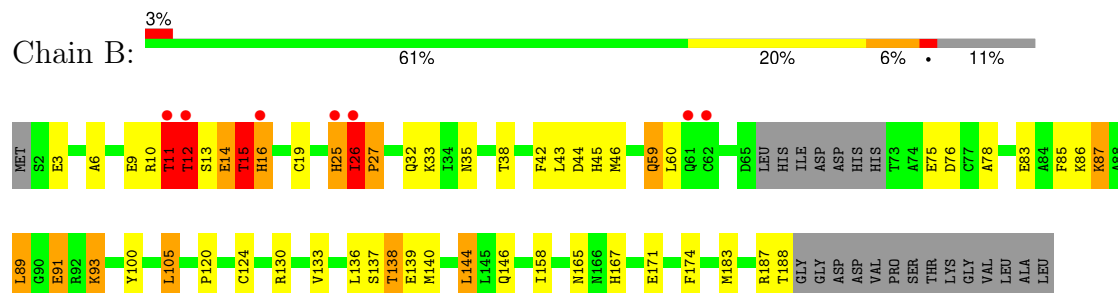
3 Residue-property plots

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: Imidazole glycerol phosphate dehydratase



- Molecule 1: Imidazole glycerol phosphate dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	105.28Å 105.28Å 105.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.55 – 2.30 19.55 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.4 (19.55-2.30) 97.2 (19.55-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.69 (at 2.21Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.189 , 0.228 0.183 , 0.217	Depositor DCC
R_{free} test set	878 reflections (4.45%)	wwPDB-VP
Wilson B-factor (Å ²)	36.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 80.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.044 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3141	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.95% 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: HG, EMC, ACY, SO4, 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	0.55	0/1444	0.97	3/1943 (0.2%)
1	B	0.58	0/1438	1.18	11/1936 (0.6%)
All	All	0.56	0/2882	1.08	14/3879 (0.4%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	26	ILE	CA-C-N	10.71	130.82	119.90
1	B	26	ILE	C-N-CA	10.71	130.82	119.90
1	B	16	HIS	N-CA-C	-9.00	101.04	111.03
1	B	187	ARG	N-CA-C	8.30	121.91	109.95
1	B	27	PRO	N-CA-C	7.53	123.04	111.15
1	B	11	THR	N-CA-C	-7.38	101.50	111.54
1	B	12	THR	N-CA-C	6.47	124.58	110.80
1	B	130	ARG	N-CA-C	-6.16	100.27	109.94
1	B	25	HIS	N-CA-C	-5.96	98.70	108.41
1	A	136	LEU	N-CA-C	5.88	119.14	109.72
1	B	138	THR	N-CA-C	5.65	117.44	111.28
1	A	184	ALA	N-CA-C	5.23	116.98	111.28
1	A	39	GLY	N-CA-C	-5.20	107.36	114.64
1	B	26	ILE	N-CA-C	-5.06	97.96	108.88

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	1395	0	1393	59	0
1	B	1390	0	1389	70	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	6	0	0	0	0
3	B	6	0	0	0	0
4	A	15	0	0	1	0
4	B	10	0	0	0	0
5	A	6	0	3	0	0
5	B	36	0	15	0	0
6	B	28	0	21	0	0
7	A	118	0	0	6	0
7	B	129	0	0	4	0
All	All	3141	0	2821	117	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 20.

