Full wwPDB X-ray Structure Validation Report

Feb 1, 2016 – 01:57 PM GMT

PDB ID	:	3VMK
Title	:	3-isopropylmalate dehydrogenase from Shewanella benthica DB21 MT-2
Authors	:	Nagae, T.; Watanabe, N.
Deposited on	:	2011-12-13
Resolution	:	1.48 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

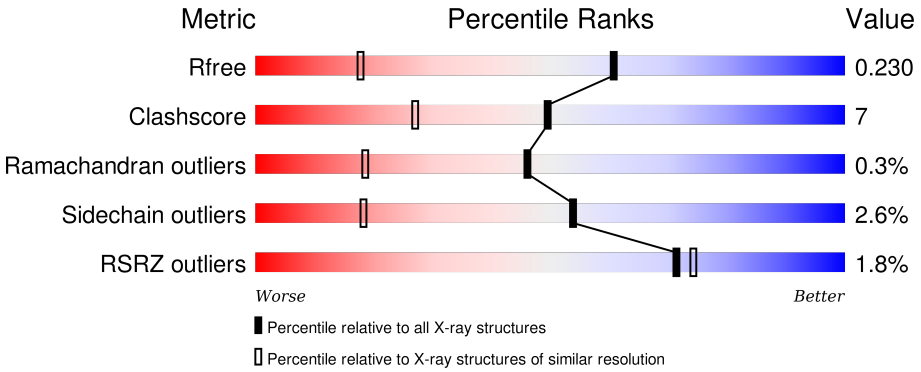
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.48 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}	91344	3129 (1.50-1.46)
Clashscore	102246	3380 (1.50-1.46)
Ramachandran outliers	100387	3310 (1.50-1.46)
Sidechain outliers	100360	3308 (1.50-1.46)
RSRZ outliers	91569	3133 (1.50-1.46)

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. 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 <=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	375	2% 80% 16%
1	B	375	2% 82% 13%

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	MG	B	402	-	-	-	X

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 6391 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.

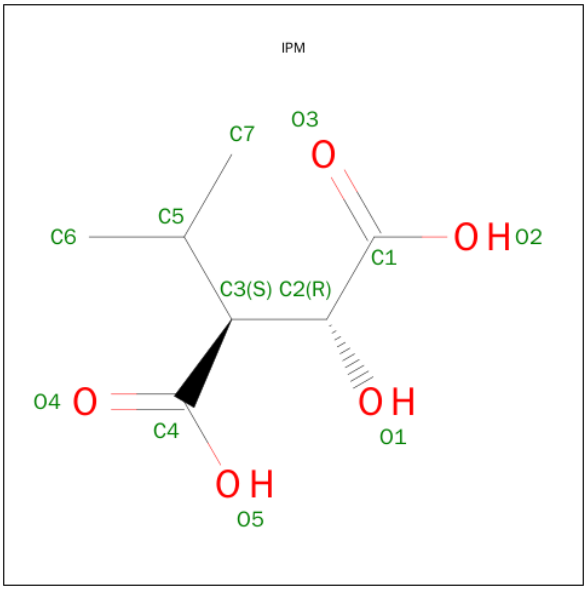
- Molecule 1 is a protein called 3-isopropylmalate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	S	0	7	0
			2832	1778	490	543	21			
1	B	368	Total	C	N	O	S	0	8	0
			2822	1772	487	541	22			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	EXPRESSION TAG	UNP D2YZL2
A	-9	ARG	-	EXPRESSION TAG	UNP D2YZL2
A	-8	GLY	-	EXPRESSION TAG	UNP D2YZL2
A	-7	SER	-	EXPRESSION TAG	UNP D2YZL2
A	-6	HIS	-	EXPRESSION TAG	UNP D2YZL2
A	-5	HIS	-	EXPRESSION TAG	UNP D2YZL2
A	-4	HIS	-	EXPRESSION TAG	UNP D2YZL2
A	-3	HIS	-	EXPRESSION TAG	UNP D2YZL2
A	-2	HIS	-	EXPRESSION TAG	UNP D2YZL2
A	-1	HIS	-	EXPRESSION TAG	UNP D2YZL2
A	0	GLY	-	EXPRESSION TAG	UNP D2YZL2
A	1	SER	-	EXPRESSION TAG	UNP D2YZL2
B	-10	MET	-	EXPRESSION TAG	UNP D2YZL2
B	-9	ARG	-	EXPRESSION TAG	UNP D2YZL2
B	-8	GLY	-	EXPRESSION TAG	UNP D2YZL2
B	-7	SER	-	EXPRESSION TAG	UNP D2YZL2
B	-6	HIS	-	EXPRESSION TAG	UNP D2YZL2
B	-5	HIS	-	EXPRESSION TAG	UNP D2YZL2
B	-4	HIS	-	EXPRESSION TAG	UNP D2YZL2
B	-3	HIS	-	EXPRESSION TAG	UNP D2YZL2
B	-2	HIS	-	EXPRESSION TAG	UNP D2YZL2
B	-1	HIS	-	EXPRESSION TAG	UNP D2YZL2
B	0	GLY	-	EXPRESSION TAG	UNP D2YZL2
B	1	SER	-	EXPRESSION TAG	UNP D2YZL2

