



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 06:39 PM GMT

PDB ID : 2B3Y
Title : Structure of a monoclinic crystal form of human cytosolic aconitase (IRP1)
Authors : Dupuy, J.; Fontecilla-Camps, J.C.; Volbeda, A.
Deposited on : 2005-09-22
Resolution : 1.85 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

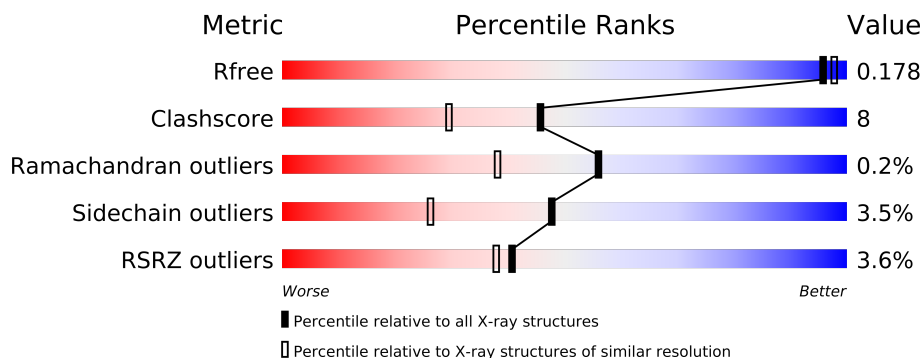
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.85 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}	66092	1269 (1.86-1.86)
Clashscore	79885	1470 (1.86-1.86)
Ramachandran outliers	78287	1451 (1.86-1.86)
Sidechain outliers	78261	1451 (1.86-1.86)
RSRZ outliers	66119	1269 (1.86-1.86)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	888	
1	B	888	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	ACT	A	890	-	X
2	ACT	B	891	-	X
4	GOL	A	1001	-	X
4	GOL	A	1002	-	X

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 14870 atoms, of which 0 are hydrogen and 0 are deuterium.

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-responsive element binding protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	888	Total	C	N	O	S	0	1	0
			6929	4431	1183	1286	29			
1	B	888	Total	C	N	O	S	0	0	0
			6928	4431	1183	1285	29			

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

- Molecule 3 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe₄S₄).



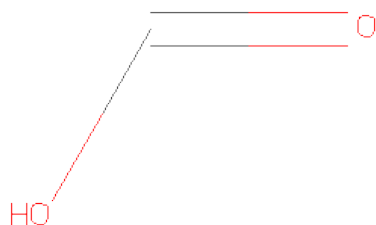
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	S	0	0
			8	4	4		
3	B	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



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

- Molecule 6 is water.

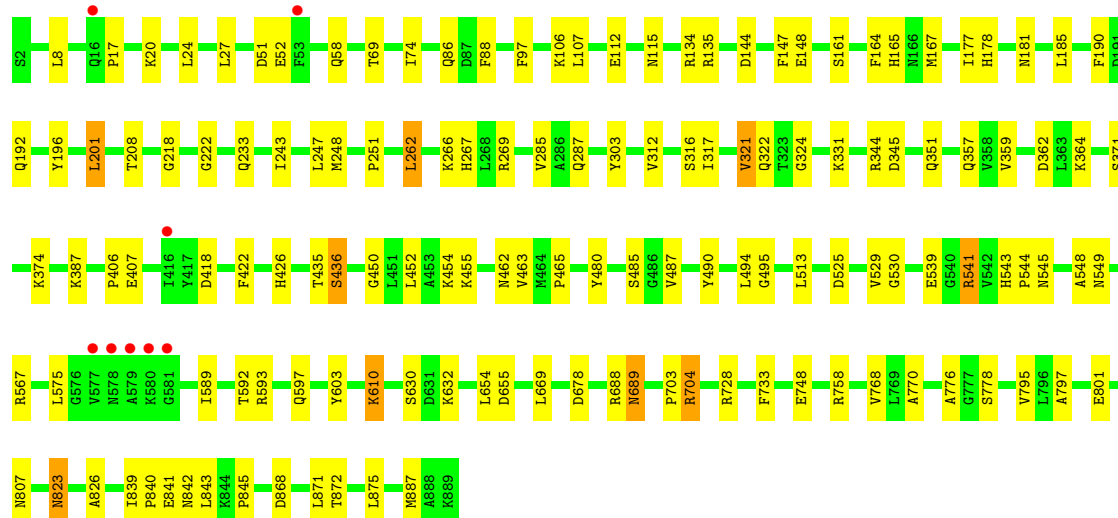
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	520	Total	O	0	0
			520	520		
6	B	450	Total	O	0	0
			450	450		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

