



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:40 PM UTC

PDB ID : 1FX7 / pdb_00001fx7
Title : CRYSTAL STRUCTURE OF THE IRON-DEPENDENT REGULATOR (IDER) FROM MYCOBACTERIUM TUBERCULOSIS
Authors : Feese, M.D.; Ingason, B.P.; Goranson-Siekierke, J.; Holmes, R.K.; Hol, W.J.G.
Deposited on : 2000-09-25
Resolution : 2.00 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

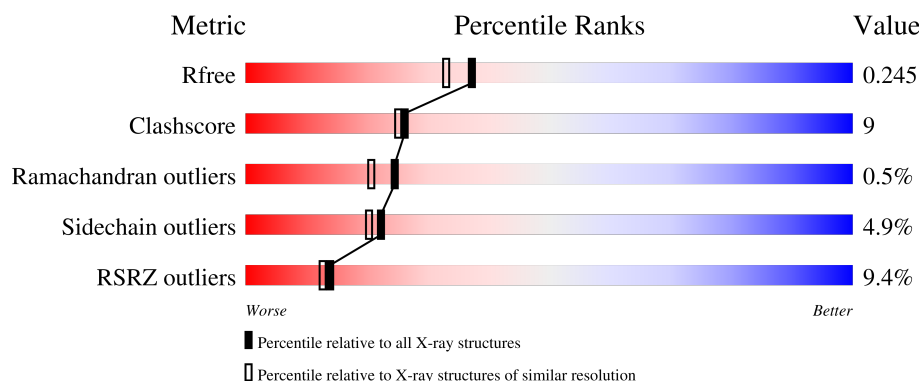
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	230	8% 82% 17% .
1	B	230	7% 78% 15% ..
1	C	230	17% 70% 23% ..
1	D	230	4% 69% 25% ..

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
2	SO4	B	1005	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7817 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 IRON-DEPENDENT REPRESSOR IDER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	230	Total	C	N	O	S	0	3	0
			1777	1109	322	336	10			
1	B	220	Total	C	N	O	S	0	3	0
			1718	1074	311	323	10			
1	C	220	Total	C	N	O	S	0	4	0
			1716	1074	308	323	11			
1	D	220	Total	C	N	O	S	0	4	0
			1716	1074	308	323	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	136	ASP	VAL	conflict	UNP P0A672
B	136	ASP	VAL	conflict	UNP P0A672
C	136	ASP	VAL	conflict	UNP P0A672
D	136	ASP	VAL	conflict	UNP P0A672

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Co	0	0
			3	3		
3	B	3	Total	Co	0	0
			3	3		
3	C	3	Total	Co	0	0
			3	3		
3	D	4	Total	Co	0	0
			4	4		

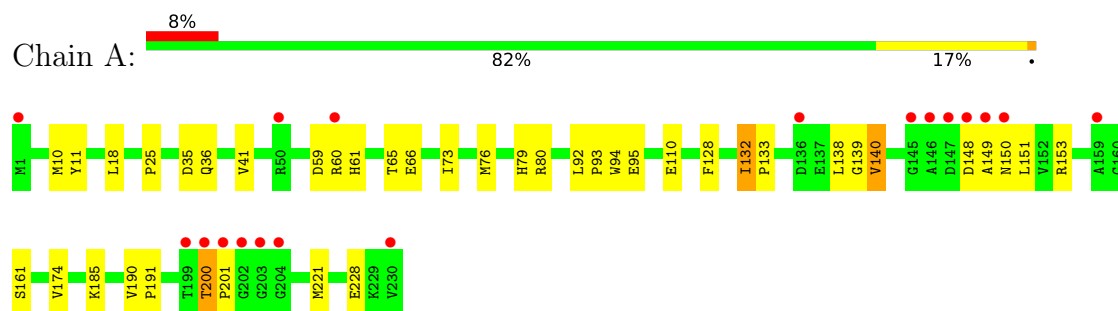
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	263	Total	O	0	0
			263	263		
4	B	208	Total	O	0	0
			208	208		
4	C	139	Total	O	0	0
			139	139		
4	D	182	Total	O	0	0
			182	182		