- Molecule 2 is 3-ISOPROPYLMALIC ACID (three-letter code: IPM) (formula: $C_7H_{12}O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 12 7 5	0	0
2	B	1	Total C O 12 7 5	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Cl 1 1	0	0
4	A	1	Total Cl 1 1	0	0

- Molecule 5 is water.

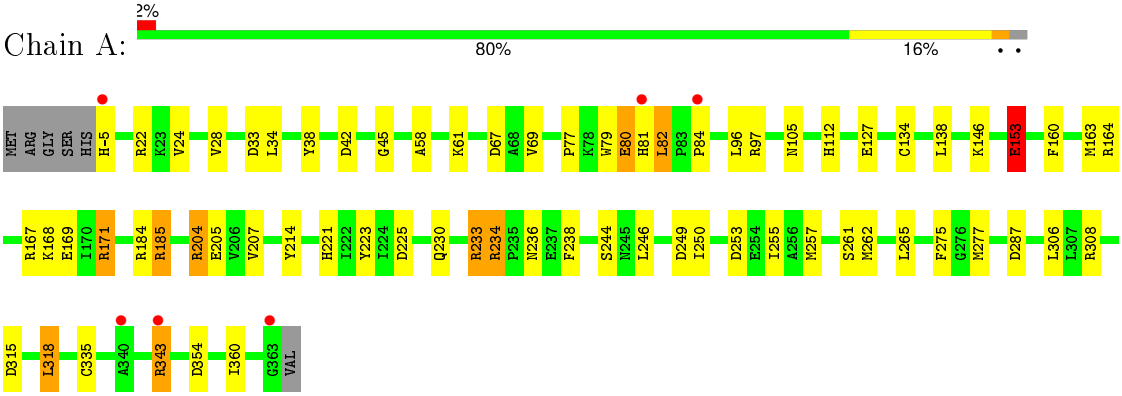
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	375	Total O 375 375	0	0
5	B	334	Total O 334 334	0	0

3 Residue-property plots (i)

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.

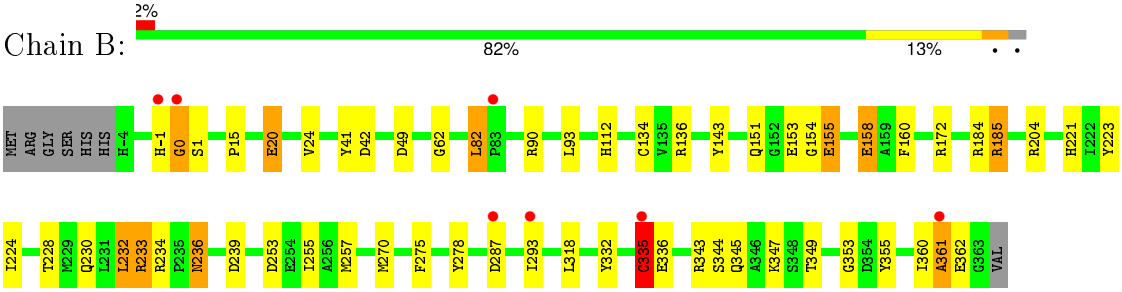
- Molecule 1: 3-isopropylmalate dehydrogenase

Chain A:



- Molecule 1: 3-isopropylmalate dehydrogenase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	51.97Å 59.21Å 119.78Å 90.00° 95.12° 90.00°	Depositor
Resolution (Å)	25.77 – 1.48 25.77 – 1.48	Depositor EDS
% Data completeness (in resolution range)	98.0 (25.77-1.48) 97.9 (25.77-1.48)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.41 (at 1.48Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.193 , 0.232 0.192 , 0.230	Depositor DCC
R_{free} test set	5928 reflections (5.29%)	DCC
Wilson B-factor (Å^2)	14.6	Xtriage
Anisotropy	0.484	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 118063 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	6391	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.64% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, CL, IPM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|>5$	RMSZ	#$\|Z\|>5$
1	A	1.41	9/2907 (0.3%)	1.38	31/3925 (0.8%)
1	B	1.40	12/2899 (0.4%)	1.28	22/3914 (0.6%)
All	All	1.41	21/5806 (0.4%)	1.33	53/7839 (0.7%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	38	TYR	CE1-CZ	9.46	1.50	1.38
1	B	154	GLY	N-CA	7.97	1.58	1.46
1	A	261	SER	CB-OG	7.84	1.52	1.42
1	A	223	TYR	CD1-CE1	7.38	1.50	1.39
1	B	335[A]	CYS	CB-SG	6.89	1.94	1.82
1	B	335[B]	CYS	CB-SG	6.89	1.94	1.82
1	B	223	TYR	CE2-CZ	-6.65	1.29	1.38
1	A	207	VAL	CB-CG2	6.63	1.66	1.52
1	B	0	GLY	N-CA	6.61	1.55	1.46
1	B	335[A]	CYS	CA-CB	6.38	1.68	1.53
1	B	335[B]	CYS	CA-CB	6.38	1.68	1.53
1	B	278	TYR	CG-CD1	5.97	1.47	1.39
1	B	155	GLU	C-O	5.91	1.34	1.23
1	B	158	GLU	CB-CG	-5.83	1.41	1.52
1	A	164	ARG	CZ-NH1	5.50	1.40	1.33
1	B	336	GLU	CD-OE1	5.50	1.31	1.25
1	A	335	CYS	CB-SG	-5.49	1.72	1.81
1	B	0	GLY	CA-C	5.42	1.60	1.51
1	A	214	TYR	CG-CD2	5.38	1.46	1.39
1	A	146	LYS	CE-NZ	5.10	1.61	1.49
1	A	45	GLY	N-CA	5.09	1.53	1.46