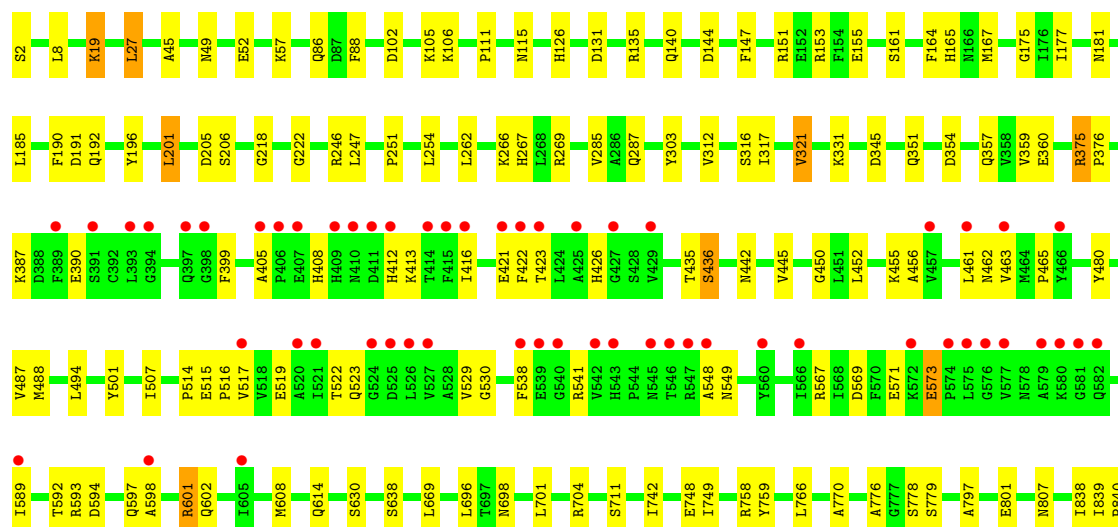
- Molecule 1: Iron-responsive element binding protein 1

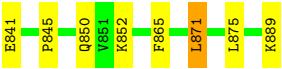
Chain A: 



- Molecule 1: Iron-responsive element binding protein 1

Chain B: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.91Å 215.47Å 64.01Å 90.00° 72.00° 90.00°	Depositor
Resolution (Å)	29.50 – 1.85 29.29 – 1.85	Depositor EDS
% Data completeness (in resolution range)	94.0 (29.50-1.85) 98.8 (29.29-1.85)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.86 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.211 , 0.261 0.157 , 0.178	Depositor DCC
R_{free} test set	6949 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	20.0	Xtriage
Anisotropy	0.601	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 63.6	EDS
Estimated twinning fraction	0.408 for -l,-k,-h	Xtriage
L-test for twinning	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Outliers	0 of 138110 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14870	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SF4, FMT, ACT