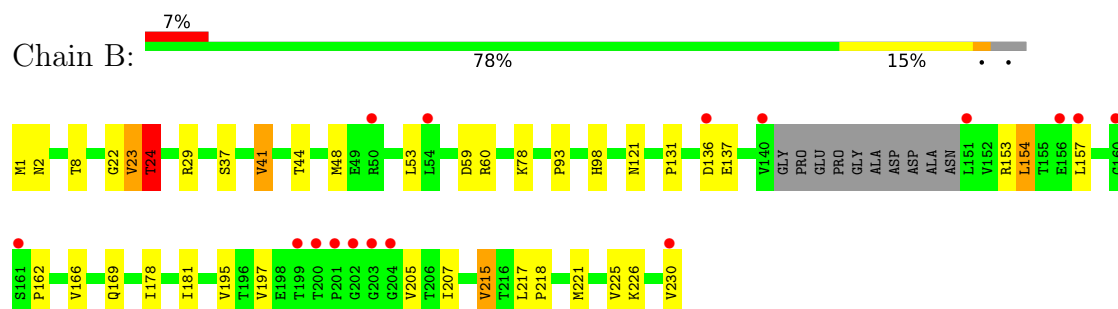
3 Residue-property plots [i](#)

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.

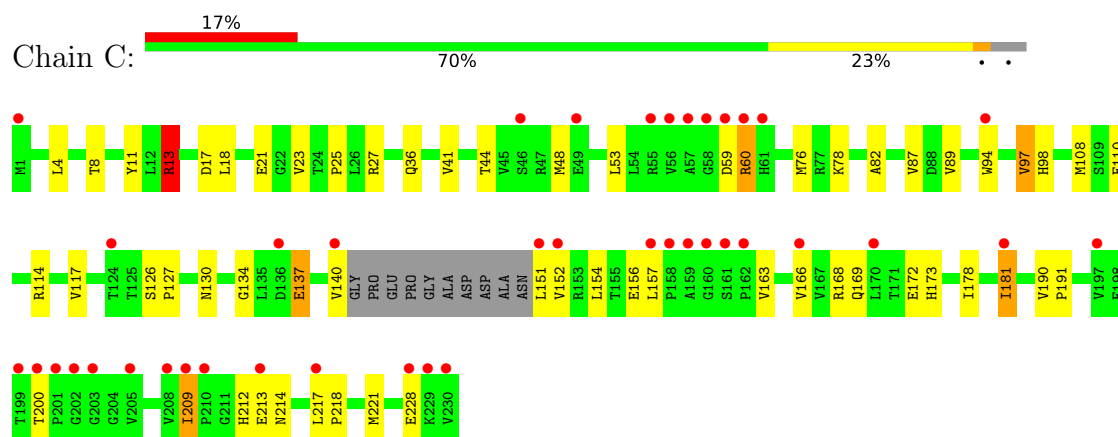
• Molecule 1: IRON-DEPENDENT REPRESSOR IDER



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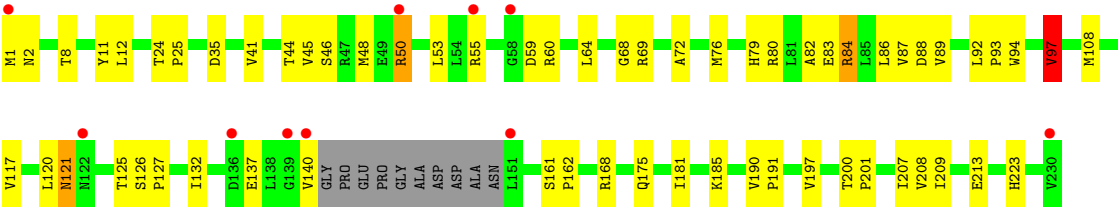


