Full wwPDB X-ray Structure Validation Report (i)

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PDB ID	:	1FI4
Title	:	THE X-RAY CRYSTAL STRUCTURE OF MEVALONATE 5-DIPHOSPHATE DECARBOXYLASE AT 2.3 ANGSTROM RESOLUTION.
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Deposited on	:	2000-08-03
Resolution	:	2.27 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:

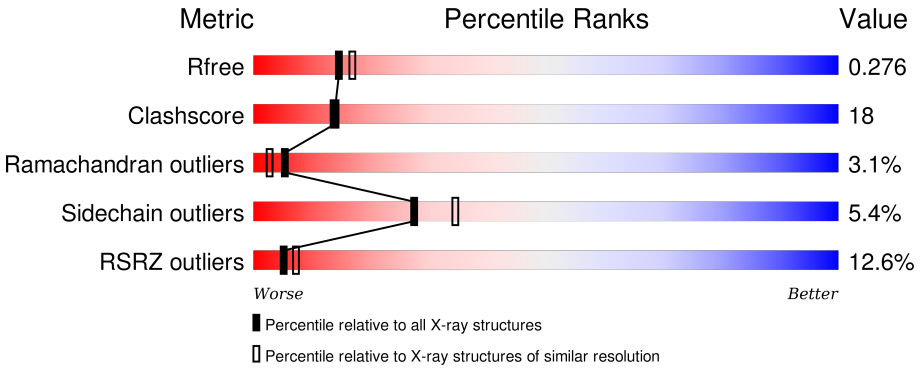
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.27 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	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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	416	12% 65% 24% • • 6%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3226 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 MEVALONATE 5-DIPHOSPHATE DECARBOXYLASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	Se	0	0	0
			3062	1935	516	597	5	9			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	see remark 999	UNP P32377
A	-18	GLY	-	see remark 999	UNP P32377
A	-17	SER	-	see remark 999	UNP P32377
A	-16	SER	-	see remark 999	UNP P32377
A	-15	HIS	-	see remark 999	UNP P32377
A	-14	HIS	-	see remark 999	UNP P32377
A	-13	HIS	-	see remark 999	UNP P32377
A	-12	HIS	-	see remark 999	UNP P32377
A	-11	HIS	-	see remark 999	UNP P32377
A	-10	HIS	-	see remark 999	UNP P32377
A	-9	SER	-	see remark 999	UNP P32377
A	-8	SER	-	see remark 999	UNP P32377
A	-7	GLY	-	see remark 999	UNP P32377
A	-6	LEU	-	see remark 999	UNP P32377
A	-5	VAL	-	see remark 999	UNP P32377
A	-4	PRO	-	see remark 999	UNP P32377
A	-3	ARG	-	see remark 999	UNP P32377
A	-2	GLY	-	see remark 999	UNP P32377
A	-1	SER	-	see remark 999	UNP P32377
A	0	HIS	-	see remark 999	UNP P32377
A	1	MSE	MET	MODIFIED RESIDUE	UNP P32377
A	89	MSE	MET	MODIFIED RESIDUE	UNP P32377
A	169	MSE	MET	MODIFIED RESIDUE	UNP P32377
A	179	MSE	MET	MODIFIED RESIDUE	UNP P32377
A	192	MSE	MET	MODIFIED RESIDUE	UNP P32377
A	212	MSE	MET	MODIFIED RESIDUE	UNP P32377
A	237	MSE	MET	MODIFIED RESIDUE	UNP P32377

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Chain	Residue	Modelled	Actual	Comment	Reference
A	254	MSE	MET	MODIFIED RESIDUE	UNP P32377
A	255	MSE	MET	MODIFIED RESIDUE	UNP P32377
A	274	MSE	MET	MODIFIED RESIDUE	UNP P32377

- Molecule 2 is water.