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	0/7095	0.65	1/9634 (0.0%)
1	B	0.54	0/7089	0.62	1/9626 (0.0%)
All	All	0.54	0/14184	0.64	2/19260 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	887	MET	CG-SD-CE	5.71	109.33	100.20
1	B	871	LEU	CA-CB-CG	5.04	126.89	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6929	0	6921	107	0
1	B	6928	0	6920	115	0
2	A	4	0	3	2	0
2	B	8	0	6	11	0
3	A	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	8	0	0	0	0
4	A	12	0	16	1	0
5	B	3	0	1	10	0
6	A	520	0	0	21	0
6	B	450	0	0	15	0
All	All	14870	0	13867	229	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (229) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:891:ACT:CH3	5:B:1001:FMT:H	1.72	1.17
1:B:450:GLY:HA3	1:B:487:VAL:HG11	1.33	1.09
1:A:450:GLY:HA3	1:A:487:VAL:HG11	1.36	1.08
1:B:450:GLY:CA	1:B:487:VAL:HG11	1.85	1.06
1:A:450:GLY:CA	1:A:487:VAL:HG11	1.85	1.05
2:B:891:ACT:CH3	5:B:1001:FMT:C	2.37	1.03
2:B:891:ACT:H3	5:B:1001:FMT:H	1.41	0.98
2:B:890:ACT:H2	2:B:891:ACT:H2	1.46	0.94
1:A:222:GLY:HA3	1:A:807:ASN:HD21	1.35	0.91
2:B:891:ACT:H2	5:B:1001:FMT:H	1.50	0.91
1:B:345:ASP:H	1:B:351:GLN:HE22	1.22	0.83
1:B:222:GLY:HA3	1:B:807:ASN:HD21	1.45	0.82
2:B:891:ACT:H2	5:B:1001:FMT:C	2.05	0.81
1:B:845:PRO:HD3	6:B:1384:HOH:O	1.80	0.80
1:A:539:GLU:HG3	1:A:728:ARG:HD3	1.62	0.79
1:B:838:ILE:HD11	1:B:852:LYS:HD3	1.65	0.77
1:B:748:GLU:OE1	1:B:758:ARG:NH1	2.16	0.77
1:B:102:ASP:O	1:B:106:LYS:HG2	1.85	0.76
1:A:135:ARG:NH2	6:A:1104:HOH:O	2.18	0.76
2:B:890:ACT:H2	2:B:891:ACT:CH3	2.16	0.76
1:B:450:GLY:CA	1:B:487:VAL:CG1	2.65	0.74
1:B:598:ALA:O	1:B:602:GLN:HG2	1.88	0.73
1:A:345:ASP:H	1:A:351:GLN:HE22	1.36	0.73
1:A:112:GLU:OE2	1:A:632:LYS:HD2	1.87	0.72
1:A:842:ASN:HB3	6:A:1519:HOH:O	1.90	0.72
1:A:58:GLN:OE1	6:A:1017:HOH:O	2.08	0.71
1:A:251:PRO:O	1:A:593:ARG:NH2	2.22	0.71
1:B:375:ARG:HD3	1:B:399:PHE:CD2	2.26	0.71
1:B:865:PHE:HB2	1:B:871:LEU:HD23	1.72	0.71
1:B:593:ARG:HG2	1:B:597:GLN:NE2	2.06	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:742:ILE:HG12	1:B:749:ILE:CD1	2.22	0.70
1:B:865:PHE:CB	1:B:871:LEU:HD23	2.22	0.70
1:B:742:ILE:HG12	1:B:749:ILE:HD12	1.75	0.69
1:A:841:GLU:HB3	6:A:1416:HOH:O	1.93	0.69
1:A:450:GLY:C	1:A:487:VAL:HG11	2.13	0.69
1:B:450:GLY:C	1:B:487:VAL:HG11	2.15	0.68
1:B:698:ASN:ND2	6:B:1315:HOH:O	2.27	0.67
1:B:779:SER:H	5:B:1001:FMT:C	2.08	0.66
1:A:324:GLY:O	1:A:610:LYS:HE3	1.95	0.66
1:A:344:ARG:HH11	1:A:351:GLN:HE21	1.44	0.65
1:B:421:GLU:N	1:B:421:GLU:OE1	2.29	0.65
4:A:1001:GOL:H2	6:A:1194:HOH:O	1.95	0.65
1:A:733:PHE:O	6:A:1365:HOH:O	2.14	0.65
1:A:454:LYS:HD2	1:A:490:TYR:CE1	2.33	0.64
1:B:357:GLN:OE1	6:B:1148:HOH:O	2.15	0.64
1:A:422:PHE:CZ	1:A:465:PRO:HD2	2.33	0.64
1:B:455:LYS:CB	1:B:589:ILE:HG23	2.29	0.63
1:A:776:ALA:HA	1:A:801:GLU:HG3	1.81	0.63
1:A:262:LEU:HG	1:A:371:SER:OG	1.97	0.63
1:A:455:LYS:HD2	1:A:589:ILE:HA	1.80	0.62
1:A:450:GLY:CA	1:A:487:VAL:CG1	2.70	0.62
1:A:655:ASP:OD2	6:A:1345:HOH:O	2.16	0.62
1:B:838:ILE:HD12	1:B:838:ILE:N	2.16	0.61
1:B:205:ASP:HA	5:B:1001:FMT:O1	2.00	0.60
1:B:115:ASN:ND2	1:B:165:HIS:H	1.99	0.60
1:B:375:ARG:NH1	6:B:1199:HOH:O	2.22	0.60
1:B:569:ASP:OD2	1:B:571:GLU:HG2	2.00	0.60
2:B:891:ACT:H3	5:B:1001:FMT:C	2.21	0.60
1:B:456:ALA:HA	1:B:589:ILE:HD13	1.82	0.60
1:A:107:LEU:HD13	1:A:654:LEU:HD13	1.84	0.60
1:B:131:ASP:O	6:B:1057:HOH:O	2.17	0.59
1:B:461:LEU:HD12	1:B:589:ILE:CD1	2.32	0.59
1:B:455:LYS:HB3	1:B:589:ILE:HG23	1.85	0.58
1:B:405:ALA:HB3	1:B:408:HIS:HD2	1.67	0.58
1:A:267:HIS:CE1	1:A:359:VAL:CG1	2.86	0.58
1:B:52:GLU:OE1	1:B:57:LYS:N	2.36	0.58
1:B:267:HIS:CE1	1:B:359:VAL:CG1	2.87	0.58
1:A:247:LEU:HD12	1:A:285:VAL:HG22	1.85	0.57
