



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 01:19 PM UTC

PDB ID : 4QGK / pdb_00004qgk
Title : Structure of the Human Sjogren Larsson Syndrome enzyme fatty aldehyde dehydrogenase (FALDH)
Authors : Zander, U.; Keller, M.; Marquez, J.A.
Deposited on : 2014-05-23
Resolution : 2.10 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
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

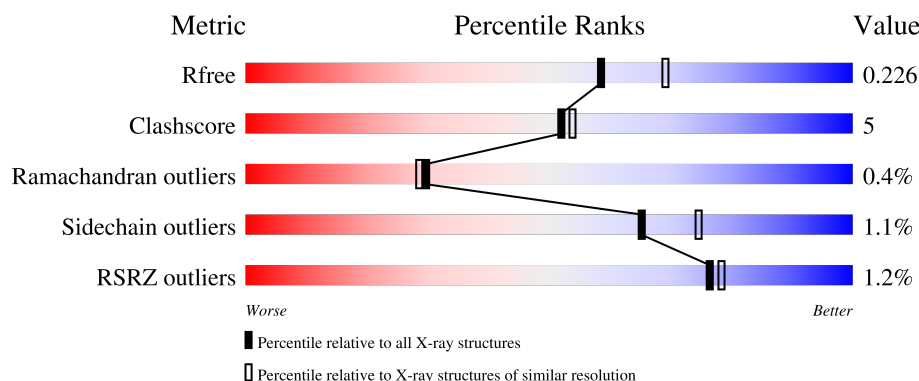
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

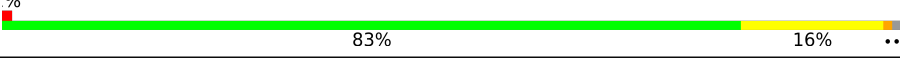
The reported resolution of this entry is 2.10 Å.

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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	461	 2% 85% 12% .
1	B	461	 % 83% 16% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7666 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 Fatty aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	461	Total	C	N	O	S	0	2	0
			3689	2370	628	673	18			
1	B	458	Total	C	N	O	S	0	2	0
			3657	2351	620	668	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	expression tag	UNP P51648
B	0	ALA	-	expression tag	UNP P51648

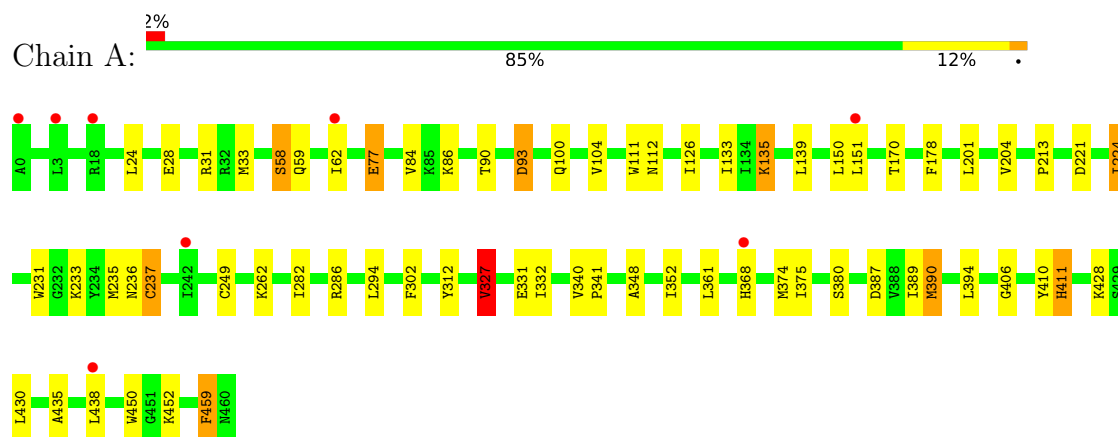
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	152	Total	O	0	0
			152	152		
2	B	168	Total	O	0	0
			168	168		