All (117) 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:91:GLU:HB2	1:A:93:LYS:HG3	1.48	0.94
1:B:6:ALA:HB1	1:B:87:LYS:HG2	1.53	0.90
1:A:14:GLU:HG3	1:B:87:LYS:HZ1	1.42	0.84
1:B:43:LEU:HB2	1:B:136[B]:LEU:HD23	1.59	0.83
1:B:6:ALA:CB	1:B:87:LYS:HG2	2.09	0.82
1:B:9:GLU:HB3	1:B:12:THR:HB	1.61	0.82
1:A:136:LEU:HD22	1:A:137:SER:N	1.98	0.79
1:A:133[A]:VAL:HG23	1:A:138:THR:HG22	1.64	0.78
1:B:11:THR:HG21	7:B:9584:HOH:O	1.86	0.76
1:B:10:ARG:NH1	1:B:76[B]:ASP:OD2	2.21	0.74
1:B:9:GLU:CB	1:B:12:THR:HB	2.23	0.69
1:B:139:GLU:N	1:B:139:GLU:OE2	2.25	0.68
1:A:130:ARG:HG2	1:A:131:GLU:H	1.59	0.67
1:A:14:GLU:CG	1:B:87:LYS:HZ1	2.07	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:GLU:CG	1:B:87:LYS:NZ	2.58	0.67
1:A:40:ILE:HG21	1:A:136:LEU:HB2	1.79	0.66
1:A:35:ASN:HB3	4:A:305:SO4:O3	1.96	0.65
1:B:133[A]:VAL:HG23	1:B:138:THR:HG22	1.78	0.64
1:A:75:GLU:HG3	1:A:140[A]:MET:CE	2.29	0.63
1:A:75:GLU:HG3	1:A:140[A]:MET:HE2	1.79	0.63
1:A:15:THR:HG22	7:A:5562:HOH:O	1.99	0.63
1:B:35:ASN:HB2	1:B:59:GLN:HE21	1.63	0.63
1:B:75:GLU:HG3	1:B:140:MET:HE1	1.82	0.62
1:A:136:LEU:HD22	1:A:137:SER:H	1.63	0.62
1:A:100:TYR:CZ	1:A:183[B]:MET:HE1	2.35	0.61
1:A:14:GLU:HG2	1:B:87:LYS:NZ	2.16	0.61
1:B:136[B]:LEU:HD12	1:B:137:SER:O	2.00	0.61
1:A:9:GLU:CD	1:B:12:THR:HG23	2.26	0.60
1:B:105:LEU:HD23	1:B:171:GLU:HB3	1.84	0.60
1:B:100:TYR:CZ	1:B:183[A]:MET:HE1	2.37	0.60
1:A:9:GLU:OE2	1:B:13:SER:HB2	2.02	0.59
1:B:14:GLU:O	1:B:16:HIS:N	2.36	0.59
1:B:14:GLU:O	1:B:15:THR:C	2.45	0.58
1:B:14:GLU:O	1:B:14:GLU:CG	2.52	0.57
1:B:15:THR:HG21	7:B:9606:HOH:O	2.05	0.57
1:A:130:ARG:CG	1:A:131:GLU:H	2.18	0.56
1:B:14:GLU:O	1:B:14:GLU:OE2	2.23	0.56
1:A:26:ILE:CG2	1:A:29:VAL:HB	2.36	0.56
1:B:33:LYS:NZ	1:B:35:ASN:ND2	2.54	0.56
1:A:14:GLU:HG3	1:B:87:LYS:NZ	2.17	0.55
1:A:66:LEU:HD23	1:A:66:LEU:N	2.21	0.55
1:A:132:LYS:NZ	1:A:132:LYS:HB3	2.21	0.55
1:A:66:LEU:N	1:A:66:LEU:CD2	2.69	0.55
1:B:38:THR:HG23	1:B:44:ASP:OD1	2.07	0.55
1:A:26:ILE:HG22	1:A:29:VAL:HB	1.88	0.54
1:A:9:GLU:OE1	1:B:12:THR:HG23	2.07	0.54
1:A:11:THR:OG1	1:B:12:THR:HG21	2.08	0.54
1:B:140:MET:O	1:B:144:LEU:HD13	2.08	0.53
1:B:26:ILE:CG2	1:B:26:ILE:O	2.57	0.53
1:B:136[B]:LEU:CD1	1:B:137:SER:N	2.72	0.52
1:A:42:PHE:CZ	1:A:133[A]:VAL:HG11	2.44	0.52
1:A:136:LEU:C	1:A:136:LEU:HD13	2.34	0.52
1:A:9:GLU:O	1:B:11:THR:N	2.37	0.51
1:B:14:GLU:O	1:B:14:GLU:CD	2.53	0.51
1:A:63:LYS:HE2	7:A:5573:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:146:GLN:NE2	7:B:9547:HOH:O	2.44	0.51
1:B:35:ASN:HD22	1:B:59:GLN:NE2	2.09	0.50
1:A:29:VAL:HG12	1:A:30:THR:HG23	1.92	0.50
