



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 1, 2023 – 09:13 AM EDT

PDB ID : 3T5B
Title : Crystal structure of N-terminal domain of FACL13 from Mycobacterium tuberculosis
Authors : Goyal, A.; Sankaranarayanan, R.
Deposited on : 2011-07-27
Resolution : 2.35 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

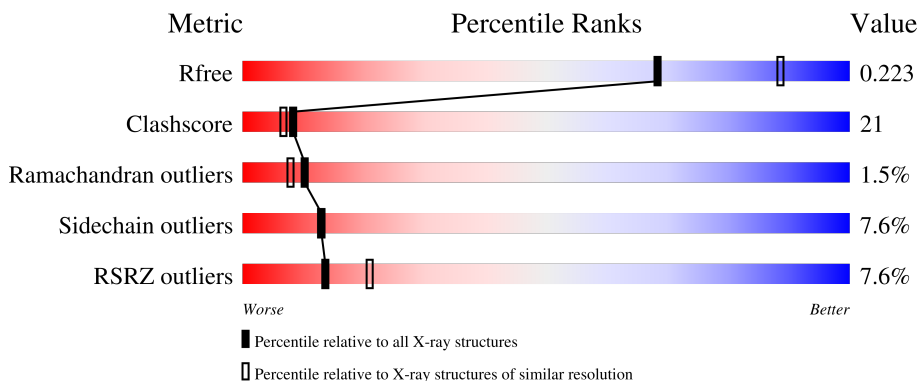
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	396	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3294 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 PROBABLE CHAIN-FATTY-ACID-CoA LIGASE FADD13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	396	3007	1904	515	570	18	0	0	0

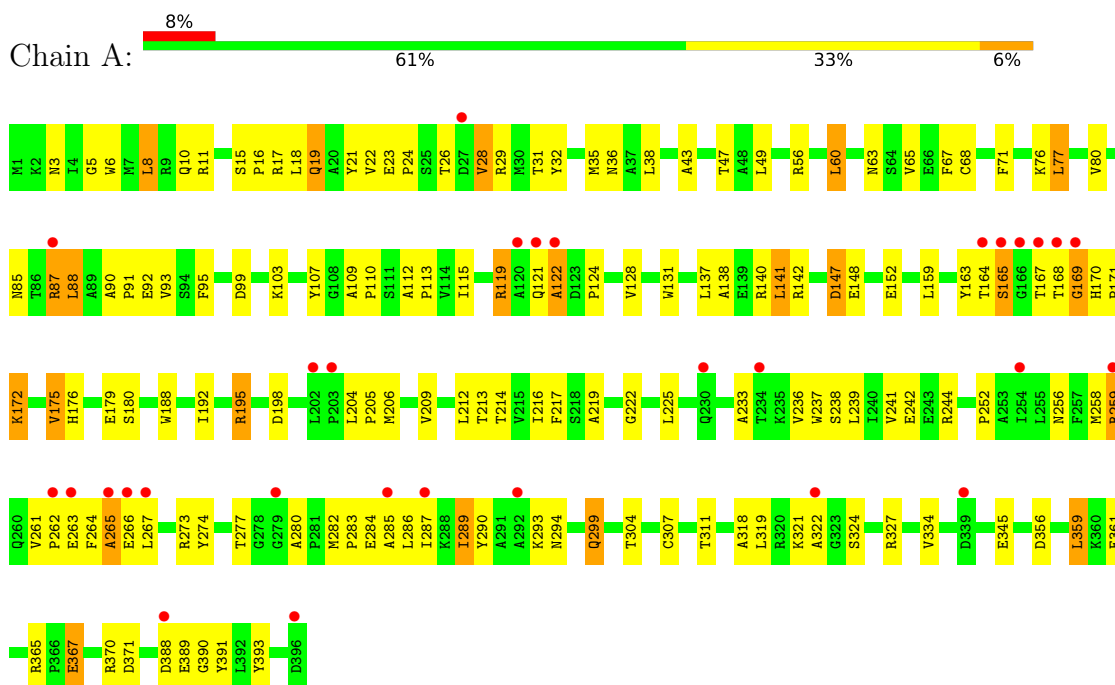
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	287	Total	O	0	0
			287	287		

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: PROBABLE CHAIN-FATTY-ACID-CoA LIGASE FADD13



4 Data and refinement statistics i

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	84.88Å 92.84Å 96.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.35 24.91 – 2.34	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.35) 99.1 (24.91-2.34)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.33Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.210 , 0.276 0.214 , 0.223	Depositor DCC
R_{free} test set	829 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	28.2	Xtrriage
Anisotropy	0.301	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.021 for -h,-l,-k	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3294	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.35	0/3070	0.62	0/4175

There are no bond length outliers.