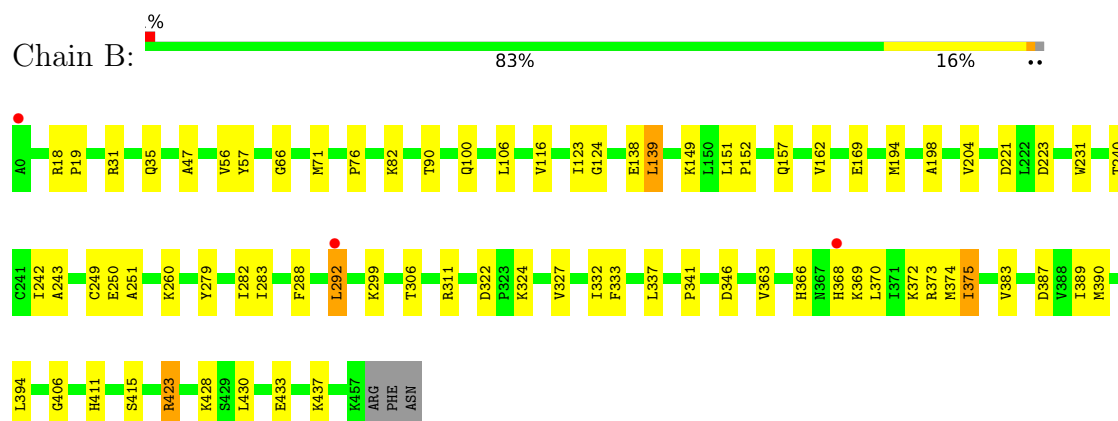
3 Residue-property plots [i](#)

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: Fatty aldehyde dehydrogenase



- Molecule 1: Fatty aldehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.14Å 98.58Å 145.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 30.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.1 (30.00-2.10) 97.1 (30.00-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.198 , 0.223 0.206 , 0.226	Depositor DCC
R_{free} test set	3355 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	33.3	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 25.1	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	7666	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 3.98% 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	1.38	13/3766 (0.3%)	1.14	10/5096 (0.2%)
1	B	1.40	15/3733 (0.4%)	1.18	11/5053 (0.2%)
All	All	1.39	28/7499 (0.4%)	1.16	21/10149 (0.2%)

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	331	GLU	C-O	-6.97	1.15	1.24
1	A	204	VAL	C-O	-6.73	1.17	1.24
1	B	66	GLY	N-CA	6.56	1.54	1.45
1	B	337	LEU	N-CA	-6.54	1.39	1.46
1	A	58	SER	C-O	-6.12	1.16	1.24
1	B	322	ASP	CA-C	6.10	1.59	1.52
1	A	340	VAL	CA-C	-5.97	1.48	1.53
1	B	106	LEU	C-O	-5.92	1.16	1.24
1	B	35	GLN	C-O	-5.70	1.17	1.24
1	B	139	LEU	C-O	-5.70	1.16	1.24
1	B	76	PRO	CA-CB	5.57	1.61	1.53
1	B	423	ARG	CD-NE	-5.47	1.38	1.46
1	A	93	ASP	C-O	-5.46	1.16	1.23
1	B	19	PRO	CA-CB	5.43	1.60	1.53
1	A	86	LYS	C-O	-5.41	1.16	1.23
1	A	100	GLN	CA-C	-5.41	1.47	1.52
1	B	279	TYR	C-O	-5.36	1.17	1.24
1	B	169	GLU	CA-C	5.34	1.59	1.52
1	A	282	ILE	C-O	-5.28	1.18	1.24
1	A	77	GLU	C-O	-5.25	1.17	1.24
1	B	47	ALA	C-O	-5.16	1.18	1.24
1	A	31	ARG	C-O	-5.14	1.18	1.24
1	B	204	VAL	C-O	-5.12	1.18	1.24
1	A	450	TRP	CA-C	5.11	1.59	1.52
1	B	198	ALA	C-O	-5.04	1.18	1.24
1	B	292	LEU	N-CA	5.03	1.52	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	312	TYR	N-CA	-5.02	1.40	1.46
1	A	100	GLN	C-O	-5.00	1.18	1.24

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	56	VAL	N-CA-C	-6.31	104.36	110.42
1	B	283	ILE	N-CA-C	6.28	117.03	110.62
1	A	327	VAL	CB-CA-C	-5.99	104.05	112.14
1	B	251	ALA	N-CA-C	5.92	118.49	111.33
1	A	406	GLY	N-CA-C	5.89	119.33	110.97
1	A	459	PHE	N-CA-C	5.68	118.51	107.71
1	A	84	VAL	CB-CA-C	-5.62	103.49	111.19
1	B	82	LYS	CA-C-N	5.51	125.52	119.90
1	B	82	LYS	C-N-CA	5.51	125.52	119.90
1	B	406	GLY	N-CA-C	5.42	117.87	111.63
1	A	237	CYS	CB-CA-C	-5.35	103.05	111.51
1	B	18	ARG	CG-CD-NE	-5.33	100.27	112.00
1	A	126	ILE	CB-CA-C	-5.24	105.27	111.97
1	B	157	GLN	N-CA-C	5.21	118.89	112.54
1	A	224	ILE	CB-CA-C	-5.19	105.33	111.97
1	A	104	VAL	CB-CA-C	-5.15	104.73	110.96
1	A	150	LEU	N-CA-C	5.14	118.02	111.69
1	B	116	VAL	CB-CA-C	-5.14	105.39	111.97
1	A	112	ASN	N-CA-C	5.05	116.87	111.36
1	B	322	ASP	CA-C-N	5.04	124.76	119.82
1	B	322	ASP	C-N-CA	5.04	124.76	119.82