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	185	ARG	NE-CZ-NH2	-11.21	114.69	120.30
1	A	185	ARG	NE-CZ-NH1	9.89	125.25	120.30
1	A	204	ARG	NE-CZ-NH1	9.84	125.22	120.30
1	A	234	ARG	NE-CZ-NH1	9.53	125.06	120.30
1	A	22	ARG	NE-CZ-NH2	9.19	124.90	120.30
1	A	164	ARG	NE-CZ-NH2	-9.06	115.77	120.30
1	A	42	ASP	CB-CG-OD2	-9.00	110.20	118.30
1	A	22	ARG	NE-CZ-NH1	-8.39	116.10	120.30
1	A	308	ARG	NE-CZ-NH2	-8.21	116.19	120.30
1	A	308	ARG	NE-CZ-NH1	7.82	124.21	120.30
1	A	354	ASP	CB-CG-OD1	7.63	125.16	118.30
1	A	318	LEU	CB-CG-CD1	7.22	123.28	111.00
1	B	136	ARG	NE-CZ-NH2	7.14	123.87	120.30
1	A	42	ASP	CB-CG-OD1	7.08	124.67	118.30
1	A	167	ARG	NE-CZ-NH2	-7.02	116.79	120.30
1	A	253	ASP	CB-CG-OD2	-6.73	112.24	118.30
1	A	249	ASP	CB-CG-OD2	-6.55	112.40	118.30
1	B	136	ARG	NE-CZ-NH1	6.46	123.53	120.30
1	A	184	ARG	NE-CZ-NH2	-6.37	117.11	120.30
1	B	143	TYR	CB-CG-CD2	-6.35	117.19	121.00
1	B	0	GLY	N-CA-C	6.32	128.90	113.10
1	B	49	ASP	CB-CG-OD2	-6.32	112.62	118.30
1	B	275	PHE	CB-CG-CD1	6.28	125.20	120.80
1	B	184	ARG	NE-CZ-NH2	-6.27	117.16	120.30
1	B	136	ARG	NH1-CZ-NH2	-6.20	112.59	119.40
1	B	42	ASP	CB-CG-OD1	6.12	123.81	118.30
1	A	233	ARG	NE-CZ-NH1	-6.09	117.25	120.30
1	A	205	GLU	OE1-CD-OE2	5.83	130.29	123.30
1	B	318	LEU	CA-CB-CG	5.79	128.61	115.30
1	A	38	TYR	CG-CD1-CE1	5.76	125.91	121.30
1	A	315	ASP	CB-CG-OD2	-5.71	113.16	118.30
1	B	239	ASP	CB-CG-OD2	-5.70	113.17	118.30
1	A	275	PHE	CB-CG-CD2	5.67	124.77	120.80
1	B	270	MET	CA-CB-CG	5.63	122.88	113.30
1	A	153	GLU	N-CA-CB	5.59	120.67	110.60
1	A	225	ASP	CB-CG-OD1	-5.55	113.30	118.30
1	B	93	LEU	CB-CG-CD1	-5.55	101.56	111.00
1	A	171	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	A	169	GLU	OE1-CD-OE2	-5.46	116.75	123.30
1	B	232	LEU	CA-CB-CG	-5.45	102.76	115.30
1	B	184	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	B	0	GLY	C-N-CA	5.37	135.13	121.70
1	A	34	LEU	CB-CG-CD2	5.35	120.09	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	257	MET	CG-SD-CE	5.34	108.74	100.20
1	A	225	ASP	CB-CG-OD2	5.27	123.04	118.30
1	B	82	LEU	CA-CB-CG	5.25	127.37	115.30
1	B	143	TYR	CB-CG-CD1	5.20	124.12	121.00
1	A	265	LEU	CB-CG-CD1	-5.12	102.30	111.00
1	B	20	GLU	OE1-CD-OE2	5.12	129.44	123.30
1	A	246	LEU	CB-CG-CD2	-5.11	102.31	111.00
1	B	172	ARG	NE-CZ-NH1	-5.11	117.74	120.30
1	B	185	ARG	NE-CZ-NH2	5.03	122.81	120.30
1	A	160	PHE	CB-CG-CD2	-5.03	117.28	120.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2832	0	2814	49	0
1	B	2822	0	2811	40	0
2	A	12	0	9	0	0
2	B	12	0	9	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	375	0	0	5	0
5	B	334	0	0	3	0
All	All	6391	0	5643	80	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 (80) 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:185:ARG:HH11	1:B:185:ARG:HG3	1.09	1.10
1:B:185:ARG:NH1	1:B:185:ARG:HG3	1.80	0.94
1:A:127:GLU:HG2	5:A:766:HOH:O	1.75	0.86
1:B:361:ALA:O	1:B:362:GLU:HG3	1.83	0.78
1:A:171:ARG:NH1	1:B:155:GLU:CD	2.38	0.77
5:A:525:HOH:O	1:B:158:GLU:HG3	1.86	0.76
1:B:-1:HIS:O	1:B:1:SER:N	2.18	0.75
1:A:112:HIS:HD2	5:A:622:HOH:O	1.68	0.75
1:A:171:ARG:NH1	1:B:155:GLU:OE1	2.21	0.73
1:A:69:VAL:HG11	1:A:277[B]:MET:HE2	1.70	0.73
1:A:77:PRO:HA	1:A:80:GLU:CG	2.19	0.73