2:B:890:ACT:CH3	2:B:891:ACT:H2	2.29	0.57
1:A:106:LYS:HD2	6:A:1092:HOH:O	2.05	0.57
1:A:688:ARG:CZ	1:A:703:PRO:HG2	2.35	0.56
1:A:192:GLN:HE22	1:B:630:SER:HB2	1.70	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:345:ASP:H	1:B:351:GLN:NE2	1.96	0.56
1:B:201:LEU:O	1:B:218:GLY:HA2	2.06	0.56
1:A:115:ASN:ND2	1:A:165:HIS:H	2.04	0.56
1:A:322:GLN:NE2	6:A:1161:HOH:O	2.39	0.55
1:B:106:LYS:HD3	1:B:106:LYS:N	2.21	0.55
1:A:778:SER:OG	2:A:890:ACT:H3	2.06	0.55
1:B:416:ILE:HG13	1:B:416:ILE:O	2.07	0.55
1:A:17:PRO:O	1:A:20:LYS:HE3	2.07	0.54
1:B:206:SER:N	5:B:1001:FMT:O1	2.33	0.54
1:A:823:ASN:ND2	1:A:826:ALA:H	2.04	0.54
1:B:529:VAL:HG22	1:B:548:ALA:HB3	1.89	0.54
1:A:454:LYS:NZ	6:A:1229:HOH:O	2.41	0.54
1:B:19:LYS:HE2	1:B:191:ASP:OD1	2.08	0.53
1:A:748:GLU:OE2	1:A:758:ARG:NH2	2.41	0.53
1:A:248:MET:HG3	1:A:362:ASP:HA	1.91	0.53
1:A:748:GLU:CD	1:A:758:ARG:HH12	2.12	0.53
1:A:529:VAL:HG22	1:A:548:ALA:HB3	1.89	0.53
1:A:487:VAL:HG12	1:A:487:VAL:O	2.09	0.53
1:A:539:GLU:HG2	6:A:1204:HOH:O	2.09	0.53
1:A:593:ARG:HD3	1:A:597:GLN:OE1	2.09	0.52
1:B:776:ALA:HA	1:B:801:GLU:HG3	1.91	0.52
1:B:251:PRO:O	1:B:593:ARG:NH1	2.41	0.52
1:B:105:LYS:HB2	1:B:111:PRO:HG3	1.91	0.52
1:B:597:GLN:O	1:B:601:ARG:HB2	2.09	0.52
1:A:688:ARG:NH2	1:A:703:PRO:HG2	2.24	0.52
1:B:541:ARG:O	1:B:541:ARG:HG2	2.09	0.51
1:A:243:ILE:HD12	1:A:357:GLN:HG2	1.92	0.51
1:A:190:PHE:O	1:A:196:TYR:HA	2.10	0.51
1:B:450:GLY:C	1:B:487:VAL:CG1	2.79	0.51
1:B:375:ARG:HD3	1:B:399:PHE:CG	2.45	0.51
1:B:140:GLN:NE2	1:B:144:ASP:OD1	2.43	0.51
1:A:374:LYS:HE2	6:A:1231:HOH:O	2.11	0.51
1:B:246:ARG:HB3	1:B:360:GLU:HG2	1.93	0.50
1:B:269:ARG:HG2	1:B:303:TYR:HA	1.93	0.50
1:B:450:GLY:O	1:B:487:VAL:HG13	2.11	0.50
1:A:222:GLY:HA3	1:A:807:ASN:ND2	2.16	0.50
1:B:541:ARG:NH2	6:B:1180:HOH:O	2.45	0.50
1:B:126:HIS:NE2	2:B:890:ACT:H1	2.27	0.50
1:A:406:PRO:HD2	1:A:407:GLU:OE2	2.11	0.50
1:A:454:LYS:HD2	1:A:490:TYR:CZ	2.46	0.49
1:B:190:PHE:O	1:B:196:TYR:HA	2.11	0.49
1:A:543:HIS:HE1	1:A:545:ASN:HD22	1.60	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:251:PRO:HG2	1:A:593:ARG:HH21	1.77	0.49
1:B:696:LEU:HD22	1:B:701:LEU:HD12	1.94	0.49
1:B:455:LYS:HD2	1:B:589:ILE:O	2.11	0.49
1:A:435:THR:O	1:A:436:SER:HB3	2.12	0.49
1:B:569:ASP:O	1:B:573:GLU:HG3	2.12	0.49
1:B:770:ALA:O	1:B:797:ALA:HA	2.13	0.49
1:B:614:GLN:HG3	6:B:1242:HOH:O	2.12	0.49
1:B:593:ARG:HG2	1:B:597:GLN:HE21	1.75	0.48
1:B:115:ASN:HD22	1:B:165:HIS:H	1.58	0.48
1:B:865:PHE:HB2	1:B:871:LEU:CD2	2.43	0.48
1:A:530:GLY:O	1:A:549:ASN:HA	2.14	0.48
1:A:768:VAL:HB	1:A:795:VAL:HG22	1.96	0.48
1:A:362:ASP:OD1	1:A:364:LYS:HG2	2.14	0.48
1:A:317:ILE:O	1:A:321:VAL:HG12	2.13	0.48
1:B:838:ILE:H	1:B:838:ILE:HD12	1.79	0.47
1:B:530:GLY:O	1:B:549:ASN:HA	2.14	0.47
1:B:488:MET:O	1:B:488:MET:HE3	2.15	0.47
1:B:778:SER:OG	2:B:891:ACT:H3	2.15	0.47
1:A:450:GLY:C	1:A:487:VAL:CG1	2.82	0.47
1:B:507:ILE:HG13	1:B:507:ILE:O	2.15	0.47
1:A:418:ASP:OD2	6:A:1247:HOH:O	2.20	0.47
1:B:8:LEU:HD11	1:B:27:LEU:CD1	2.45	0.47
1:B:19:LYS:HA	1:B:19:LYS:HD2	1.52	0.46
1:A:69:THR:HB	1:A:74:ILE:HD12	1.98	0.46
1:A:539:GLU:HG3	1:A:728:ARG:CD	2.39	0.46
1:B:115:ASN:ND2	1:B:164:PHE:HA	2.31	0.46
1:B:161:SER:HA	1:B:167:MET:HE2	1.97	0.46
1:B:345:ASP:N	1:B:351:GLN:HE22	2.02	0.45
1:A:201:LEU:O	1:A:218:GLY:HA2	2.15	0.45
1:B:435:THR:O	1:B:436:SER:CB	2.65	0.45
1:B:317:ILE:O	1:B:321:VAL:HG12	2.17	0.45
1:A:8:LEU:HD11	1:A:27:LEU:HD11	1.99	0.45
1:A:487:VAL:O	1:A:487:VAL:CG1	2.65	0.45
1:A:426:HIS:HB2	1:A:567:ARG:NH1	2.31	0.44
1:A:704:ARG:HG3	6:A:1500:HOH:O	2.17	0.44
1:A:689:ASN:N	1:A:689:ASN:OD1	2.47	0.44
1:B:871:LEU:O	1:B:875:LEU:HG	2.17	0.44
1:B:514:PRO:HG2	1:B:517:VAL:CG2	2.48	0.44
1:A:287:GLN:HG2	6:A:1157:HOH:O	2.16	0.44