1:A:42:PHE:H	1:A:135:ASP:HB3	1.76	0.50
1:A:116:ILE:HG22	1:A:116:ILE:O	2.12	0.50
1:B:136[B]:LEU:HD13	1:B:137:SER:N	2.28	0.49
1:B:78:ALA:HB2	1:B:144:LEU:HB2	1.94	0.49
1:A:14:GLU:HG2	1:B:87:LYS:HZ3	1.77	0.49
1:A:124:CYS:HA	1:A:158:ILE:O	2.13	0.49
1:A:165:ASN:OD1	1:A:167:HIS:HB2	2.12	0.48
1:B:46:MET:HA	1:B:46:MET:HE2	1.95	0.48
1:B:85:PHE:CE2	1:B:89:LEU:HD21	2.48	0.48
1:B:42:PHE:O	1:B:45:HIS:HB3	2.13	0.47
1:A:122:PHE:CZ	1:A:124:CYS:HB3	2.49	0.47
1:A:29:VAL:HG12	1:A:30:THR:N	2.29	0.46
1:B:26:ILE:HA	1:B:27:PRO:HD2	1.65	0.46
1:B:33:LYS:HZ3	1:B:35:ASN:ND2	2.13	0.46
1:B:33:LYS:HZ3	1:B:35:ASN:HD21	1.63	0.46
1:A:46:MET:HE2	1:A:46:MET:HA	1.97	0.46
1:B:19:CYS:HA	1:B:59:GLN:O	2.15	0.46
1:A:75:GLU:OE1	1:A:140[A]:MET:HE1	2.16	0.46
1:B:133[A]:VAL:HG23	1:B:138:THR:CG2	2.46	0.46
1:B:136[B]:LEU:HD13	1:B:137:SER:H	1.80	0.45
1:B:137:SER:OG	1:B:139:GLU:OE2	2.29	0.45
1:A:27:PRO:HA	7:A:5563:HOH:O	2.16	0.45
1:A:130:ARG:HG2	1:A:131:GLU:N	2.29	0.45
1:A:140[A]:MET:HE3	1:A:143:HIS:HB2	1.97	0.45
1:B:10:ARG:HD2	1:B:83:GLU:OE2	2.17	0.45
7:A:5535:HOH:O	1:B:86:LYS:HE2	2.17	0.45
1:A:130:ARG:CG	1:A:131:GLU:N	2.81	0.44
1:B:91:GLU:CD	1:B:91:GLU:H	2.25	0.44
1:A:136:LEU:HD21	1:A:140[B]:MET:HB3	1.99	0.44
1:B:165:ASN:OD1	1:B:167:HIS:HB2	2.18	0.44
1:B:188:THR:CB	7:B:9605:HOH:O	2.66	0.43
1:A:98:TYR:HD2	7:A:5595:HOH:O	2.00	0.43
1:B:144:LEU:HD21	1:B:174:PHE:HZ	1.83	0.43
1:A:136:LEU:HD21	1:A:140[A]:MET:HB3	1.99	0.43
1:B:60:LEU:C	1:B:60:LEU:HD23	2.43	0.43
1:A:40:ILE:HG23	1:A:135:ASP:OD1	2.18	0.43
1:B:25:HIS:C	1:B:27:PRO:HD3	2.44	0.43
1:B:6:ALA:HB2	1:B:87:LYS:HG2	1.95	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:GLU:CD	1:A:16:HIS:HE2	2.27	0.42
1:A:132:LYS:HA	1:A:137:SER:HA	2.01	0.42
1:A:19:CYS:HA	1:A:59:GLN:O	2.19	0.42
1:B:35:ASN:HB2	1:B:59:GLN:NE2	2.31	0.42
1:A:73:THR:HB	1:A:74:ALA:H	1.60	0.42
1:A:30:THR:HG21	7:A:5619:HOH:O	2.18	0.42
1:B:12:THR:H	1:B:12:THR:HG22	1.53	0.42
1:A:37:SER:HA	1:A:44:ASP:OD2	2.20	0.42
1:A:66:LEU:CD2	1:A:66:LEU:H	2.33	0.41
1:A:10:ARG:HG2	1:B:10:ARG:HG3	2.02	0.41
1:B:91:GLU:CG	1:B:93:LYS:HG3	2.51	0.41
1:A:4:ARG:HB2	1:A:23:LEU:HB2	2.02	0.41
1:B:12:THR:O	1:B:13:SER:C	2.63	0.41
1:B:85:PHE:O	1:B:89:LEU:HD22	2.21	0.40
1:B:25:HIS:CD2	1:B:32:GLN:NE2	2.89	0.40
1:B:124:CYS:HA	1:B:158:ILE:O	2.22	0.40
1:B:15:THR:O	1:B:15:THR:HG22	2.21	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	181/202 (90%)	172 (95%)	8 (4%)	1 (1%)	21	27
1	B	181/202 (90%)	170 (94%)	10 (6%)	1 (1%)	21	27
All	All	362/404 (90%)	342 (94%)	18 (5%)	2 (1%)	21	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	29	VAL