• Molecule 1: IRON-DEPENDENT REPRESSOR IDER



• Molecule 1: IRON-DEPENDENT REPRESSOR IDER





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.31Å 113.97Å 236.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.6 (20.00-2.00) 95.8 (20.00-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.87 (at 1.80Å)	Xtriage
Refinement program	REFMAC, TNT	Depositor
R, R_{free}	0.214 , 0.272 0.199 , 0.245	Depositor DCC
R_{free} test set	4839 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	29.5	Xtriage
Anisotropy	0.853	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 72.6	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	0.95	EDS
Total number of atoms	7817	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: CO, 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	0.57	0/1818	1.39	11/2474 (0.4%)
1	B	0.52	0/1756	1.30	8/2386 (0.3%)
1	C	0.47	0/1758	1.27	3/2389 (0.1%)
1	D	0.49	0/1758	1.31	6/2389 (0.3%)
All	All	0.51	0/7090	1.32	28/9638 (0.3%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	24	THR	CA-C-O	9.11	126.06	119.32
1	A	61	HIS	CA-CB-CG	8.99	122.79	113.80
1	C	98	HIS	CA-CB-CG	-7.41	106.39	113.80
1	A	128	PHE	CA-CB-CG	-7.29	106.52	113.80
1	D	208	VAL	N-CA-CB	-6.96	101.79	111.25
1	A	140	VAL	N-CA-CB	6.93	118.27	110.72
1	A	10	MET	CA-C-O	6.66	127.63	119.97
1	A	139	GLY	CA-C-N	-6.53	113.95	122.37
1	A	139	GLY	C-N-CA	-6.53	113.95	122.37
1	A	132	ILE	CA-C-O	6.33	122.66	119.12
1	D	97	VAL	N-CA-CB	-6.17	102.15	110.54
1	A	94	TRP	CA-CB-CG	6.10	125.19	113.60
1	B	217	LEU	CB-CA-C	-5.62	103.44	109.85
1	A	66	GLU	CA-C-N	5.58	127.69	120.44
1	A	66	GLU	C-N-CA	5.58	127.69	120.44
1	D	97	VAL	N-CA-C	5.53	116.26	110.62
1	B	131	PRO	CA-C-N	-5.47	118.88	122.60
1	B	131	PRO	C-N-CA	-5.47	118.88	122.60
1	A	153	ARG	CA-C-O	-5.47	115.60	121.56
1	B	98	HIS	CA-CB-CG	-5.46	108.34	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	132	ILE	CA-C-O	5.37	122.32	119.15
1	D	223	HIS	CA-CB-CG	5.35	119.15	113.80
1	C	13	ARG	CG-CD-NE	5.31	123.67	112.00
1	C	13	ARG	CB-CG-CD	5.27	123.42	111.30
1	B	23	VAL	CA-C-O	-5.14	116.22	121.67
1	D	168	ARG	CD-NE-CZ	5.10	131.53	124.40
1	B	93	PRO	CA-C-N	5.01	127.31	120.54
1	B	93	PRO	C-N-CA	5.01	127.31	120.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1777	0	1806	25	0
1	B	1718	0	1759	26	0
1	C	1716	0	1753	43	0
1	D	1716	0	1753	39	0
2	A	30	0	0	0	0
2	B	25	0	0	2	0
2	C	20	0	0	0	0
2	D	10	0	0	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
3	D	4	0	0	0	0
4	A	263	0	0	1	0
4	B	208	0	0	3	0
4	C	139	0	0	2	0
4	D	182	0	0	1	0
All	All	7817	0	7071	132	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 9.