1:A:450:GLY:O	1:A:487:VAL:HG13	2.17	0.44
1:A:704:ARG:CG	6:A:1500:HOH:O	2.66	0.44
1:B:452:LEU:C	1:B:452:LEU:HD23	2.38	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:312:VAL:HG13	1:B:316:SER:HB2	1.99	0.44
1:A:344:ARG:HD2	1:A:351:GLN:NE2	2.33	0.44
1:A:20:LYS:HD2	6:A:1024:HOH:O	2.16	0.44
1:A:463:VAL:HG23	1:A:494:LEU:HD22	1.99	0.44
1:A:839:ILE:HA	1:A:840:PRO:HD3	1.85	0.44
1:A:161:SER:HA	1:A:167:MET:HE2	2.00	0.44
1:B:507:ILE:HD12	1:B:541:ARG:NH2	2.33	0.44
1:B:181:ASN:HA	1:B:185:LEU:HB2	1.99	0.44
1:B:778:SER:OG	5:B:1001:FMT:H	2.18	0.43
1:A:845:PRO:HG3	1:A:871:LEU:HD12	2.00	0.43
1:B:839:ILE:HA	1:B:840:PRO:HD3	1.80	0.43
1:B:593:ARG:O	1:B:597:GLN:HG3	2.18	0.43
1:B:455:LYS:HB2	1:B:589:ILE:HG23	1.99	0.43
1:A:840:PRO:HD2	1:A:843:LEU:CD2	2.48	0.43
1:B:376:PRO:HD3	1:B:538:PHE:CE1	2.53	0.43
1:A:269:ARG:HG2	1:A:303:TYR:HA	2.01	0.43
1:B:704:ARG:NH1	6:B:1272:HOH:O	2.52	0.43
1:B:602:GLN:HB3	6:B:1422:HOH:O	2.17	0.43
1:A:192:GLN:HE22	1:B:630:SER:CB	2.31	0.43
1:B:515:GLU:N	1:B:516:PRO:CD	2.82	0.43
1:B:461:LEU:CD1	1:B:589:ILE:CD1	2.96	0.43
1:A:776:ALA:CA	1:A:801:GLU:HG3	2.48	0.43
1:A:452:LEU:C	1:A:452:LEU:HD23	2.38	0.43
1:A:868:ASP:O	1:A:872:THR:HG23	2.18	0.43
1:A:575:LEU:CD1	1:A:589:ILE:HD11	2.49	0.43
1:A:51:ASP:O	1:A:52:GLU:HB2	2.19	0.43
1:B:463:VAL:HG23	1:B:494:LEU:HD22	2.01	0.43
1:B:106:LYS:HE2	6:B:1405:HOH:O	2.19	0.42
1:A:134:ARG:HG3	1:A:495:GLY:HA2	2.01	0.42
1:A:543:HIS:CG	1:A:544:PRO:HD2	2.54	0.42
1:A:97:PHE:HB3	1:A:164:PHE:CZ	2.55	0.42
1:A:115:ASN:HD21	1:A:164:PHE:HA	1.84	0.42
1:B:357:GLN:HB2	6:B:1155:HOH:O	2.18	0.42
1:A:178:HIS:CE1	1:A:208:THR:HA	2.55	0.42
1:B:426:HIS:HB2	1:B:567:ARG:NH1	2.33	0.42
1:B:254:LEU:HB2	6:B:1142:HOH:O	2.18	0.42
1:A:513:LEU:HG	1:A:543:HIS:CE1	2.54	0.42
1:A:24:LEU:HA	1:A:27:LEU:HD13	2.02	0.42
1:B:45:ALA:O	1:B:49:ASN:HB2	2.20	0.42
1:B:222:GLY:HA3	1:B:807:ASN:ND2	2.24	0.42
1:A:312:VAL:HG13	1:A:316:SER:HB2	2.02	0.42
1:A:875:LEU:CD1	6:A:1496:HOH:O	2.68	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:144:ASP:O	1:A:148:GLU:HG3	2.20	0.42
1:B:175:GLY:HA2	1:B:501:TYR:HB3	2.02	0.42
1:B:115:ASN:HD21	1:B:164:PHE:HA	1.85	0.42
1:A:465:PRO:HG2	6:A:1256:HOH:O	2.20	0.41
1:A:267:HIS:HE1	1:A:359:VAL:CG1	2.32	0.41
1:A:704:ARG:H	1:A:704:ARG:HG3	1.47	0.41
1:A:630:SER:HB2	1:B:192:GLN:HE22	1.85	0.41
1:B:140:GLN:HB2	6:B:1087:HOH:O	2.19	0.41
1:A:8:LEU:HD11	1:A:27:LEU:CD1	2.50	0.41
1:B:422:PHE:CZ	1:B:465:PRO:HD2	2.55	0.41
1:B:442:ASN:HB3	1:B:445:VAL:HG22	2.02	0.41
1:A:192:GLN:NE2	6:A:1070:HOH:O	2.53	0.41
1:A:770:ALA:O	1:A:797:ALA:HA	2.20	0.41
1:A:845:PRO:HD3	6:A:1326:HOH:O	2.20	0.41
1:A:541:ARG:HD2	1:A:541:ARG:HA	1.90	0.41
1:A:485:SER:HA	1:A:603:TYR:CZ	2.55	0.41
1:B:247:LEU:HD12	1:B:285:VAL:HG22	2.01	0.41
1:B:153:ARG:HD2	1:B:711:SER:O	2.21	0.41
1:B:106:LYS:CE	6:B:1405:HOH:O	2.69	0.41
1:B:151:ARG:O	1:B:155:GLU:HG3	2.21	0.41
1:B:19:LYS:HB3	1:B:196:TYR:CE2	2.56	0.41
1:B:450:GLY:HA3	1:B:487:VAL:CG1	2.23	0.41
1:A:97:PHE:HB3	1:A:164:PHE:CE2	2.56	0.40
1:A:181:ASN:HA	1:A:185:LEU:HB2	2.02	0.40
1:A:115:ASN:HD22	1:A:165:HIS:H	1.69	0.40
1:A:778:SER:OG	2:A:890:ACT:CH3	2.69	0.40
1:B:412:HIS:HB2	1:B:423:THR:HG22	2.03	0.40
1:A:115:ASN:ND2	1:A:164:PHE:HA	2.35	0.40
1:B:594:ASP:HB2	6:B:1425:HOH:O	2.21	0.40
1:B:105:LYS:HG2	1:B:106:LYS:HD3	2.03	0.40
1:B:759:TYR:CG	1:B:766:LEU:HD21	2.57	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	887/888 (100%)	857 (97%)	28 (3%)	2 (0%)	56	38
1	B	886/888 (100%)	857 (97%)	27 (3%)	2 (0%)	56	38
All	All	1773/1776 (100%)	1714 (97%)	55 (3%)	4 (0%)	56	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	436	SER
1	B	436	SER
1	B	177	ILE
1	A	177	ILE