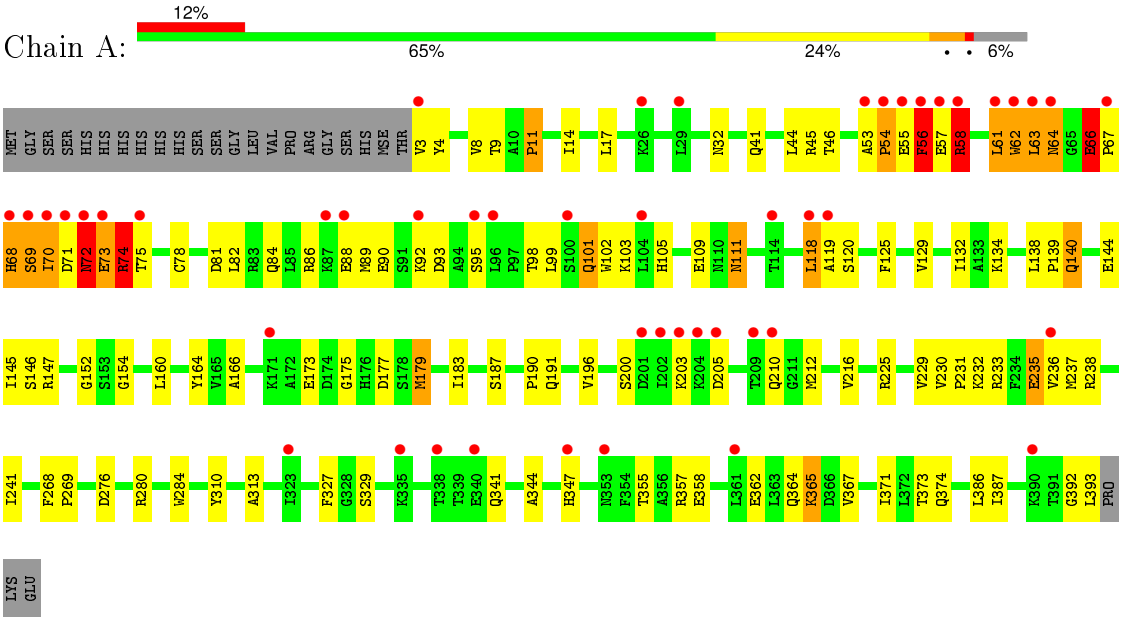
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	164	Total	O	0	0
			164	164		

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: MEVALONATE 5-DIPHOSPHATE DECARBOXYLASE

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	78.80Å 126.40Å 47.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.27 18.00 – 2.28	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.00-2.27) 99.9 (18.00-2.28)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.63 (at 2.28Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.239 , 0.268 0.248 , 0.276	Depositor DCC
R_{free} test set	2187 reflections (9.86%)	DCC
Wilson B-factor (Å^2)	36.0	Xtriage
Anisotropy	0.668	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 39.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	1 of 41454 reflections (0.002%)	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	3226	wwPDB-VP
Average B, all atoms (Å^2)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.54% 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

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	0.42	0/3117	0.76	10/4210 (0.2%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	71	ASP	N-CA-C	-7.54	90.63	111.00
1	A	63	LEU	N-CA-C	-7.24	91.46	111.00
1	A	72	ASN	CA-C-N	-7.08	101.61	117.20
1	A	74	ARG	N-CA-C	-6.81	92.61	111.00
1	A	56	PHE	N-CA-C	-6.71	92.89	111.00
1	A	72	ASN	CB-CA-C	-6.56	97.28	110.40
1	A	66	GLU	N-CA-C	6.54	128.66	111.00
1	A	57	GLU	N-CA-C	6.17	127.66	111.00
1	A	55	GLU	N-CA-C	5.54	125.95	111.00
1	A	72	ASN	C-N-CA	5.53	135.53	121.70

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	3062	0	3009	111	0
2	A	164	0	0	4	0
All	All	3226	0	3009	111	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 18.

