



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:25 AM UTC

PDB ID : 1FJ2 / pdb_00001fj2
Title : Crystal structure of the human acyl protein thioesterase 1 at 1.5 Å resolution
Authors : Devedjiev, Y.; Dauter, Z.; Kuznetsov, S.; Jones, T.; Derewenda, Z.
Deposited on : 2000-08-07
Resolution : 1.50 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

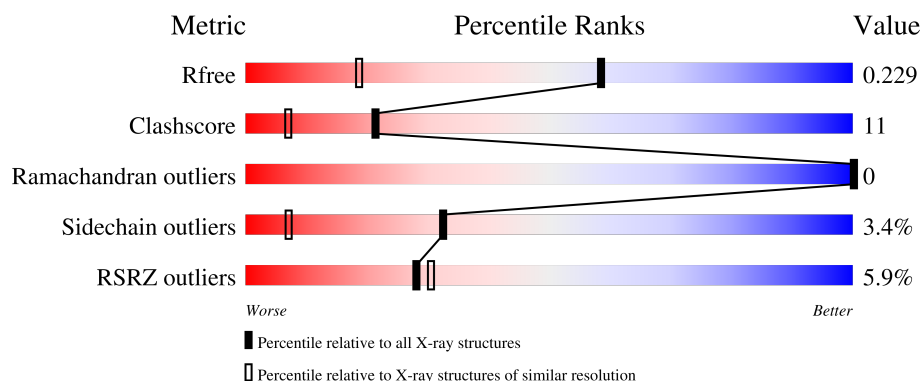
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	232	5% 75% 20% ..
1	B	232	6% 79% 16% ..

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
2	BR	B	839	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3959 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 PROTEIN (ACYL PROTEIN THIOESTERASE 1).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	S	0	0	0
			1727	1102	292	319	14			
1	B	229	Total	C	N	O	S	0	0	0
			1727	1102	292	319	14			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	see remark 999	UNP O75608
A	-6	ALA	-	see remark 999	UNP O75608
A	-5	MET	-	see remark 999	UNP O75608
A	-4	ASP	-	see remark 999	UNP O75608
A	-3	PRO	-	see remark 999	UNP O75608
A	-2	GLU	-	see remark 999	UNP O75608
A	-1	PHE	-	see remark 999	UNP O75608

- Molecule 2 is BROMIDE ION (CCD ID: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	19	Total	Br	0	0
			19	19		
2	B	21	Total	Br	0	0
			21	21		

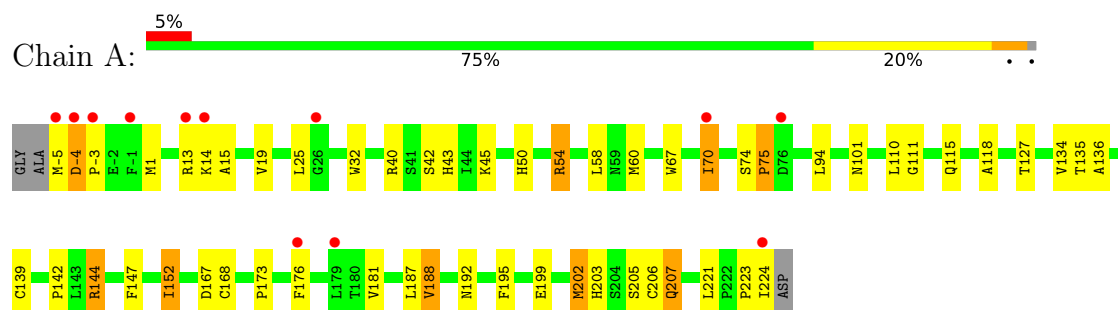
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	222	Total	O	6	0
			222	222		
3	B	243	Total	O	9	0
			243	243		

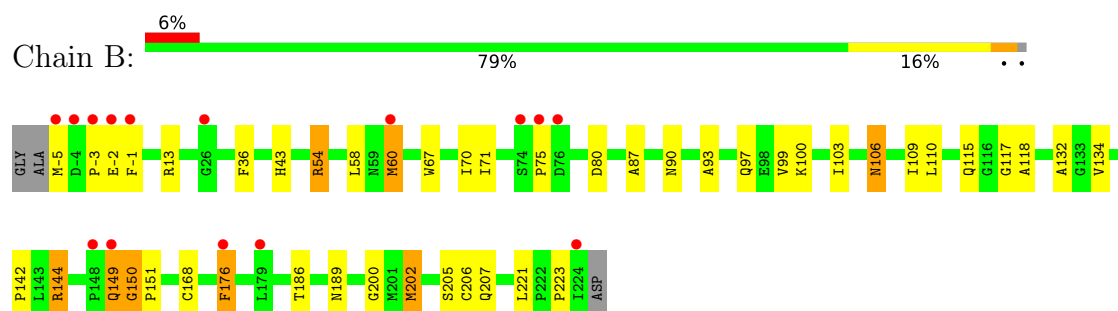
3 Residue-property plots [i](#)