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	752/751 (100%)	731 (97%)	21 (3%)	56	37
1	B	751/751 (100%)	719 (96%)	32 (4%)	40	19
All	All	1503/1502 (100%)	1450 (96%)	53 (4%)	48	26

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86	GLN
1	A	88	PHE
1	A	147	PHE
1	A	201	LEU
1	A	233	GLN
1	A	262	LEU
1	A	266	LYS
1	A	321	VAL
1	A	331	LYS
1	A	387	LYS
1	A	462	ASN
1	A	480	TYR
1	A	525	ASP

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Mol	Chain	Res	Type
1	A	541	ARG
1	A	592	THR
1	A	610	LYS
1	A	669	LEU
1	A	678	ASP
1	A	689	ASN
1	A	704	ARG
1	A	823	ASN
1	B	2	SER
1	B	19	LYS
1	B	27	LEU
1	B	86	GLN
1	B	88	PHE
1	B	135	ARG
1	B	147	PHE
1	B	201	LEU
1	B	262	LEU
1	B	266	LYS
1	B	287	GLN
1	B	321	VAL
1	B	331	LYS
1	B	354	ASP
1	B	375	ARG
1	B	387	LYS
1	B	390	GLU
1	B	413	LYS
1	B	462	ASN
1	B	480	TYR
1	B	519	GLU
1	B	522	THR
1	B	523	GLN
1	B	573	GLU
1	B	592	THR
1	B	601	ARG
1	B	608	MET
1	B	638	SER
1	B	669	LEU
1	B	841	GLU
1	B	850	GLN
1	B	889	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	25	ASN
1	A	115	ASN
1	A	192	GLN
1	A	233	GLN
1	A	322	GLN
1	A	351	GLN
1	A	397	GLN
1	A	408	HIS
1	A	462	ASN
1	A	493	GLN
1	A	523	GLN
1	A	545	ASN
1	A	549	ASN
1	A	602	GLN
1	A	807	ASN
1	A	823	ASN
1	A	832	GLN
1	B	16	GLN
1	B	25	ASN
1	B	115	ASN
1	B	192	GLN
1	B	233	GLN
1	B	322	GLN
1	B	351	GLN
1	B	357	GLN
1	B	397	GLN
1	B	408	HIS
1	B	462	ASN
1	B	493	GLN
1	B	545	ASN
1	B	549	ASN
1	B	597	GLN
1	B	602	GLN
1	B	698	ASN
1	B	807	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