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Mol	Chain	Res	Type
1	B	15	THR

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	152/165 (92%)	142 (93%)	10 (7%)	15	21
1	B	152/165 (92%)	139 (91%)	13 (9%)	10	13
All	All	304/330 (92%)	281 (92%)	23 (8%)	11	16

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	VAL
1	A	14	GLU
1	A	15	THR
1	A	23	LEU
1	A	60	LEU
1	A	61	GLN
1	A	66	LEU
1	A	89	LEU
1	A	91	GLU
1	A	136	LEU
1	B	3	GLU
1	B	11	THR
1	B	12	THR
1	B	14	GLU
1	B	15	THR
1	B	26	ILE
1	B	59	GLN
1	B	87	LYS
1	B	89	LEU
1	B	91	GLU
1	B	93	LYS
1	B	105	LEU

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Mol	Chain	Res	Type
1	B	144	LEU

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

Mol	Chain	Res	Type
1	A	32	GLN
1	B	25	HIS
1	B	32	GLN
1	B	35	ASN
1	B	59	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 25 ligands modelled in this entry, 2 are monoatomic - leaving 23 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	EMC	B	9502	1	1,2,2	0.52	0	-		
5	GOL	B	324	-	5,5,5	4.41	5 (100%)	5,5,5	5.34	4 (80%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	ACY	B	313	-	3,3,3	1.37	1 (33%)	3,3,3	1.59	1 (33%)
5	GOL	B	325	-	5,5,5	4.35	5 (100%)	5,5,5	5.23	4 (80%)
3	EMC	A	5502	1	1,2,2	0.23	0	-		
4	SO4	B	303	-	4,4,4	0.37	0	6,6,6	0.09	0
6	ACY	B	312	-	3,3,3	1.38	1 (33%)	3,3,3	1.72	1 (33%)
3	EMC	B	9501	1	1,2,2	0.45	0	-		
4	SO4	A	305	-	4,4,4	0.37	0	6,6,6	0.09	0
4	SO4	A	304	-	4,4,4	0.37	0	6,6,6	0.10	0
5	GOL	B	321	-	5,5,5	4.42	5 (100%)	5,5,5	5.25	4 (80%)
6	ACY	B	316	-	3,3,3	1.41	1 (33%)	3,3,3	1.66	1 (33%)
3	EMC	A	5501	1	1,2,2	0.38	0	-		
6	ACY	B	317	-	3,3,3	1.36	1 (33%)	3,3,3	1.57	1 (33%)
6	ACY	B	311	-	3,3,3	1.40	1 (33%)	3,3,3	1.71	1 (33%)
5	GOL	B	327	-	5,5,5	4.18	4 (80%)	5,5,5	5.26	4 (80%)
5	GOL	B	322	-	5,5,5	4.41	5 (100%)	5,5,5	5.30	4 (80%)
4	SO4	A	301	-	4,4,4	0.38	0	6,6,6	0.15	0
5	GOL	A	326	-	5,5,5	4.31	5 (100%)	5,5,5	5.26	4 (80%)
5	GOL	B	323	-	5,5,5	4.57	5 (100%)	5,5,5	5.20	4 (80%)
6	ACY	B	314	-	3,3,3	1.41	1 (33%)	3,3,3	1.61	1 (33%)
6	ACY	B	315	-	3,3,3	1.39	1 (33%)	3,3,3	1.72	1 (33%)
4	SO4	B	302	-	4,4,4	0.42	0	6,6,6	0.10	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	GOL	B	322	-	-	2/4/4/4	-
5	GOL	B	324	-	-	3/4/4/4	-
5	GOL	A	326	-	-	2/4/4/4	-
5	GOL	B	323	-	-	3/4/4/4	-
5	GOL	B	321	-	-	2/4/4/4	-
5	GOL	B	325	-	-	2/4/4/4	-
5	GOL	B	327	-	-	3/4/4/4	-