All (111) 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:138:LEU:HB3	1:A:140:GLN:HE21	1.20	1.05
1:A:62:TRP:HD1	1:A:64:ASN:HA	1.20	1.01
1:A:62:TRP:CD1	1:A:64:ASN:HA	1.96	0.99
1:A:329:SER:H	1:A:364:GLN:NE2	1.68	0.91
1:A:70:ILE:C	1:A:72:ASN:H	1.70	0.90
1:A:75:THR:OG1	1:A:125:PHE:HE1	1.60	0.83
1:A:58:ARG:H	1:A:103:LYS:HG3	1.42	0.83
1:A:138:LEU:HB3	1:A:140:GLN:NE2	1.91	0.83
1:A:70:ILE:N	1:A:73:GLU:HB2	1.99	0.78
1:A:173:GLU:CD	1:A:173:GLU:H	1.87	0.77
1:A:362:GLU:O	1:A:365:LYS:HG2	1.85	0.77
1:A:125:PHE:CE2	1:A:152:GLY:HA3	2.20	0.76
1:A:75:THR:OG1	1:A:125:PHE:CE1	2.37	0.76
1:A:387:ILE:HG23	1:A:392:GLY:HA2	1.69	0.75
1:A:61:LEU:HB2	1:A:68:HIS:HE1	1.53	0.73
1:A:84:GLN:O	1:A:88:GLU:HG2	1.88	0.73
1:A:45:ARG:H	1:A:111:ASN:HD21	1.35	0.73
1:A:103:LYS:HA	1:A:103:LYS:HE2	1.72	0.72
1:A:11:PRO:HB3	1:A:41:GLN:HG3	1.74	0.70
1:A:66:GLU:CD	1:A:66:GLU:H	1.96	0.70
1:A:329:SER:H	1:A:364:GLN:HE22	1.40	0.69
1:A:140:GLN:HG2	1:A:145:ILE:HD11	1.75	0.69
1:A:3:VAL:HA	1:A:53:ALA:HB2	1.75	0.68
1:A:17:LEU:HD13	1:A:233:ARG:HG2	1.77	0.66
1:A:69:SER:OG	1:A:73:GLU:HG3	1.96	0.65
1:A:82:LEU:HD13	1:A:132:ILE:HG21	1.80	0.64
1:A:61:LEU:HB2	1:A:68:HIS:CE1	2.32	0.64
1:A:327:PHE:CE2	1:A:367:VAL:HG21	2.34	0.63
1:A:62:TRP:HD1	1:A:64:ASN:CA	2.05	0.63
1:A:139:PRO:HD2	1:A:140:GLN:NE2	2.15	0.61
1:A:70:ILE:H	1:A:73:GLU:HB2	1.64	0.61
1:A:56:PHE:O	1:A:56:PHE:HD2	1.83	0.59
1:A:8:VAL:HG12	1:A:9:THR:N	2.17	0.59
1:A:357:ARG:HG2	1:A:358:GLU:N	2.18	0.59
1:A:4:TYR:CD2	1:A:99:LEU:HD21	2.37	0.59
1:A:101:GLN:HE21	1:A:101:GLN:N	2.01	0.59
1:A:357:ARG:HG2	1:A:358:GLU:H	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:PHE:O	1:A:129:VAL:HG23	2.03	0.58
1:A:177:ASP:HA	1:A:179:MSE:HE2	1.84	0.58
1:A:63:LEU:O	1:A:64:ASN:HB2	2.04	0.58
1:A:58:ARG:N	1:A:103:LYS:HG3	2.16	0.57
1:A:140:GLN:CD	1:A:140:GLN:H	2.06	0.57
1:A:111:ASN:H	1:A:111:ASN:HD22	1.53	0.57
1:A:69:SER:O	1:A:70:ILE:HB	2.04	0.57
1:A:70:ILE:CA	1:A:73:GLU:HB2	2.34	0.57
1:A:212:MSE:O	1:A:216:VAL:HG23	2.06	0.55
1:A:14:ILE:HB	1:A:310:TYR:OH	2.07	0.55
1:A:118:LEU:O	1:A:120:SER:N	2.40	0.55
1:A:387:ILE:CG2	1:A:392:GLY:HA2	2.36	0.55
1:A:99:LEU:HA	1:A:102:TRP:CE3	2.43	0.54
1:A:41:GLN:HA	1:A:44:LEU:O	2.08	0.54
1:A:82:LEU:HD13	1:A:132:ILE:CG2	2.39	0.53
1:A:173:GLU:CD	1:A:173:GLU:N	2.61	0.53
1:A:364:GLN:HB2	2:A:1028:HOH:O	2.09	0.53
1:A:144:GLU:HG3	1:A:147:ARG:NH2	2.23	0.52
1:A:175:GLY:HA2	2:A:1073:HOH:O	2.11	0.51