All (132) 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:48[A]:MET:HG2	1:B:53:LEU:HD12	1.45	0.95
1:A:76[A]:MET:HE2	1:A:80:ARG:HG3	1.48	0.95
1:D:76[A]:MET:HE2	1:D:80:ARG:NE	1.88	0.89
1:C:209:ILE:HD11	1:C:212:HIS:HB2	1.55	0.89
1:A:76[A]:MET:CE	1:A:80:ARG:HG3	2.07	0.84
1:D:11:TYR:HB2	1:D:48[B]:MET:HE1	1.61	0.80
1:A:76[A]:MET:HE3	1:A:79:HIS:HB3	1.65	0.79
1:A:76[A]:MET:HE2	1:A:80:ARG:CG	2.13	0.78
1:C:130:ASN:HD22	1:C:169:GLN:HE21	1.28	0.78
1:A:95:GLU:OE2	1:A:185:LYS:HD3	1.87	0.74
1:B:59:ASP:O	1:B:60:ARG:HB2	1.88	0.73
1:C:11:TYR:HB2	1:C:48[B]:MET:HE1	1.70	0.71
1:D:83:GLU:HG2	1:D:97:VAL:HG13	1.74	0.70
1:A:174:VAL:HA	1:A:221[B]:MET:HE2	1.75	0.69
1:D:59:ASP:O	1:D:60:ARG:HB2	1.92	0.69
1:B:48[A]:MET:HG2	1:B:53:LEU:CD1	2.23	0.68
1:A:76[A]:MET:HE2	1:A:80:ARG:NE	2.09	0.67
1:A:76[A]:MET:HE1	1:A:79:HIS:CD2	2.31	0.66
1:C:89:VAL:HG11	1:D:89:VAL:HG11	1.78	0.64
1:C:209:ILE:H	1:C:214:ASN:ND2	1.95	0.64
1:A:59:ASP:O	1:A:60:ARG:HB2	1.97	0.62
1:C:17:ASP:O	1:C:21:GLU:HG2	1.99	0.62
1:A:18:LEU:HD13	1:A:25:PRO:HA	1.83	0.60
1:C:48[B]:MET:HG2	1:C:53:LEU:HD12	1.82	0.60
1:C:87:VAL:HG23	1:C:97:VAL:HG21	1.82	0.60
1:C:130:ASN:HD22	1:C:169:GLN:NE2	1.99	0.60
1:C:76[A]:MET:HE1	1:C:172:GLU:HG3	1.82	0.60
1:C:154:LEU:HD12	1:C:217:LEU:HD22	1.83	0.60
1:C:18:LEU:HD13	1:C:25:PRO:HA	1.83	0.59
1:D:200:THR:HB	1:D:201:PRO:CD	2.33	0.58
1:C:59:ASP:O	1:C:60:ARG:HB2	2.02	0.58
1:B:1:MET:HE2	1:B:2:ASN:O	2.03	0.58
1:D:87:VAL:HG13	1:D:97:VAL:HG21	1.86	0.58
1:B:22:GLY:HA3	1:B:153:ARG:NH1	2.20	0.57
1:D:200:THR:HB	1:D:201:PRO:HD2	1.86	0.57
1:A:76[A]:MET:HE2	1:A:80:ARG:CD	2.35	0.57
1:C:11:TYR:OH	1:C:36:GLN:NE2	2.38	0.57
1:A:200:THR:HG23	1:A:201:PRO:HD2	1.87	0.57
1:C:209:ILE:H	1:C:214:ASN:HD22	1.53	0.57
1:C:218:PRO:HD2	1:C:221[B]:MET:HG3	1.87	0.55
1:D:190:VAL:HB	1:D:191:PRO:HD2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:117:VAL:HG11	1:C:140:VAL:HG11	1.87	0.55
1:A:190:VAL:HB	1:A:191:PRO:HD2	1.89	0.54
1:D:8:THR:HG23	1:D:48[B]:MET:HE2	1.89	0.54
1:C:178:ILE:O	1:C:181:ILE:HG22	2.08	0.54
1:D:44:THR:HG22	1:D:48[B]:MET:HE3	1.89	0.53
1:C:134:GLY:HA2	1:C:137:GLU:OE1	2.09	0.53