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	3689	0	3759	41	0
1	B	3657	0	3732	48	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	152	0	0	2	0
2	B	168	0	0	8	0
All	All	7666	0	7491	76	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 5.

All (76) 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:288:PHE:CE1	1:B:292:LEU:HD11	2.01	0.95
1:A:201:LEU:HA	1:B:194:MET:HE1	1.55	0.85
1:B:151:LEU:HD12	1:B:162:VAL:CG2	2.16	0.76
1:A:390:MET:HA	1:A:390:MET:HE2	1.68	0.75
1:A:58:SER:HA	1:A:62:ILE:HD12	1.67	0.75
1:A:111:TRP:CD1	1:A:139:LEU:HD12	2.24	0.72
1:B:250:GLU:HG2	2:B:628:HOH:O	1.93	0.67
1:B:390:MET:HA	1:B:390:MET:HE2	1.75	0.66
1:A:24:LEU:O	1:A:28:GLU:HG2	1.95	0.66
1:B:151:LEU:HD12	1:B:162:VAL:HG23	1.78	0.66
1:A:201:LEU:HD23	1:B:194:MET:CE	2.25	0.66
1:B:327:VAL:O	1:B:332:ILE:HD11	1.98	0.63
1:B:288:PHE:HE1	1:B:292:LEU:HD11	1.59	0.62
1:B:149:LYS:HE2	2:B:556:HOH:O	2.00	0.62
1:A:327:VAL:HG12	1:A:332:ILE:HD11	1.84	0.60
1:B:31:ARG:HD2	2:B:622:HOH:O	2.03	0.58
1:B:223:ASP:OD1	1:B:260:LYS:HE3	2.05	0.57
1:A:294:LEU:O	1:A:327:VAL:HG22	2.05	0.57
1:B:231:TRP:HH2	1:B:390:MET:HE3	1.69	0.57
1:B:324:LYS:HD2	2:B:629:HOH:O	2.04	0.56
1:A:380:SER:O	1:B:423:ARG:HD3	2.05	0.56
1:A:33:MET:CE	1:A:151:LEU:HA	2.38	0.54
1:B:363:VAL:CG2	1:B:374:MET:HG2	2.38	0.54
1:A:33:MET:HE1	1:A:151:LEU:HA	1.91	0.53
1:A:428:LYS:HD3	1:B:387:ASP:HB2	1.91	0.52
1:B:100:GLN:HG2	2:B:623:HOH:O	2.10	0.52
1:A:262:LYS:HG3	1:A:302:PHE:CE1	2.45	0.52
1:A:387:ASP:HB2	1:B:428:LYS:HD3	1.93	0.51
1:A:390:MET:HA	1:A:390:MET:CE	2.39	0.51
1:B:363:VAL:HG21	1:B:374:MET:HG2	1.92	0.51
1:A:201:LEU:HD23	1:B:194:MET:HE3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:ILE:O	1:B:423:ARG:NH2	2.44	0.51
1:A:135:LYS:HD2	1:A:135:LYS:C	2.36	0.51
1:A:90:THR:HG21	1:B:394:LEU:HD11	1.92	0.51
1:B:231:TRP:CH2	1:B:390:MET:HE3	2.48	0.49
1:A:236:ASN:O	1:A:237:CYS:HB2	2.11	0.49
1:B:292:LEU:CD2	1:B:306:THR:HG21	2.44	0.46
1:A:133:ILE:HD12	1:A:178:PHE:HE1	1.80	0.46
1:A:135:LYS:HE2	1:A:170:THR:OG1	2.15	0.46
1:A:435:ALA:O	1:A:438:LEU:HD13	2.15	0.46
1:B:437:LYS:HD2	2:B:647:HOH:O	2.15	0.46
1:A:452:LYS:HG2	1:B:57:TYR:OH	2.16	0.46
1:B:369:LYS:HD2	2:B:585:HOH:O	2.16	0.46
1:A:348:ALA:O	1:A:352[A]:ILE:HG12	2.17	0.45
1:A:224:ILE:HD12	1:B:433:GLU:HB2	1.98	0.45
1:A:294:LEU:O	1:A:327:VAL:CG2	2.64	0.45
1:B:282:ILE:O	1:B:311:ARG:NH1	2.39	0.45
1:A:430:LEU:HD23	1:B:366:HIS:CE1	2.51	0.45
1:A:233:LYS:HD2	1:A:233:LYS:O	2.16	0.45
1:A:249:CYS:O	1:A:341:PRO:HA	2.17	0.45
1:B:221:ASP:OD1	1:B:221:ASP:C	2.58	0.45
1:B:375:ILE:CD1	1:B:383:VAL:HG11	2.46	0.45
1:B:375:ILE:HD11	1:B:383:VAL:HG11	1.98	0.44
1:A:394:LEU:HD11	1:B:90:THR:HG21	2.00	0.44
1:B:151:LEU:HB2	1:B:152:PRO:HD3	2.00	0.44
1:B:346:ASP:OD1	1:B:373:ARG:NH2	2.43	0.44
1:A:327:VAL:HG12	1:A:332:ILE:CD1	2.47	0.43
1:A:59:GLN:NE2	1:A:235:MET:HE2	2.34	0.43
1:B:368:HIS:O	1:B:372:LYS:HG2	2.19	0.43
1:B:370:LEU:C	1:B:370:LEU:HD23	2.43	0.42
1:A:93:ASP:CG	1:A:428:LYS:HG2	2.45	0.42
1:A:410:TYR:O	1:A:411:HIS:HB2	2.20	0.42
1:B:249:CYS:O	1:B:341:PRO:HA	2.20	0.42
1:B:124:GLY:HA3	1:B:415:SER:HB3	2.01	0.42
1:B:437:LYS:CE	2:B:647:HOH:O	2.66	0.42
1:B:240:THR:HB	1:B:243:ALA:HB2	2.01	0.42
1:B:242[A]:ILE:HD12	1:B:242[A]:ILE:HG23	1.74	0.41
1:A:231:TRP:HH2	1:A:390:MET:HE3	1.86	0.41
1:A:213:PRO:HG3	1:A:352[A]:ILE:HD12	2.02	0.41
1:B:138:GLU:HG3	1:B:139:LEU:HD23	2.03	0.41
1:A:286:ARG:CD	2:A:606:HOH:O	2.68	0.41
1:A:368:HIS:CE1	2:A:623:HOH:O	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:71:MET:HG3	1:B:123:ILE:CG2	2.52	0.40
1:A:374:MET:HE2	1:A:374:MET:HB3	1.84	0.40
1:B:242[A]:ILE:O	1:B:242[A]:ILE:HG22	2.21	0.40
1:A:221:ASP:H	1:B:430:LEU:HD13	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	461/461 (100%)	455 (99%)	4 (1%)	2 (0%)	30	28
1	B	458/461 (99%)	448 (98%)	8 (2%)	2 (0%)	30	28
All	All	919/922 (100%)	903 (98%)	12 (1%)	4 (0%)	30	28