There are no bond angle outliers.

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	3007	0	2984	126	0
2	A	287	0	0	10	0
All	All	3294	0	2984	126	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 21.

All (126) 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:23:GLU:HB3	1:A:28:VAL:HG13	1.39	1.02
1:A:282:MET:H	1:A:299:GLN:HE22	1.09	0.99
1:A:168:THR:HG22	1:A:169:GLY:H	1.33	0.93
1:A:287:ILE:HD13	1:A:318:ALA:HB1	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:ARG:HG2	1:A:244:ARG:HH11	1.48	0.79
1:A:168:THR:HG22	1:A:169:GLY:N	1.99	0.77
1:A:23:GLU:HB3	1:A:28:VAL:CG1	2.13	0.76
1:A:49:LEU:HD11	1:A:140:ARG:NH1	2.02	0.73
1:A:87:ARG:HA	1:A:87:ARG:HE	1.53	0.73
1:A:11:ARG:HD2	1:A:219:ALA:O	1.90	0.71
1:A:266:GLU:HA	1:A:293:LYS:NZ	2.05	0.71
1:A:282:MET:H	1:A:299:GLN:NE2	1.87	0.71
1:A:22:VAL:HG13	1:A:29:ARG:HG2	1.71	0.70
1:A:122:ALA:O	1:A:124:PRO:HD3	1.92	0.69
1:A:121:GLN:HB2	2:A:659:HOH:O	1.92	0.68
1:A:259:ARG:CZ	1:A:289:ILE:HB	2.23	0.68
1:A:284:GLU:HG3	1:A:319:LEU:HD22	1.76	0.66
1:A:17:ARG:HA	1:A:31:THR:HG21	1.78	0.65
1:A:327:ARG:HD2	1:A:391:TYR:CZ	2.33	0.64
1:A:63:ASN:HB3	1:A:206:MET:HE1	1.80	0.64
1:A:138:ALA:O	1:A:142:ARG:HG3	1.99	0.62
1:A:179:GLU:HG2	2:A:407:HOH:O	1.99	0.62
1:A:180:SER:HA	1:A:356:ASP:O	2.00	0.62
1:A:259:ARG:NH1	1:A:264:PHE:CZ	2.69	0.61
1:A:264:PHE:O	1:A:265:ALA:C	2.37	0.61
1:A:43:ALA:O	1:A:47:THR:HG23	2.00	0.61
1:A:213:THR:HG21	1:A:307:CYS:O	2.03	0.58
1:A:259:ARG:NH1	1:A:289:ILE:CG2	2.67	0.57
1:A:259:ARG:HD3	1:A:289:ILE:HD12	1.85	0.57
1:A:261:VAL:HG23	1:A:262:PRO:HD2	1.85	0.57
1:A:311:THR:CG2	1:A:334:VAL:HG11	2.34	0.57
1:A:152:GLU:HG2	2:A:620:HOH:O	2.05	0.57
1:A:311:THR:HG22	1:A:334:VAL:HG11	1.87	0.57
1:A:256:ASN:HA	1:A:286:LEU:HD11	1.86	0.57
1:A:282:MET:N	1:A:299:GLN:HE22	1.90	0.56
1:A:11:ARG:HD3	1:A:222:GLY:HA2	1.86	0.56
1:A:112:ALA:N	1:A:113:PRO:HD2	2.21	0.55
1:A:238:SER:O	1:A:242:GLU:HG3	2.07	0.55
1:A:76:LYS:NZ	1:A:148:GLU:OE2	2.37	0.55
1:A:261:VAL:CG2	1:A:262:PRO:HD2	2.37	0.55
1:A:233:ALA:HA	1:A:236:VAL:HG23	1.90	0.54
1:A:252:PRO:HD3	1:A:277:THR:HG23	1.90	0.54
1:A:175:VAL:HG13	1:A:361:GLU:O	2.09	0.53
1:A:163:TYR:HA	1:A:172:LYS:O	2.09	0.53
1:A:261:VAL:HG21	1:A:263:GLU:OE1	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ASN:HB3	1:A:206:MET:CE	2.39	0.52
1:A:109:ALA:N	1:A:110:PRO:HD2	2.25	0.52
1:A:8:LEU:CD1	1:A:36:ASN:HA	2.40	0.52
1:A:367:GLU:O	1:A:370:ARG:HG2	2.10	0.52
1:A:16:PRO:O	1:A:31:THR:HG21	2.09	0.51
1:A:88:LEU:HD13	1:A:93:VAL:HG23	1.92	0.51
1:A:168:THR:CG2	1:A:169:GLY:H	2.06	0.51
1:A:16:PRO:O	1:A:31:THR:CG2	2.58	0.51
1:A:209:VAL:HG11	1:A:304:THR:HA	1.91	0.51
1:A:68:CYS:HA	1:A:71:PHE:CE2	2.45	0.51
1:A:99:ASP:OD2	1:A:365:ARG:NH2	2.43	0.51
1:A:389:GLU:HG3	2:A:507:HOH:O	2.10	0.51
1:A:252:PRO:HG2	1:A:280:ALA:O	2.11	0.50
1:A:168:THR:O	1:A:170:HIS:N	2.44	0.50
1:A:85:ASN:HB3	1:A:88:LEU:HG	1.93	0.50