1:B:230:GLN:HE22	1:B:233:ARG:HH11	1.35	0.71
1:A:84:PRO:HB3	5:B:695:HOH:O	1.91	0.70
1:A:69:VAL:HG11	1:A:277[B]:MET:CE	2.23	0.69
1:B:112:HIS:HD2	5:B:619:HOH:O	1.76	0.69
1:B:344:SER:OG	1:B:345:GLN:OE1	2.10	0.67
1:B:151:GLN:OE1	1:B:158:GLU:OE1	2.12	0.67
1:B:185:ARG:CG	1:B:185:ARG:NH1	2.57	0.66
1:A:168:LYS:NZ	1:A:168:LYS:HB2	2.11	0.66
1:A:171:ARG:HH12	1:B:155:GLU:CD	1.99	0.66
1:A:343:ARG:HA	1:A:343:ARG:NE	2.12	0.65
1:A:-5:HIS:HD2	1:A:67:ASP:OD1	1.78	0.65
1:A:24:VAL:HG13	1:A:360:ILE:HD12	1.79	0.64
1:A:168:LYS:NZ	1:B:155:GLU:HB3	2.12	0.64
1:A:230:GLN:HE21	1:A:234:ARG:HE	1.46	0.64
1:A:81:HIS:CE1	1:A:82:LEU:HD13	2.33	0.63
1:A:153:GLU:OE2	1:A:153:GLU:CA	2.47	0.62
1:B:230:GLN:HE21	1:B:234:ARG:HE	1.48	0.60
1:A:77:PRO:HA	1:A:80:GLU:HG2	1.82	0.59
1:A:343:ARG:HE	1:A:343:ARG:HA	1.67	0.59
1:B:332:TYR:CD2	1:B:347:LYS:HG3	2.37	0.58
1:A:230:GLN:HE22	1:A:233:ARG:HH11	1.50	0.58
1:A:28:VAL:HG23	1:A:360:ILE:HD13	1.87	0.56
1:A:96:LEU:HD13	1:A:277[B]:MET:HE1	1.88	0.54
1:A:168:LYS:HZ2	1:A:168:LYS:HB2	1.73	0.53
1:B:151:GLN:HG2	1:B:158:GLU:HG2	1.91	0.52
1:A:69:VAL:CG1	1:A:277[B]:MET:HE2	2.40	0.51
1:B:230:GLN:NE2	1:B:234:ARG:HE	2.09	0.51
1:A:134[B]:CYS:SG	1:A:255:ILE:HD12	2.51	0.50
1:B:185:ARG:CG	1:B:185:ARG:HH11	1.95	0.49
1:A:-5:HIS:CD2	1:A:67:ASP:OD1	2.64	0.49
1:B:20:GLU:OE2	1:B:353:GLY:HA3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:ALA:HA	1:A:61:LYS:HE3	1.95	0.48
1:A:257:MET:CE	1:A:262[A]:MET:CE	2.92	0.48
1:B:293:ILE:HG23	1:B:335[A]:CYS:HB3	1.96	0.48
1:B:24:VAL:HG13	1:B:360:ILE:CD1	2.43	0.47
1:A:230:GLN:NE2	1:A:233:ARG:HH11	2.12	0.47
1:B:134[B]:CYS:SG	1:B:255:ILE:HD12	2.54	0.47
1:A:171:ARG:NH1	1:B:155:GLU:OE2	2.47	0.47
1:B:335[B]:CYS:SG	1:B:343:ARG:HD2	2.55	0.47
1:A:97:ARG:HD3	1:A:138:LEU:HG	1.97	0.47
1:A:79:TRP:O	1:A:82:LEU:HB2	2.16	0.46
1:A:28:VAL:CG2	1:A:360:ILE:HD13	2.46	0.46
1:A:250:ILE:HG23	1:B:228:THR:HG21	1.97	0.46
1:B:343:ARG:HG2	1:B:343:ARG:HH11	1.81	0.45
1:A:171:ARG:NH2	1:B:155:GLU:OE2	2.49	0.45
1:B:41:TYR:CZ	1:B:62:GLY:HA3	2.52	0.45
1:B:224:ILE:O	1:B:228:THR:HG23	2.16	0.45
1:B:236:ASN:ND2	5:B:565:HOH:O	2.50	0.44
1:A:168:LYS:HG3	5:A:549:HOH:O	2.17	0.44
1:A:24:VAL:HG13	1:A:360:ILE:CD1	2.46	0.43
1:B:355:TYR:C	1:B:355:TYR:CD2	2.91	0.43
1:A:287:ASP:OD2	1:A:287:ASP:N	2.51	0.43
1:B:287:ASP:N	1:B:287:ASP:OD1	2.50	0.43
1:A:257:MET:CE	1:A:262[A]:MET:HE1	2.48	0.43
1:B:335[B]:CYS:SG	1:B:343:ARG:CD	3.07	0.43
1:A:257:MET:HE1	1:A:262[A]:MET:CE	2.50	0.42
1:A:230:GLN:NE2	1:A:234:ARG:HE	2.16	0.42
1:B:20:GLU:HG3	1:B:349:THR:HG22	2.03	0.41
1:B:232:LEU:HD23	1:B:232:LEU:HA	1.83	0.41
1:A:230:GLN:HE22	1:A:233:ARG:HE	1.68	0.41
1:B:204:ARG:HG3	1:B:221:HIS:CE1	2.56	0.41
1:A:244:SER:OG	5:A:747:HOH:O	1.94	0.41
1:A:306:LEU:HD23	1:A:306:LEU:C	2.42	0.40
1:B:232:LEU:H	1:B:232:LEU:HG	1.82	0.40
1:A:163:MET:O	1:B:160:PHE:HA	2.20	0.40
1:A:168:LYS:HZ3	1:B:155:GLU:HB3	1.82	0.40
1:A:105:ASN:HB2	1:A:138:LEU:HD22	2.01	0.40
1:A:204:ARG:HG3	1:A:221:HIS:CE1	2.56	0.40
1:A:185:ARG:NH2	1:A:238:PHE:O	2.50	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	374/375 (100%)	360 (96%)	14 (4%)	0	100 100
1	B	374/375 (100%)	357 (96%)	15 (4%)	2 (0%)	34 10
All	All	748/750 (100%)	717 (96%)	29 (4%)	2 (0%)	46 18