1:A:139:PRO:HD2	1:A:140:GLN:HE22	1.74	0.50
1:A:86:ARG:CZ	1:A:99:LEU:HD12	2.42	0.50
1:A:164:TYR:N	1:A:164:TYR:CD1	2.80	0.50
1:A:90:GLU:C	1:A:92:LYS:H	2.14	0.50
1:A:347:HIS:HB2	2:A:1128:HOH:O	2.11	0.49
1:A:191:GLN:HG2	1:A:313:ALA:H	1.77	0.48
1:A:93:ASP:OD2	1:A:95:SER:HB3	2.13	0.48
1:A:70:ILE:HA	1:A:73:GLU:HB2	1.94	0.48
1:A:111:ASN:ND2	1:A:111:ASN:H	2.11	0.48
1:A:78:CYS:O	1:A:81:ASP:HB2	2.14	0.47
1:A:86:ARG:O	1:A:89:MSE:HB2	2.15	0.47
1:A:9:THR:HA	1:A:46:THR:O	2.15	0.46
1:A:134:LYS:HE2	2:A:1153:HOH:O	2.14	0.46
1:A:72:ASN:ND2	1:A:75:THR:HB	2.30	0.46
1:A:11:PRO:HB3	1:A:41:GLN:CG	2.44	0.46
1:A:225:ARG:HA	1:A:229:VAL:HB	1.98	0.46
1:A:69:SER:HA	1:A:73:GLU:HG2	1.97	0.46
1:A:166:ALA:HB2	1:A:183:ILE:HD11	1.98	0.45
1:A:45:ARG:H	1:A:111:ASN:ND2	2.09	0.45
1:A:355:THR:O	1:A:355:THR:HG22	2.16	0.45
1:A:8:VAL:CG1	1:A:9:THR:N	2.79	0.45
1:A:146:SER:HA	1:A:160:LEU:HD12	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:SER:HA	1:A:190:PRO:HG3	1.99	0.45
1:A:69:SER:HA	1:A:73:GLU:CG	2.47	0.45
1:A:61:LEU:HG	1:A:62:TRP:H	1.82	0.44
1:A:276:ASP:O	1:A:280:ARG:HG3	2.18	0.44
1:A:70:ILE:CG2	1:A:70:ILE:O	2.65	0.44
1:A:365:LYS:NZ	1:A:365:LYS:HB3	2.32	0.44
1:A:373:THR:OG1	1:A:374:GLN:N	2.51	0.44
1:A:63:LEU:O	1:A:63:LEU:HD13	2.19	0.43
1:A:329:SER:H	1:A:364:GLN:HE21	1.58	0.43
1:A:177:ASP:CA	1:A:179:MSE:HE2	2.49	0.43
1:A:140:GLN:CG	1:A:145:ILE:HD11	2.47	0.43
1:A:70:ILE:HD13	1:A:70:ILE:HA	1.91	0.43
1:A:75:THR:O	1:A:78:CYS:N	2.52	0.43
1:A:268:PHE:HA	1:A:269:PRO:C	2.40	0.43
1:A:32:ASN:HB2	1:A:230:VAL:HG11	2.00	0.43
1:A:196:VAL:HB	1:A:371:ILE:HB	2.00	0.43
1:A:203:LYS:HE3	1:A:205:ASP:OD1	2.18	0.43
1:A:284:TRP:CZ2	1:A:357:ARG:HD2	2.55	0.42
1:A:89:MSE:O	1:A:92:LYS:HB2	2.19	0.42
1:A:232:LYS:O	1:A:236:VAL:HG23	2.20	0.42
1:A:237:MSE:O	1:A:241:ILE:HG13	2.19	0.42
1:A:393:LEU:CD2	1:A:393:LEU:O	2.68	0.42
1:A:70:ILE:H	1:A:73:GLU:CB	2.31	0.42
1:A:66:GLU:HA	1:A:67:PRO:HD3	1.46	0.42
1:A:109:GLU:HG2	1:A:386:LEU:HD11	2.01	0.42
1:A:9:THR:O	1:A:41:GLN:NE2	2.53	0.41
1:A:341:GLN:O	1:A:344:ALA:HB3	2.20	0.41
1:A:62:TRP:CD1	1:A:64:ASN:CA	2.86	0.41
1:A:68:HIS:O	1:A:69:SER:HB2	2.20	0.41
1:A:191:GLN:CG	1:A:313:ALA:HB3	2.51	0.41
1:A:230:VAL:HB	1:A:231:PRO:HD3	2.01	0.41
1:A:235:GLU:OE2	1:A:238:ARG:NH2	2.54	0.40
1:A:70:ILE:HG22	1:A:70:ILE:O	2.20	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	389/416 (94%)	348 (90%)	29 (8%)	12 (3%)	5 3