1:A:259:ARG:NH2	1:A:290:TYR:HA	2.27	0.50
1:A:327:ARG:HD2	1:A:391:TYR:CE1	2.47	0.50
1:A:147:ASP:HB3	2:A:678:HOH:O	2.11	0.49
1:A:261:VAL:HG22	1:A:263:GLU:H	1.77	0.49
1:A:122:ALA:C	1:A:124:PRO:HD3	2.32	0.49
1:A:244:ARG:HG2	1:A:244:ARG:NH1	2.17	0.49
1:A:29:ARG:HG3	1:A:29:ARG:HH11	1.77	0.49
1:A:285:ALA:O	1:A:289:ILE:HG13	2.13	0.49
1:A:258:MET:O	1:A:261:VAL:HG12	2.14	0.48
1:A:195:ARG:HG2	1:A:198:ASP:OD2	2.12	0.48
1:A:119:ARG:HG3	1:A:119:ARG:NH1	2.27	0.48
1:A:171:PRO:HG2	2:A:520:HOH:O	2.14	0.48
1:A:90:ALA:N	1:A:91:PRO:HD2	2.29	0.48
1:A:18:LEU:HD12	1:A:19:GLN:H	1.79	0.47
1:A:23:GLU:CB	1:A:28:VAL:HG13	2.27	0.47
1:A:176:HIS:CE1	1:A:307:CYS:HB2	2.50	0.47
1:A:26:THR:HG22	1:A:26:THR:O	2.14	0.47
1:A:31:THR:HG22	1:A:32:TYR:N	2.29	0.47
1:A:261:VAL:HG13	1:A:264:PHE:H	1.80	0.47
1:A:21:TYR:HB2	1:A:35:MET:CE	2.45	0.47
1:A:259:ARG:NH1	1:A:289:ILE:HG22	2.30	0.47
1:A:345:GLU:HG2	2:A:624:HOH:O	2.15	0.46
1:A:21:TYR:HB2	1:A:35:MET:HE1	1.96	0.46
1:A:259:ARG:CD	1:A:289:ILE:HD12	2.44	0.46
1:A:237:TRP:O	1:A:241:VAL:HG13	2.16	0.46
1:A:119:ARG:HG3	1:A:119:ARG:HH11	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:TRP:O	1:A:192:ILE:HG12	2.15	0.46
1:A:214:THR:HA	1:A:217:PHE:CE2	2.52	0.45
1:A:267:LEU:HD12	1:A:267:LEU:HA	1.81	0.45
1:A:47:THR:CG2	1:A:77:LEU:HD21	2.48	0.44
1:A:6:TRP:O	1:A:10:GLN:HG3	2.18	0.44
1:A:112:ALA:N	1:A:113:PRO:CD	2.81	0.44
1:A:321:LYS:HG2	1:A:391:TYR:CD2	2.53	0.44
1:A:35:MET:CE	1:A:65:VAL:HG23	2.48	0.43
1:A:91:PRO:HG2	1:A:92:GLU:OE1	2.18	0.43
1:A:164:THR:O	1:A:165:SER:C	2.55	0.43
1:A:212:LEU:O	1:A:216:ILE:HG12	2.18	0.43
1:A:273:ARG:HB3	1:A:274:TYR:HD1	1.82	0.43
1:A:327:ARG:HG2	1:A:390:GLY:O	2.19	0.43
1:A:204:LEU:HG	1:A:205:PRO:HD2	1.99	0.43
1:A:60:LEU:O	1:A:107:TYR:HA	2.18	0.43
1:A:24:PRO:HG2	1:A:239:LEU:CD1	2.49	0.42
1:A:266:GLU:HA	1:A:293:LYS:HZ1	1.81	0.42
1:A:169:GLY:O	1:A:170:HIS:CG	2.72	0.42
1:A:95:PHE:CE2	1:A:171:PRO:HD2	2.54	0.42
1:A:233:ALA:HA	1:A:236:VAL:CG2	2.50	0.42
1:A:238:SER:HA	1:A:241:VAL:HG22	2.01	0.42
1:A:284:GLU:HG2	2:A:458:HOH:O	2.19	0.42
1:A:128:VAL:HG11	1:A:131:TRP:CE2	2.55	0.42
1:A:60:LEU:HD13	1:A:115:ILE:HD11	2.02	0.41
1:A:103:LYS:HE2	2:A:423:HOH:O	2.19	0.41
1:A:359:LEU:H	1:A:359:LEU:CD1	2.33	0.41
1:A:31:THR:HG22	1:A:32:TYR:H	1.85	0.41
1:A:137:LEU:HG	1:A:141:LEU:HD22	2.02	0.41
1:A:259:ARG:NH1	1:A:259:ARG:HB2	2.35	0.41
1:A:370:ARG:HG3	1:A:371:ASP:N	2.35	0.41
1:A:56:ARG:HA	1:A:80:VAL:O	2.21	0.41
1:A:264:PHE:CE1	1:A:293:LYS:HG3	2.55	0.41
1:A:88:LEU:HD13	1:A:93:VAL:CG2	2.50	0.41
1:A:283:PRO:HG2	1:A:286:LEU:HB2	2.02	0.41
1:A:324:SER:HB2	1:A:393:TYR:CZ	2.55	0.41
1:A:110:PRO:HG2	2:A:558:HOH:O	2.21	0.41
1:A:172:LYS:HD3	1:A:365:ARG:HH11	1.86	0.41
1:A:121:GLN:O	1:A:122:ALA:HB3	2.22	0.40
1:A:3:ASN:OD1	1:A:5:GLY:N	2.54	0.40
1:A:225:LEU:HD12	1:A:225:LEU:HA	1.87	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	394/396 (100%)	368 (93%)	20 (5%)	6 (2%)	10 8