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	0	GLY
1	B	361	ALA

5.3.2 Protein sidechains (i)

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	300/298 (101%)	293 (98%)	7 (2%)	58 23
1	B	300/298 (101%)	291 (97%)	9 (3%)	48 14
All	All	600/596 (101%)	584 (97%)	16 (3%)	54 17

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	ASP
1	A	80	GLU
1	A	82	LEU
1	A	153	GLU

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Mol	Chain	Res	Type
1	A	236	ASN
1	A	318	LEU
1	A	343	ARG
1	B	15	PRO
1	B	82	LEU
1	B	90	ARG
1	B	153	GLU
1	B	233	ARG
1	B	236	ASN
1	B	253	ASP
1	B	335[A]	CYS
1	B	335[B]	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-5	HIS
1	A	51	HIS
1	A	81	HIS
1	A	112	HIS
1	A	230	GLN
1	B	4	GLN
1	B	105	ASN
1	B	112	HIS
1	B	151	GLN
1	B	230	GLN
1	B	236	ASN
1	B	273	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

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	IPM	A	401	3	4,11,11	0.76	0	6,15,15	2.28	1 (16%)
2	IPM	B	401	3	4,11,11	0.95	0	6,15,15	1.44	2 (33%)

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
2	IPM	A	401	3	-	0/8/16/16	0/0/0/0
2	IPM	B	401	3	-	0/8/16/16	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	401	IPM	C1-C2-C3	-5.49	107.16	112.34
2	B	401	IPM	C7-C5-C6	-2.31	103.88	110.67
2	B	401	IPM	C1-C2-C3	-2.01	110.45	112.34

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

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, 95^{th} 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	369/375 (98%)	-0.31	6 (1%)	74	78	7, 16, 33, 51	0
1	B	368/375 (98%)	-0.23	7 (1%)	70	72	10, 19, 33, 46	0
All	All	737/750 (98%)	-0.27	13 (1%)	71	74	7, 17, 33, 51	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	84	PRO	4.9
1	B	293	ILE	4.0
1	A	-5	HIS	3.5
1	B	0	GLY	3.4
1	A	363	GLY	3.2
1	A	81	HIS	2.9
1	B	335[A]	CYS	2.8
1	B	361	ALA	2.8
1	B	-1	HIS	2.5
1	A	343	ARG	2.4
1	B	287	ASP	2.3
1	A	340	ALA	2.2
1	B	83	PRO	2.1

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, 95^{th} 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	LLDF	B-factors(Å^2)	Q<0.9
3	MG	B	402	1/1	0.99	0.07	2.44	12,12,12,12	0
4	CL	A	403	1/1	1.00	0.07	0.53	10,10,10,10	0
2	IPM	B	401	12/12	0.97	0.06	-0.40	12,13,14,15	0
4	CL	B	403	1/1	0.99	0.05	-0.92	11,11,11,11	0
2	IPM	A	401	12/12	0.98	0.06	-1.11	9,10,11,12	0
3	MG	A	402	1/1	0.97	0.06	-1.90	7,7,7,7	0

6.5 Other polymers

There are no such residues in this entry.