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	324	GOL	C1-C2	-5.67	1.30	1.51
5	B	323	GOL	C1-C2	-5.45	1.31	1.51
5	B	321	GOL	C1-C2	-5.42	1.31	1.51
5	B	325	GOL	C1-C2	-5.33	1.31	1.51
5	B	323	GOL	C3-C2	-5.33	1.31	1.51
5	B	322	GOL	C1-C2	-5.22	1.31	1.51
5	A	326	GOL	C1-C2	-5.10	1.32	1.51
5	B	322	GOL	C3-C2	-5.10	1.32	1.51
5	A	326	GOL	C3-C2	-5.02	1.32	1.51
5	B	321	GOL	C3-C2	-5.01	1.32	1.51
5	B	327	GOL	C3-C2	-4.93	1.33	1.51
5	B	325	GOL	C3-C2	-4.86	1.33	1.51
5	B	324	GOL	C3-C2	-4.71	1.33	1.51
5	B	327	GOL	C1-C2	-4.59	1.34	1.51
5	B	327	GOL	O1-C1	4.52	1.61	1.42
5	B	324	GOL	O3-C3	4.50	1.61	1.42
5	B	325	GOL	O3-C3	4.50	1.61	1.42
5	B	321	GOL	O3-C3	4.47	1.61	1.42
5	B	322	GOL	O3-C3	4.34	1.60	1.42
5	A	326	GOL	O1-C1	4.34	1.60	1.42
5	B	322	GOL	O1-C1	4.26	1.60	1.42
5	B	323	GOL	O1-C1	4.26	1.60	1.42
5	B	327	GOL	O3-C3	4.24	1.60	1.42
5	B	324	GOL	O1-C1	4.24	1.60	1.42
5	A	326	GOL	O3-C3	4.21	1.60	1.42
5	B	323	GOL	O3-C3	4.18	1.60	1.42
5	B	325	GOL	O1-C1	4.18	1.60	1.42
5	B	321	GOL	O1-C1	4.09	1.59	1.42
5	B	323	GOL	O2-C2	-3.29	1.33	1.43
5	B	322	GOL	O2-C2	-2.63	1.35	1.43
5	B	321	GOL	O2-C2	-2.57	1.35	1.43
6	B	314	ACY	O-C	2.23	1.32	1.22
5	A	326	GOL	O2-C2	-2.22	1.36	1.43
5	B	325	GOL	O2-C2	-2.20	1.37	1.43
6	B	315	ACY	O-C	2.19	1.31	1.22
6	B	316	ACY	O-C	2.19	1.31	1.22
6	B	311	ACY	O-C	2.18	1.31	1.22
6	B	312	ACY	O-C	2.16	1.31	1.22
5	B	324	GOL	O2-C2	-2.15	1.37	1.43
6	B	313	ACY	O-C	2.13	1.31	1.22
6	B	317	ACY	O-C	2.07	1.31	1.22

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	324	GOL	O3-C3-C2	7.68	144.95	110.38
5	B	322	GOL	O1-C1-C2	7.47	144.02	110.38
5	B	327	GOL	O1-C1-C2	7.40	143.71	110.38
5	B	321	GOL	O1-C1-C2	7.40	143.69	110.38
5	B	327	GOL	O3-C3-C2	7.40	143.69	110.38
5	B	323	GOL	O3-C3-C2	7.39	143.66	110.38
5	A	326	GOL	O1-C1-C2	7.38	143.60	110.38
5	B	322	GOL	O3-C3-C2	7.37	143.54	110.38
5	B	325	GOL	O3-C3-C2	7.35	143.46	110.38
5	B	324	GOL	O1-C1-C2	7.30	143.22	110.38
5	A	326	GOL	O3-C3-C2	7.29	143.22	110.38
5	B	321	GOL	O3-C3-C2	7.29	143.18	110.38
5	B	325	GOL	O1-C1-C2	7.25	143.01	110.38
5	B	323	GOL	O1-C1-C2	7.24	142.95	110.38
5	B	325	GOL	O2-C2-C3	4.03	125.86	109.18
5	B	321	GOL	O2-C2-C3	3.99	125.69	109.18
5	A	326	GOL	O2-C2-C3	3.92	125.42	109.18
5	B	322	GOL	O2-C2-C3	3.92	125.42	109.18
5	B	324	GOL	O2-C2-C3	3.89	125.29	109.18
5	B	324	GOL	O2-C2-C1	3.87	125.22	109.18
5	A	326	GOL	O2-C2-C1	3.85	125.10	109.18
5	B	327	GOL	O2-C2-C1	3.85	125.10	109.18
5	B	323	GOL	O2-C2-C3	3.82	125.00	109.18
5	B	322	GOL	O2-C2-C1	3.79	124.87	109.18
5	B	321	GOL	O2-C2-C1	3.70	124.48	109.18
5	B	327	GOL	O2-C2-C3	3.69	124.47	109.18
5	B	323	GOL	O2-C2-C1	3.67	124.39	109.18
5	B	325	GOL	O2-C2-C1	3.67	124.37	109.18
6	B	312	ACY	O-C-CH3	-2.33	112.96	122.53
6	B	311	ACY	O-C-CH3	-2.33	112.98	122.53
6	B	315	ACY	O-C-CH3	-2.33	112.99	122.53
6	B	316	ACY	O-C-CH3	-2.25	113.28	122.53
6	B	314	ACY	O-C-CH3	-2.18	113.59	122.53
6	B	313	ACY	O-C-CH3	-2.15	113.73	122.53
6	B	317	ACY	O-C-CH3	-2.14	113.76	122.53