1:C:8:THR:HA	1:C:48[B]:MET:HE2	1.91	0.52
1:C:117:VAL:HG11	1:C:140:VAL:HG21	1.92	0.52
1:C:152:VAL:CG2	1:C:156:GLU:HB2	2.40	0.52
1:B:23:VAL:HG23	2:B:1005:SO4:S	2.50	0.51
1:D:44:THR:CG2	1:D:48[B]:MET:HE3	2.40	0.51
1:D:94:TRP:CE3	1:D:185:LYS:HB2	2.46	0.51
1:A:174:VAL:CA	1:A:221[B]:MET:HE2	2.40	0.50
1:A:190:VAL:HB	1:A:191:PRO:CD	2.41	0.50
1:D:161:SER:HB3	1:D:162:PRO:HD2	1.92	0.50
1:A:11:TYR:OH	1:A:36:GLN:NE2	2.44	0.50
1:B:166:VAL:HG23	1:B:230:VAL:HG22	1.94	0.50
1:B:162:PRO:HA	1:B:197:VAL:O	2.12	0.50
1:D:46:SER:O	1:D:50:ARG:HG2	2.11	0.50
1:D:84:ARG:HD2	1:D:125:THR:O	2.12	0.50
1:B:178:ILE:HA	1:B:181:ILE:HG22	1.94	0.49
1:D:8:THR:HA	1:D:48[B]:MET:CE	2.42	0.49
1:C:173:HIS:HD2	4:C:2026:HOH:O	1.96	0.49
1:B:44:THR:HG22	1:B:48[A]:MET:CE	2.43	0.49
1:D:175:GLN:HA	1:D:181:ILE:HG13	1.94	0.49
1:B:48[A]:MET:CG	1:B:53:LEU:HD12	2.32	0.48
1:B:169:GLN:HB2	1:B:226:LYS:HB2	1.95	0.48
1:C:27:ARG:NH2	1:C:60:ARG:HG2	2.28	0.48
1:B:207:ILE:HB	1:B:215:VAL:HG13	1.95	0.48
1:C:11:TYR:CB	1:C:48[B]:MET:HE1	2.42	0.48
1:B:37:SER:O	1:B:41:VAL:HG13	2.14	0.47
1:D:120:LEU:O	1:D:121:ASN:HB2	2.15	0.47
1:B:23:VAL:HG23	2:B:1005:SO4:O3	2.15	0.47
1:D:12:LEU:HB3	1:D:72:ALA:HB2	1.97	0.47
1:C:78:LYS:HE3	4:C:2048:HOH:O	2.14	0.47
1:C:110:GLU:OE1	1:C:114:ARG:NH2	2.47	0.47
1:D:41:VAL:O	1:D:45:VAL:HG23	2.14	0.47
1:D:76[A]:MET:HE2	1:D:80:ARG:CZ	2.45	0.47
1:D:87:VAL:CG1	1:D:97:VAL:HG21	2.45	0.47
1:D:190:VAL:HB	1:D:191:PRO:CD	2.44	0.47
1:D:1:MET:HE2	1:D:2:ASN:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:76[A]:MET:HE1	1:C:172:GLU:CG	2.44	0.46
1:B:154:LEU:HD22	1:B:225:VAL:HB	1.96	0.46
1:C:82:ALA:HB2	1:C:108[B]:MET:SD	2.56	0.46
1:A:65:THR:HB	4:A:2180:HOH:O	2.14	0.46
1:C:8:THR:HA	1:C:48[B]:MET:CE	2.46	0.46
1:A:190:VAL:O	1:A:191:PRO:C	2.59	0.46
1:C:44:THR:HG22	1:C:48[B]:MET:HE3	1.98	0.45
1:A:80:ARG:HB2	1:A:132:ILE:HG12	1.97	0.45
1:D:117:VAL:HG11	1:D:140:VAL:HG21	1.97	0.45
1:B:195:VAL:HG21	1:B:207:ILE:HG23	1.99	0.45
1:D:11:TYR:CB	1:D:48[B]:MET:HE1	2.39	0.45
1:C:130:ASN:ND2	1:C:169:GLN:HE21	2.04	0.45
1:A:151:LEU:HD12	1:A:228:GLU:HG3	1.99	0.45
1:C:190:VAL:HB	1:C:191:PRO:HD2	1.98	0.45