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: PROTEIN (ACYL PROTEIN THIOESTERASE 1)



• Molecule 1: PROTEIN (ACYL PROTEIN THIOESTERASE 1)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.59Å 127.89Å 39.66Å 90.00° 102.80° 90.00°	Depositor
Resolution (Å)	20.00 – 1.50 20.00 – 1.50	Depositor EDS
% Data completeness (in resolution range)	80.0 (20.00-1.50) 70.2 (20.00-1.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 1.48Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.183 , 0.237 0.179 , 0.229	Depositor DCC
R_{free} test set	1303 reflections (2.57%)	wwPDB-VP
Wilson B-factor (Å ²)	11.6	Xtriage
Anisotropy	0.226	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 48.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.119 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3959	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.86% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BR

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	1.21	2/1768 (0.1%)	1.80	30/2404 (1.2%)
1	B	1.20	1/1768 (0.1%)	1.76	18/2404 (0.7%)
All	All	1.20	3/3536 (0.1%)	1.78	48/4808 (1.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	202	MET	SD-CE	-10.74	1.52	1.79
1	A	202	MET	SD-CE	-9.17	1.56	1.79
1	A	203	HIS	ND1-CE1	6.64	1.39	1.32

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	VAL	N-CA-CB	-18.39	91.85	112.45
1	B	223	PRO	CA-C-N	13.96	146.82	121.70
1	B	223	PRO	C-N-CA	13.96	146.82	121.70
1	A	223	PRO	CA-C-N	11.39	142.21	121.70
1	A	223	PRO	C-N-CA	11.39	142.21	121.70
1	A	188	VAL	CB-CA-C	9.37	124.45	111.34
1	A	-4	ASP	CA-CB-CG	9.35	121.95	112.60
1	A	40	ARG	CG-CD-NE	8.28	130.21	112.00
1	A	127	THR	CA-C-N	-8.01	109.84	122.73
1	A	127	THR	C-N-CA	-8.01	109.84	122.73
1	A	32	TRP	CA-C-O	-7.37	112.74	120.55
1	A	54	ARG	CA-CB-CG	7.32	128.75	114.10
1	A	40	ARG	NE-CZ-NH2	7.06	125.56	119.20
1	A	45	LYS	CA-CB-CG	7.02	128.14	114.10
1	A	54	ARG	CB-CG-CD	6.88	127.12	111.30
1	B	149	GLN	N-CA-C	-6.87	101.44	110.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	189	ASN	OD1-CG-ND2	6.71	129.31	122.60
1	B	207	GLN	OE1-CD-NE2	6.55	129.15	122.60
1	A	188	VAL	O-C-N	-6.54	116.05	122.92
1	B	144	ARG	NE-CZ-NH1	-6.39	115.11	121.50
1	B	70	ILE	O-C-N	6.20	129.96	123.26
1	A	147	PHE	CA-CB-CG	-6.08	107.72	113.80
1	B	80	ASP	CA-CB-CG	6.01	118.61	112.60
1	A	195	PHE	CA-CB-CG	5.86	119.66	113.80
1	B	87	ALA	CA-C-O	-5.81	114.39	120.55
1	A	224	ILE	N-CA-CB	5.77	121.31	111.50
1	A	19	VAL	CA-C-O	-5.61	114.55	120.39
1	A	192	ASN	OD1-CG-ND2	5.58	128.18	122.60
1	A	75	PRO	CA-C-N	5.47	132.24	121.58
1	A	75	PRO	C-N-CA	5.47	132.24	121.58
1	B	149	GLN	CB-CG-CD	5.46	121.88	112.60
1	A	40	ARG	NE-CZ-NH1	-5.45	116.05	121.50
1	A	101	ASN	CA-CB-CG	-5.45	107.15	112.60
1	B	189	ASN	CB-CG-ND2	-5.42	108.27	116.40
1	B	150	GLY	CA-C-N	5.38	125.30	119.76
1	B	150	GLY	C-N-CA	5.38	125.30	119.76
1	A	224	ILE	CA-CB-CG2	-5.35	101.40	110.50
1	A	50	HIS	CA-CB-CG	-5.34	108.46	113.80
1	A	144	ARG	NE-CZ-NH2	-5.34	114.40	119.20
1	B	117	GLY	O-C-N	5.33	127.36	122.19
1	A	207	GLN	CA-C-O	-5.33	115.22	120.82
1	B	60	MET	CA-C-N	5.33	131.62	123.47
1	B	60	MET	C-N-CA	5.33	131.62	123.47
1	A	224	ILE	CB-CA-C	-5.29	101.01	111.60
1	B	36	PHE	CA-CB-CG	-5.27	108.53	113.80
1	A	50	HIS	CA-C-O	-5.19	115.04	120.80
1	B	186	THR	CA-CB-OG1	5.12	117.28	109.60
1	A	54	ARG	O-C-N	-5.00	117.93	121.83