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	389	ILE
1	B	411	HIS
1	A	411	HIS
1	B	389	ILE

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	405/403 (100%)	399 (98%)	6 (2%)	57	65
1	B	402/403 (100%)	399 (99%)	3 (1%)	76	83
All	All	807/806 (100%)	798 (99%)	9 (1%)	65	74

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	GLU
1	A	135	LYS
1	A	327	VAL
1	A	361	LEU
1	A	390	MET
1	A	459	PHE
1	B	299	LYS
1	B	333	PHE
1	B	375	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	200	HIS
1	B	100	GLN
1	B	421	HIS

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 oligosaccharides 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	461/461 (100%)	-0.04	8 (1%) 69 71	16, 41, 63, 99	2 (0%)
1	B	458/461 (99%)	-0.14	3 (0%) 84 86	18, 40, 60, 85	2 (0%)
All	All	919/922 (99%)	-0.09	11 (1%) 76 78	16, 41, 62, 99	4 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	151	LEU	3.2
1	B	0	ALA	2.7
1	A	62	ILE	2.5
1	A	242	ILE	2.5
1	A	0	ALA	2.4
1	A	3	LEU	2.4
1	A	368	HIS	2.4
1	A	438	LEU	2.3
1	B	292	LEU	2.2
1	A	18	ARG	2.0
1	B	368	HIS	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.