1:C:44:THR:CG2	1:C:48[B]:MET:HE3	2.48	0.44
1:C:166:VAL:HG23	1:C:228:GLU:HB3	2.00	0.44
1:C:13:ARG:HD2	1:C:76[A]:MET:SD	2.58	0.44
1:C:152:VAL:HG23	1:C:156:GLU:HB2	2.00	0.44
1:D:8:THR:HA	1:D:48[B]:MET:HE2	1.99	0.44
1:C:130:ASN:ND2	1:C:169:GLN:HB3	2.32	0.44
1:B:23:VAL:HG21	4:B:2182:HOH:O	2.17	0.44
1:D:197:VAL:HG12	1:D:207:ILE:HD13	2.00	0.43
1:A:92:LEU:HD12	1:A:93:PRO:HD2	1.99	0.43
1:D:24:THR:HA	1:D:25:PRO:HD3	1.87	0.43
1:D:87:VAL:HG23	1:D:88:ASP:N	2.34	0.43
1:D:82:ALA:HB2	1:D:108[B]:MET:SD	2.58	0.43
1:D:83:GLU:O	1:D:87:VAL:HG22	2.19	0.43
1:B:23:VAL:HG22	1:B:24:THR:N	2.33	0.43
1:D:86:LEU:HB3	1:D:97:VAL:HG22	2.00	0.43
1:B:218:PRO:HD2	1:B:221[B]:MET:HG3	2.01	0.42
1:A:174:VAL:HA	1:A:221[B]:MET:CE	2.46	0.42
1:B:154:LEU:O	1:B:205:VAL:HG21	2.19	0.42
1:C:59:ASP:O	1:C:60:ARG:CB	2.67	0.42
1:A:73:ILE:HG23	1:A:133:PRO:HB2	2.02	0.42
1:B:195:VAL:HG23	4:B:2103:HOH:O	2.18	0.42
1:C:126:SER:HB2	1:C:127:PRO:HD2	2.01	0.42
1:D:126:SER:HB2	1:D:127:PRO:CD	2.51	0.41
1:D:76[A]:MET:HE3	1:D:79:HIS:HB3	2.02	0.41
1:D:69:ARG:HD3	4:D:2192:HOH:O	2.20	0.41
1:D:53:LEU:O	1:D:68:GLY:HA3	2.21	0.41
1:A:110:GLU:HG3	1:A:138:LEU:CD2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:181:ILE:HG21	4:B:2031:HOH:O	2.20	0.41
1:C:166:VAL:CG2	1:C:228:GLU:HB3	2.51	0.41
1:C:190:VAL:HB	1:C:191:PRO:CD	2.50	0.41
1:D:92:LEU:HD12	1:D:93:PRO:HD2	2.03	0.41
1:B:8:THR:HG23	1:B:48[A]:MET:SD	2.61	0.40
1:B:197:VAL:HG12	1:B:207:ILE:HD13	2.04	0.40
1:C:151:LEU:HD11	1:C:168:ARG:HB2	2.03	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	231/230 (100%)	222 (96%)	7 (3%)	2 (1%)	14	9
1	B	219/230 (95%)	214 (98%)	5 (2%)	0	100	100
1	C	220/230 (96%)	214 (97%)	5 (2%)	1 (0%)	24	21
1	D	220/230 (96%)	213 (97%)	6 (3%)	1 (0%)	24	21
All	All	890/920 (97%)	863 (97%)	23 (3%)	4 (0%)	24	27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	148	ASP
1	A	149	ALA
1	C	60	ARG
1	D	121	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	196/195 (100%)	190 (97%)	6 (3%)	35	37
1	B	192/195 (98%)	183 (95%)	9 (5%)	23	22
1	C	192/195 (98%)	179 (93%)	13 (7%)	14	11
1	D	192/195 (98%)	183 (95%)	9 (5%)	23	22
All	All	772/780 (99%)	735 (95%)	37 (5%)	22	21