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	1727	0	1743	47	0
1	B	1727	0	1743	36	0
2	A	19	0	0	5	0
2	B	21	0	0	4	0
3	A	222	0	0	8	0
3	B	243	0	0	7	0
All	All	3959	0	3486	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 11.

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:A:-4:ASP:HB2	1:A:-3:PRO:HD2	1.41	1.00
1:B:75:PRO:HB3	1:B:144:ARG:HH12	1.34	0.89
1:B:75:PRO:HB2	1:B:144:ARG:HH22	1.39	0.87
1:A:25:LEU:HD11	1:A:70:ILE:HG21	1.60	0.84
1:A:42:SER:OG	3:A:950:HOH:O	1.97	0.74
1:A:168:CYS:SG	2:A:831:BR:BR	3.03	0.72
1:A:-4:ASP:HB2	1:A:-3:PRO:CD	2.19	0.72
1:A:70:ILE:HB	3:A:960:HOH:O	1.89	0.72
1:B:75:PRO:CB	1:B:144:ARG:HH12	2.02	0.72
1:A:67:TRP:H	1:A:115:GLN:NE2	1.88	0.70
1:A:67:TRP:H	1:A:115:GLN:HE22	1.39	0.69
1:A:60:MET:SD	3:B:1029:HOH:O	2.52	0.68
1:B:43:HIS:HD2	3:B:898:HOH:O	1.76	0.67
2:B:839:BR:BR	3:B:851:HOH:O	2.68	0.66
1:A:144:ARG:HD2	3:A:978:HOH:O	1.97	0.65
1:B:13:ARG:NH2	2:B:803:BR:BR	2.85	0.64
1:A:25:LEU:HD11	1:A:70:ILE:CG2	2.28	0.63
1:B:43:HIS:HE1	1:B:221:LEU:O	1.82	0.62
1:A:60:MET:HG2	3:A:1040:HOH:O	2.01	0.61
1:A:206:CYS:CA	1:B:202:MET:HE2	2.29	0.61
1:A:25:LEU:CD1	1:A:70:ILE:HG21	2.30	0.59
1:A:152:ILE:HG23	1:A:187:LEU:HD22	1.83	0.59
2:A:830:BR:BR	3:A:963:HOH:O	2.72	0.58
1:B:75:PRO:HB2	1:B:144:ARG:NH2	2.14	0.58
1:A:43:HIS:HE1	1:A:221:LEU:O	1.86	0.58
1:B:106:ASN:HD22	1:B:106:ASN:C	2.11	0.57
1:B:202:MET:SD	3:B:958:HOH:O	2.56	0.57
1:A:202:MET:HE2	1:B:206:CYS:CA	2.33	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:TRP:H	1:B:115:GLN:NE2	2.02	0.57
1:B:118:ALA:HB1	1:B:142:PRO:HG3	1.86	0.56
3:A:914:HOH:O	1:B:60:MET:HE3	2.06	0.56
1:A:-5:MET:HE1	1:A:94:LEU:HD13	1.86	0.55
1:A:206:CYS:HA	1:B:202:MET:HE2	1.88	0.55
1:B:67:TRP:H	1:B:115:GLN:HE22	1.55	0.55
1:A:173:PRO:HD2	1:A:176:PHE:CE2	2.42	0.54
1:B:110:LEU:O	1:B:134:VAL:HA	2.08	0.54
1:A:202:MET:CE	1:B:205:SER:OG	2.55	0.54
1:A:111:GLY:HA3	1:A:135:THR:HG22	1.90	0.53
1:A:202:MET:HE2	1:B:206:CYS:HA	1.90	0.53