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	SF4	A	1000	1,6	12,12,12	7.55	10 (83%)	0,24,24	0.00	-
4	GOL	A	1001	-	5,5,5	0.35	0	5,5,5	0.27	0
4	GOL	A	1002	-	5,5,5	0.31	0	5,5,5	0.35	0
2	ACT	A	890	-	1,3,3	1.40	0	0,3,3	0.00	-
3	SF4	B	1000	1,2,6	12,12,12	6.28	9 (75%)	0,24,24	0.00	-
5	FMT	B	1001	-	2,2,2	0.80	0	1,1,1	0.29	0
2	ACT	B	890	3	1,3,3	0.55	0	0,3,3	0.00	-
2	ACT	B	891	-	1,3,3	1.06	0	0,3,3	0.00	-

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
3	SF4	A	1000	1,6	-	0/0/48/48	0/0/5/5
4	GOL	A	1001	-	-	0/4/4/4	0/0/0/0
4	GOL	A	1002	-	-	0/4/4/4	0/0/0/0
2	ACT	A	890	-	-	0/0/0/0	0/0/0/0
3	SF4	B	1000	1,2,6	-	0/0/48/48	0/0/5/5
5	FMT	B	1001	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ACT	B	890	3	-	0/0/0/0	0/0/0/0
2	ACT	B	891	-	-	0/0/0/0	0/0/0/0