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ASP
1	A	41	VAL
1	A	140	VAL
1	A	150	ASN
1	A	161	SER
1	A	200	THR
1	B	24	THR
1	B	29	ARG
1	B	41	VAL
1	B	121	ASN
1	B	136	ASP
1	B	137	GLU
1	B	154	LEU
1	B	157	LEU
1	B	215	VAL
1	C	4	LEU
1	C	13	ARG
1	C	23	VAL
1	C	41	VAL
1	C	94	TRP
1	C	97	VAL
1	C	137	GLU
1	C	157	LEU
1	C	163	VAL

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Mol	Chain	Res	Type
1	C	181	ILE
1	C	200	THR
1	C	209	ILE
1	C	213	GLU
1	D	35	ASP
1	D	50	ARG
1	D	55	ARG
1	D	64	LEU
1	D	84	ARG
1	D	97	VAL
1	D	137	GLU
1	D	209	ILE
1	D	213	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	121	ASN
1	A	150	ASN
1	A	169	GLN
1	B	36	GLN
1	C	36	GLN
1	C	43	GLN
1	C	122	ASN
1	C	130	ASN
1	C	169	GLN
1	C	173	HIS
1	C	214	ASN
1	D	36	GLN
1	D	43	GLN
1	D	121	ASN

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 30 ligands modelled in this entry, 13 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	$\# Z > 2$
2	SO4	B	1005	-	4,4,4	0.68	0	6,6,6	0.28	0
2	SO4	A	1001	-	4,4,4	0.63	0	6,6,6	0.27	0
2	SO4	D	1012	-	4,4,4	0.67	0	6,6,6	0.12	0
2	SO4	D	1015	-	4,4,4	0.69	0	6,6,6	0.10	0
2	SO4	A	1017	-	4,4,4	0.71	0	6,6,6	0.08	0
2	SO4	B	1009	-	4,4,4	0.75	0	6,6,6	0.29	0
2	SO4	C	1007	-	4,4,4	0.68	0	6,6,6	0.22	0
2	SO4	C	1003	-	4,4,4	0.70	0	6,6,6	0.34	0
2	SO4	A	1013	-	4,4,4	0.74	0	6,6,6	0.27	0
2	SO4	A	1006	-	4,4,4	0.73	0	6,6,6	0.13	0
2	SO4	B	1011	-	4,4,4	0.68	0	6,6,6	0.14	0
2	SO4	B	1002	-	4,4,4	0.72	0	6,6,6	0.24	0
2	SO4	B	1014	-	4,4,4	0.69	0	6,6,6	0.24	0
2	SO4	C	1016	-	4,4,4	0.71	0	6,6,6	0.09	0
2	SO4	A	1008	-	4,4,4	0.72	0	6,6,6	0.17	0
2	SO4	C	1010	-	4,4,4	0.72	0	6,6,6	0.16	0
2	SO4	A	1004	-	4,4,4	0.55	0	6,6,6	0.32	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1005	SO4	2	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	230/230 (100%)	0.06	18 (7%)	19 18	15, 30, 67, 134	3 (1%)
1	B	220/230 (95%)	0.31	16 (7%)	21 19	16, 36, 74, 166	3 (1%)
1	C	220/230 (95%)	0.91	40 (18%)	3 3	22, 47, 108, 183	4 (1%)
1	D	220/230 (95%)	0.26	10 (4%)	38 37	20, 39, 68, 96	4 (1%)
All	All	890/920 (96%)	0.38	84 (9%)	14 13	15, 38, 79, 183	14 (1%)