1:A:14:LYS:HG3	3:A:1048:HOH:O	2.11	0.51
1:A:205:SER:OG	2:A:840:BR:BR	2.81	0.51
1:B:75:PRO:HB3	1:B:176:PHE:CZ	2.46	0.50
1:A:205:SER:OG	1:B:202:MET:CE	2.60	0.50
1:A:207:GLN:HE21	1:B:200:GLY:HA2	1.77	0.50
1:A:135:THR:O	1:A:135:THR:HG23	2.12	0.50
1:A:74:SER:HB2	1:A:75:PRO:HD2	1.93	0.49
1:A:13:ARG:HG3	1:A:42:SER:HB2	1.95	0.49
1:A:67:TRP:N	1:A:115:GLN:HE22	2.09	0.48
1:B:93:ALA:O	1:B:97:GLN:HG3	2.14	0.47
1:B:-2:GLU:OE1	1:B:54:ARG:NH1	2.48	0.47
1:A:43:HIS:HD2	3:A:950:HOH:O	1.98	0.46
1:A:136:ALA:HB1	1:A:139:CYS:SG	2.56	0.46
1:A:118:ALA:HB1	1:A:142:PRO:HG3	1.99	0.45
1:A:-5:MET:HE1	1:A:94:LEU:CD1	2.46	0.45
1:A:173:PRO:HD2	1:A:176:PHE:CD2	2.52	0.44
1:A:15:ALA:O	2:A:829:BR:BR	2.91	0.44
1:A:110:LEU:O	1:A:134:VAL:HA	2.18	0.44
1:B:99:VAL:HA	1:B:103:ILE:O	2.18	0.44
2:B:839:BR:BR	3:B:1077:HOH:O	2.76	0.43
1:B:100:LYS:HE2	2:B:839:BR:BR	2.74	0.43
1:A:167:ASP:CG	1:A:199:GLU:HA	2.44	0.42
1:A:-5:MET:HG3	1:A:1:MET:HE3	2.01	0.42
1:A:205:SER:OG	1:B:202:MET:HE3	2.18	0.42
1:A:202:MET:HE3	1:B:205:SER:OG	2.19	0.42
1:B:151:PRO:HB3	3:B:1068:HOH:O	2.20	0.42
1:A:-5:MET:SD	2:A:816:BR:BR	3.33	0.42
1:B:-1:PHE:CD1	1:B:90:ASN:ND2	2.88	0.42
1:A:202:MET:HE1	1:B:205:SER:OG	2.19	0.41
1:B:106:ASN:C	1:B:106:ASN:ND2	2.78	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:CYS:O	1:B:202:MET:HG2	2.20	0.41
1:B:150:GLY:HA2	1:B:151:PRO:HD3	1.78	0.41
1:B:-5:MET:N	1:B:90:ASN:OD1	2.48	0.41
1:A:167:ASP:OD2	1:A:199:GLU:HA	2.19	0.41
1:B:109:ILE:HG13	1:B:132:ALA:HB3	2.03	0.41
1:A:74:SER:HB2	1:A:75:PRO:CD	2.51	0.41
1:A:60:MET:HE2	3:B:1037:HOH:O	2.21	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	227/232 (98%)	220 (97%)	7 (3%)	0	100	100
1	B	227/232 (98%)	223 (98%)	4 (2%)	0	100	100
All	All	454/464 (98%)	443 (98%)	11 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	189/190 (100%)	183 (97%)	6 (3%)	34	8

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	189/190 (100%)	182 (96%)	7 (4%)	30 6
All	All	378/380 (100%)	365 (97%)	13 (3%)	32 7