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1000	SF4	S3-FE1	-12.92	2.24	2.33
3	B	1000	SF4	S1-FE2	-11.96	2.25	2.33
3	A	1000	SF4	S1-FE2	-11.39	2.25	2.33
3	B	1000	SF4	S3-FE1	-10.15	2.26	2.33
3	A	1000	SF4	S3-FE4	-9.10	2.27	2.33
3	B	1000	SF4	S4-FE3	-8.62	2.27	2.33
3	A	1000	SF4	S4-FE2	-7.89	2.27	2.33
3	A	1000	SF4	S2-FE3	-7.52	2.28	2.33
3	A	1000	SF4	S2-FE4	-7.12	2.28	2.33
3	A	1000	SF4	S4-FE1	-6.92	2.28	2.33
3	B	1000	SF4	S4-FE2	-6.80	2.28	2.33
3	A	1000	SF4	S1-FE3	-6.49	2.28	2.33
3	A	1000	SF4	S4-FE3	-6.17	2.29	2.33
3	B	1000	SF4	S3-FE4	-5.98	2.29	2.33
3	B	1000	SF4	S1-FE3	-4.73	2.30	2.33
3	B	1000	SF4	S4-FE1	-4.24	2.30	2.33
3	B	1000	SF4	S2-FE3	-3.69	2.30	2.33
3	B	1000	SF4	S2-FE4	-3.11	2.31	2.33
3	A	1000	SF4	S2-FE1	-2.16	2.31	2.33

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	888/888 (100%)	-0.04	8 (0%) 81 81	17, 26, 40, 58	6 (0%)
1	B	888/888 (100%)	0.28	56 (6%) 19 18	17, 26, 40, 58	7 (0%)
All	All	1776/1776 (100%)	0.12	64 (3%) 41 38	17, 26, 40, 58	13 (0%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	581	GLY	10.2
1	B	410	ASN	7.1
1	B	524	GLY	6.7
1	A	579	ALA	6.2
1	B	540	GLY	5.3
1	B	412	HIS	5.1
1	B	580	LYS	5.0
1	B	425	ALA	4.7
1	B	407	GLU	4.5
1	B	423	THR	4.4
1	A	580	LYS	4.2
1	B	520	ALA	4.2
1	B	576	GLY	4.1
1	B	538	PHE	3.9
1	B	542	VAL	3.9
1	B	414	THR	3.9
1	B	546	THR	3.9
1	B	429	VAL	3.7
1	B	579	ALA	3.6
1	B	525	ASP	3.5
1	B	415	PHE	3.5
1	B	526	LEU	3.4
1	B	548	ALA	3.4
1	B	422	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	574	PRO	3.3
1	B	527	VAL	3.3
1	B	545	ASN	3.3
1	B	427	GLY	3.3
1	B	575	LEU	3.2
1	B	521	ILE	3.2
1	B	582	GLN	3.2
1	B	566	ILE	3.1
1	B	397	GLN	3.1
1	A	416	ILE	3.1
1	B	406	PRO	3.0
1	B	461	LEU	3.0
1	B	539	GLU	2.9
1	B	589	ILE	2.9
1	B	517	VAL	2.8
1	B	598	ALA	2.8
1	B	547	ARG	2.7
1	A	577	VAL	2.5
1	B	421	GLU	2.5
1	B	463	VAL	2.5
1	B	543	HIS	2.5
1	B	409	HIS	2.4
1	B	457	VAL	2.4
1	A	581	GLY	2.4
1	B	389	PHE	2.4
1	B	391	SER	2.3
1	B	416	ILE	2.3
1	A	53	PHE	2.3
1	A	578	ASN	2.3
1	B	411	ASP	2.2
1	B	577	VAL	2.2
1	B	398	GLY	2.2
1	B	393	LEU	2.1
1	B	605	ILE	2.1
1	B	560	TYR	2.1
1	B	394	GLY	2.1
1	B	405	ALA	2.1
1	B	466	TYR	2.1
1	B	572	LYS	2.0
1	A	16	GLN	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 ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	GOL	A	1001	6/6	0.34	14.85	48,48,51,54	0
2	ACT	A	890	4/4	0.15	5.84	26,29,31,31	0
2	ACT	B	891	4/4	0.15	3.02	22,22,24,26	0
4	GOL	A	1002	6/6	0.29	2.84	50,51,53,54	0
3	SF4	A	1000	8/8	0.12	1.99	21,22,25,26	0
5	FMT	B	1001	3/3	0.12	1.85	28,28,30,33	0
2	ACT	B	890	4/4	0.14	1.23	19,30,31,33	0
3	SF4	B	1000	8/8	0.10	-0.33	22,25,30,31	0

6.5 Other polymers ⓘ

There are no such residues in this entry.