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	60	ARG	6.3
1	C	57	ALA	5.8
1	C	56	VAL	5.6
1	B	203	GLY	5.5
1	C	160	GLY	5.4
1	C	140	VAL	5.3
1	B	202	GLY	5.1
1	C	55	ARG	5.0
1	C	230	VAL	4.9
1	A	149	ALA	4.8
1	A	148	ASP	4.7
1	B	160	GLY	4.5
1	C	49	GLU	4.5
1	B	201	PRO	4.5
1	C	58	GLY	4.4
1	B	200	THR	4.4
1	A	150	ASN	4.3
1	C	136	ASP	4.3
1	A	1	MET	4.1
1	C	158	PRO	4.1
1	C	159	ALA	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	201	PRO	3.9
1	B	140	VAL	3.9
1	C	202	GLY	3.8
1	C	157	LEU	3.7
1	C	59	ASP	3.7
1	B	204	GLY	3.7
1	C	151	LEU	3.6
1	C	161	SER	3.6
1	C	200	THR	3.5
1	D	1	MET	3.4
1	D	230	VAL	3.4
1	C	1	MET	3.4
1	C	201	PRO	3.4
1	B	54	LEU	3.4
1	A	146	ALA	3.2
1	B	50	ARG	3.2
1	B	230	VAL	3.2
1	D	58	GLY	3.1
1	D	140	VAL	3.0
1	A	203	GLY	3.0
1	D	139	GLY	3.0
1	A	202	GLY	2.9
1	A	204	GLY	2.9
1	A	147	ASP	2.9
1	A	136	ASP	2.8
1	C	199	THR	2.8
1	B	161	SER	2.8
1	A	145	GLY	2.8
1	C	209	ILE	2.8
1	D	55	ARG	2.7
1	A	200	THR	2.7
1	B	199	THR	2.7
1	A	199	THR	2.6
1	B	157	LEU	2.6
1	D	122	ASN	2.5
1	D	136	ASP	2.5
1	D	151	LEU	2.5
1	A	230	VAL	2.4
1	C	94	TRP	2.4
1	C	166	VAL	2.4
1	D	50	ARG	2.4
1	A	159	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	152	VAL	2.4
1	C	208	VAL	2.3
1	B	151	LEU	2.3
1	A	50	ARG	2.2
1	A	60	ARG	2.2
1	C	213	GLU	2.2
1	C	124	THR	2.2
1	C	162	PRO	2.1
1	C	181	ILE	2.1
1	C	197	VAL	2.1
1	C	61	HIS	2.1
1	C	229	LYS	2.1
1	C	217	LEU	2.1
1	B	156	GLU	2.1
1	C	46	SER	2.1
1	C	170	LEU	2.0
1	C	205	VAL	2.0
1	C	228	GLU	2.0
1	B	136	ASP	2.0
1	C	210	PRO	2.0
1	C	203	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	C	1016	5/5	0.28	0.17	201,202,202,203	0
2	SO4	A	1017	5/5	0.73	0.11	94,100,103,103	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	B	1014	5/5	0.79	0.12	48,64,75,80	0
2	SO4	A	1013	5/5	0.81	0.19	38,57,66,78	0
2	SO4	D	1015	5/5	0.83	0.08	83,86,90,95	0
2	SO4	A	1008	5/5	0.87	0.11	46,61,71,74	0
2	SO4	B	1002	5/5	0.89	0.11	62,66,66,69	0
2	SO4	C	1007	5/5	0.90	0.12	50,50,63,66	0
2	SO4	B	1011	5/5	0.90	0.09	55,60,69,72	0
2	SO4	A	1006	5/5	0.90	0.18	41,58,71,75	0
2	SO4	C	1010	5/5	0.91	0.12	54,55,66,67	0
2	SO4	B	1005	5/5	0.93	0.10	52,58,65,71	0
2	SO4	A	1001	5/5	0.94	0.11	37,44,55,58	0
2	SO4	D	1012	5/5	0.94	0.10	44,48,60,65	0
2	SO4	C	1003	5/5	0.94	0.10	41,50,56,60	0
2	SO4	B	1009	5/5	0.96	0.08	42,45,48,59	0
3	CO	C	2011	1/1	0.98	0.05	44,44,44,44	0
3	CO	D	2013	1/1	0.98	0.04	49,49,49,49	0
3	CO	B	2010	1/1	0.99	0.02	44,44,44,44	0
3	CO	C	2005	1/1	0.99	0.02	32,32,32,32	0
3	CO	C	2006	1/1	0.99	0.02	31,31,31,31	0
2	SO4	A	1004	5/5	0.99	0.04	30,31,38,39	0
3	CO	D	2012	1/1	0.99	0.02	39,39,39,39	0
3	CO	B	2003	1/1	0.99	0.02	27,27,27,27	0
3	CO	A	2001	1/1	1.00	0.02	21,21,21,21	0
3	CO	B	2004	1/1	1.00	0.01	26,26,26,26	0
3	CO	D	2007	1/1	1.00	0.01	30,30,30,30	0
3	CO	D	2008	1/1	1.00	0.02	31,31,31,31	0
3	CO	A	2002	1/1	1.00	0.04	23,23,23,23	0
3	CO	A	2009	1/1	1.00	0.01	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.