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	167	THR
1	A	265	ALA
1	A	322	ALA
1	A	165	SER
1	A	122	ALA
1	A	169	GLY

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	314/314 (100%)	290 (92%)	24 (8%)	13 13

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LEU
1	A	15	SER
1	A	19	GLN
1	A	28	VAL
1	A	38	LEU

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Mol	Chain	Res	Type
1	A	60	LEU
1	A	67	PHE
1	A	77	LEU
1	A	87	ARG
1	A	88	LEU
1	A	119	ARG
1	A	141	LEU
1	A	147	ASP
1	A	159	LEU
1	A	172	LYS
1	A	175	VAL
1	A	195	ARG
1	A	259	ARG
1	A	289	ILE
1	A	294	ASN
1	A	299	GLN
1	A	359	LEU
1	A	367	GLU
1	A	388	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	256	ASN
1	A	294	ASN
1	A	299	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	396/396 (100%)	0.39	30 (7%) 13 21	13, 28, 54, 75	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	168	THR	10.5
1	A	167	THR	9.2
1	A	169	GLY	8.1
1	A	166	GLY	7.8
1	A	265	ALA	4.6
1	A	259	ARG	4.3
1	A	122	ALA	4.1
1	A	339	ASP	3.6
1	A	285	ALA	3.5
1	A	165	SER	3.2
1	A	266	GLU	3.2
1	A	287	ILE	3.1
1	A	121	GLN	3.0
1	A	396	ASP	2.8
1	A	87	ARG	2.8
1	A	254	ILE	2.7
1	A	388	ASP	2.7
1	A	279	GLY	2.7
1	A	322	ALA	2.6
1	A	164	THR	2.5
1	A	120	ALA	2.5
1	A	230	GLN	2.5
1	A	234	THR	2.5
1	A	203	PRO	2.4
1	A	27	ASP	2.3
1	A	262	PRO	2.1
1	A	267	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	263	GLU	2.0
1	A	202	LEU	2.0
1	A	292	ALA	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 monosaccharides 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.