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ARG
1	A	58	LEU
1	A	70	ILE
1	A	152	ILE
1	A	181	VAL
1	A	188	VAL
1	B	-3	PRO
1	B	54	ARG
1	B	58	LEU
1	B	71	ILE
1	B	106	ASN
1	B	149	GLN
1	B	176	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	HIS
1	A	97	GLN
1	A	101	ASN
1	A	115	GLN
1	A	207	GLN
1	B	43	HIS
1	B	78	GLN
1	B	106	ASN
1	B	115	GLN
1	B	128	GLN
1	B	129	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 40 ligands modelled in this entry, 40 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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 [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	229/232 (98%)	0.00	12 (5%) 33 35	6, 12, 29, 55	0
1	B	229/232 (98%)	0.06	15 (6%) 24 26	6, 12, 30, 50	0
All	All	458/464 (98%)	0.03	27 (5%) 28 30	6, 12, 30, 55	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	224	ILE	5.6
1	B	75	PRO	4.8
1	A	-5	MET	4.7
1	A	-1	PHE	4.4
1	A	26	GLY	4.2
1	A	176	PHE	4.2
1	B	26	GLY	3.7
1	B	76	ASP	3.6
1	B	224	ILE	3.6
1	B	-1	PHE	3.5
1	B	-5	MET	3.4
1	B	-3	PRO	3.3
1	B	60	MET	3.3
1	A	-4	ASP	3.2
1	A	14	LYS	3.1
1	A	179	LEU	3.1
1	B	179	LEU	3.1
1	B	149	GLN	3.1
1	A	76	ASP	2.7
1	A	13	ARG	2.6
1	B	-2	GLU	2.5
1	A	-3	PRO	2.5
1	B	176	PHE	2.5
1	B	148	PRO	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	-4	ASP	2.4
1	B	74	SER	2.4
1	A	70	ILE	2.1

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](#)

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. 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	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BR	B	826	1/1	0.70	0.21	53,53,53,53	1
2	BR	B	837	1/1	0.70	0.24	27,27,27,27	1
2	BR	A	831	1/1	0.71	0.30	41,41,41,41	1
2	BR	B	822	1/1	0.81	0.29	16,16,16,16	1
2	BR	A	828	1/1	0.82	0.27	45,45,45,45	1
2	BR	B	825	1/1	0.86	0.15	38,38,38,38	1
2	BR	A	817	1/1	0.87	0.11	26,26,26,26	1
2	BR	A	832	1/1	0.87	0.15	47,47,47,47	1
2	BR	A	824	1/1	0.88	0.12	50,50,50,50	0
2	BR	B	821	1/1	0.88	0.27	53,53,53,53	0
2	BR	A	816	1/1	0.89	0.17	48,48,48,48	0
2	BR	B	833	1/1	0.89	0.19	22,22,22,22	1
2	BR	A	827	1/1	0.89	0.16	28,28,28,28	1
2	BR	A	830	1/1	0.91	0.11	44,44,44,44	1
2	BR	A	804	1/1	0.93	0.08	27,27,27,27	0
2	BR	A	818	1/1	0.93	0.15	46,46,46,46	0
2	BR	B	836	1/1	0.93	0.19	34,34,34,34	1
2	BR	B	820	1/1	0.93	0.19	56,56,56,56	0
2	BR	A	829	1/1	0.94	0.10	31,31,31,31	1
2	BR	B	834	1/1	0.95	0.16	19,19,19,19	1

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BR	A	809	1/1	0.95	0.10	32,32,32,32	0
2	BR	A	812	1/1	0.95	0.13	32,32,32,32	0
2	BR	B	838	1/1	0.95	0.12	22,22,22,22	1
2	BR	B	835	1/1	0.96	0.11	23,23,23,23	1
2	BR	B	805	1/1	0.97	0.05	31,31,31,31	0
2	BR	B	811	1/1	0.97	0.08	32,32,32,32	0
2	BR	B	815	1/1	0.97	0.12	42,42,42,42	0
2	BR	B	819	1/1	0.97	0.17	43,43,43,43	0
2	BR	A	806	1/1	0.97	0.07	31,31,31,31	0
2	BR	B	803	1/1	0.98	0.04	22,22,22,22	0
2	BR	A	808	1/1	0.98	0.10	29,29,29,29	0
2	BR	B	807	1/1	0.98	0.10	32,32,32,32	0
2	BR	A	813	1/1	0.98	0.07	26,26,26,26	0
2	BR	B	814	1/1	0.98	0.11	35,35,35,35	0
2	BR	A	840	1/1	0.98	0.09	17,17,17,17	1
2	BR	A	801	1/1	0.99	0.03	14,14,14,14	0
2	BR	A	823	1/1	0.99	0.20	38,38,38,38	0
2	BR	B	839	1/1	0.99	0.06	21,21,21,21	1
2	BR	B	802	1/1	1.00	0.03	14,14,14,14	0
2	BR	B	810	1/1	1.00	0.11	28,28,28,28	0

6.5 Other polymers

There are no such residues in this entry.