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	64	ASN
1	A	70	ILE
1	A	72	ASN
1	A	119	ALA
1	A	58	ARG
1	A	69	SER
1	A	74	ARG
1	A	118	LEU
1	A	200	SER
1	A	54	PRO
1	A	66	GLU
1	A	154	GLY

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	336/348 (97%)	318 (95%)	18 (5%)	27 34

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	PRO
1	A	54	PRO
1	A	56	PHE
1	A	58	ARG
1	A	61	LEU
1	A	62	TRP
1	A	68	HIS
1	A	73	GLU
1	A	74	ARG
1	A	98	THR
1	A	101	GLN
1	A	105	HIS
1	A	111	ASN
1	A	140	GLN
1	A	179	MSE
1	A	210	GLN
1	A	235	GLU
1	A	365	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	41	GLN
1	A	68	HIS
1	A	76	GLN
1	A	101	GLN
1	A	111	ASN
1	A	140	GLN
1	A	353	ASN
1	A	364	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

There are no ligands in this entry.

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, 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	382/416 (91%)	0.66	48 (12%) 5 7	22, 41, 70, 75	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	72	ASN	8.0
1	A	202	ILE	7.1
1	A	58	ARG	6.8
1	A	55	GLU	6.4
1	A	204	LYS	5.6
1	A	64	ASN	5.6
1	A	54	PRO	4.9
1	A	57	GLU	4.7
1	A	201	ASP	4.3
1	A	69	SER	4.1
1	A	210	GLN	4.1
1	A	68	HIS	3.7
1	A	119	ALA	3.6
1	A	114	THR	3.6
1	A	70	ILE	3.5
1	A	56	PHE	3.4
1	A	353	ASN	3.4
1	A	95	SER	3.4
1	A	118	LEU	3.3
1	A	73	GLU	3.3
1	A	63	LEU	3.2
1	A	87	LYS	3.2
1	A	96	LEU	3.2
1	A	62	TRP	3.1
1	A	236	VAL	3.1
1	A	390	LYS	3.1
1	A	53	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	205	ASP	2.9
1	A	104	LEU	2.9
1	A	3	VAL	2.9
1	A	92	LYS	2.9
1	A	171	LYS	2.8
1	A	209	THR	2.8
1	A	67	PRO	2.8
1	A	88	GLU	2.7
1	A	71	ASP	2.7
1	A	100	SER	2.7
1	A	61	LEU	2.7
1	A	323	ILE	2.6
1	A	347	HIS	2.5
1	A	203	LYS	2.5
1	A	340	GLU	2.5
1	A	335	LYS	2.4
1	A	26	LYS	2.4
1	A	338	THR	2.1
1	A	75	THR	2.1
1	A	29	LEU	2.0
1	A	361	LEU	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 carbohydrates in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.