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	326	GOL	O1-C1-C2-C3
5	A	326	GOL	C1-C2-C3-O3
5	B	321	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
5	B	322	GOL	C1-C2-C3-O3
5	B	323	GOL	O1-C1-C2-C3
5	B	323	GOL	C1-C2-C3-O3
5	B	324	GOL	C1-C2-C3-O3
5	B	325	GOL	O1-C1-C2-C3
5	B	325	GOL	C1-C2-C3-O3
5	B	327	GOL	C1-C2-C3-O3
5	B	324	GOL	O1-C1-C2-C3
5	B	327	GOL	O1-C1-C2-C3
5	B	324	GOL	O1-C1-C2-O2
5	B	327	GOL	O1-C1-C2-O2
5	B	321	GOL	O1-C1-C2-O2
5	B	322	GOL	O1-C1-C2-O2
5	B	323	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	305	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	180/202 (89%)	-0.04	1 (0%) 85 86	19, 40, 78, 104	5 (2%)
1	B	180/202 (89%)	0.03	7 (3%) 43 45	18, 38, 86, 110	5 (2%)
All	All	360/404 (89%)	-0.00	8 (2%) 62 64	18, 40, 84, 110	10 (2%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	12	THR	4.3
1	B	11	THR	3.6
1	B	25	HIS	3.2
1	B	62	CYS	3.0
1	A	137	SER	2.7
1	B	16	HIS	2.5
1	B	61	GLN	2.4
1	B	26	ILE	2.3

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
6	ACY	B	312	4/4	0.38	0.23	110,113,113,113	0
6	ACY	B	311	4/4	0.46	0.16	95,96,97,98	0
6	ACY	B	316	4/4	0.49	0.15	89,94,95,96	0
6	ACY	B	317	4/4	0.63	0.11	108,110,111,112	0
5	GOL	B	321	6/6	0.64	0.13	96,101,102,104	0
5	GOL	A	326	6/6	0.65	0.13	99,109,109,111	0
5	GOL	B	324	6/6	0.69	0.14	105,109,110,111	0
5	GOL	B	322	6/6	0.70	0.09	89,94,95,95	0
5	GOL	B	327	6/6	0.70	0.10	79,87,89,89	0
5	GOL	B	325	6/6	0.73	0.11	90,92,94,95	0
4	SO4	A	305	5/5	0.76	0.11	118,120,120,120	0
6	ACY	B	315	4/4	0.78	0.13	102,103,104,105	0
4	SO4	B	303	5/5	0.78	0.16	87,92,96,98	0
6	ACY	B	313	4/4	0.78	0.13	97,98,99,99	0
5	GOL	B	323	6/6	0.84	0.12	69,72,73,73	0
4	SO4	A	304	5/5	0.84	0.15	90,93,98,99	0
4	SO4	A	301	5/5	0.86	0.08	72,76,80,80	0
6	ACY	B	314	4/4	0.88	0.08	53,54,54,56	0
4	SO4	B	302	5/5	0.93	0.09	59,67,70,74	0
3	EMC	A	5502	3/3	0.97	0.13	79,79,80,92	0
2	HG	A	5503	1/1	0.98	0.06	82,82,82,82	0
3	EMC	B	9502	3/3	0.98	0.10	55,55,63,79	0
2	HG	B	9503	1/1	0.98	0.06	77,77,77,77	0
3	EMC	A	5501	3/3	0.99	0.07	34,34,56,72	0
3	EMC	B	9501	3/3	0.99	0.08	60,60,62,70	0

6.5 Other polymers ⓘ

There are no such residues in this entry.