



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 24, 2026 – 10:11 PM EDT

PDB ID : 1G83 / pdb_00001g83
Title : CRYSTAL STRUCTURE OF FYN SH3-SH2
Authors : Arold, S.T.; Ulmer, T.S.; Mulhern, T.D.; Werner, J.M.; Ladbury, J.E.; Campbell, I.D.; Noble, M.E.M.
Deposited on : 2000-11-16
Resolution : 2.60 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

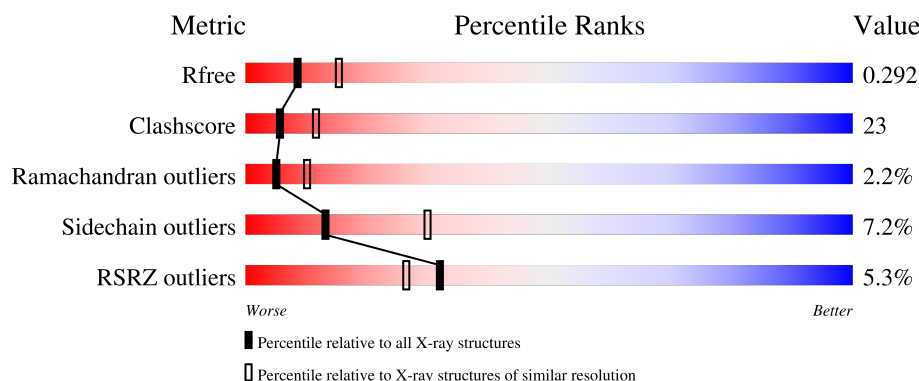
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

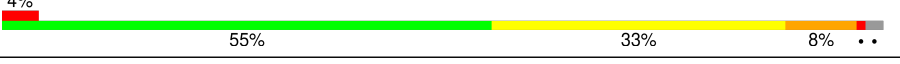
The reported resolution of this entry is 2.60 Å.

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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	165	
1	B	165	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2642 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 PROTO-ONCOGENE TYROSINE-PROTEIN KINASE FYN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	S	0	0	0
			1313	832	226	254	1			
1	B	161	Total	C	N	O	S	0	0	0
			1313	832	226	254	1			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	239	SER	CYS	engineered mutation	UNP P06241
A	240	SER	CYS	engineered mutation	UNP P06241
A	246	SER	CYS	engineered mutation	UNP P06241
B	239	SER	CYS	engineered mutation	UNP P06241
B	240	SER	CYS	engineered mutation	UNP P06241
B	246	SER	CYS	engineered mutation	UNP P06241

- Molecule 2 is water.

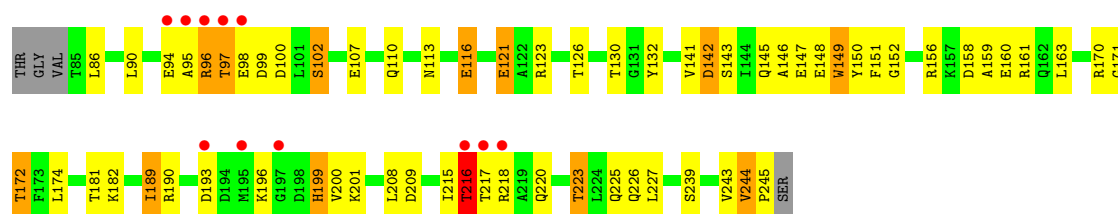
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	8	Total	O	0	0
			8	8		
2	B	8	Total	O	0	0
			8	8		

3 Residue-property plots

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.

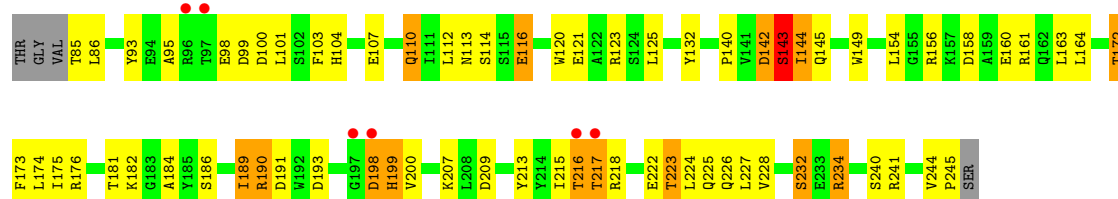
• Molecule 1: PROTO-ONCOGENE TYROSINE-PROTEIN KINASE FYN

Chain A: 



• Molecule 1: PROTO-ONCOGENE TYROSINE-PROTEIN KINASE FYN

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.97Å 89.96Å 60.27Å 90.00° 101.44° 90.00°	Depositor
Resolution (Å)	37.00 – 2.60 37.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	82.0 (37.00-2.60) 81.3 (37.00-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 2.61Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.214 , 0.277 0.228 , 0.292	Depositor DCC
R_{free} test set	506 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	29.9	Xtriage
Anisotropy	0.717	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 46.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2642	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.40% 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

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	0.77	1/1345 (0.1%)	1.18	10/1818 (0.6%)
1	B	0.67	0/1345	1.12	7/1818 (0.4%)
All	All	0.72	1/2690 (0.0%)	1.15	17/3636 (0.5%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	142	ASP	C-N	-5.68	1.26	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	98	GLU	N-CA-C	-9.63	102.33	114.56
1	A	99	ASP	N-CA-C	-8.43	102.97	113.18
1	A	199	HIS	N-CA-C	7.92	119.07	108.38
1	B	101	LEU	N-CA-C	7.21	121.60	110.20
1	B	99	ASP	N-CA-C	-7.02	103.98	112.54
1	B	143	SER	N-CA-C	6.77	125.22	110.80
1	A	98	GLU	N-CA-C	-6.01	106.11	113.50
1	B	199	HIS	N-CA-C	5.89	116.33	108.38
1	A	196	LYS	N-CA-C	-5.83	102.84	110.53
1	A	216	THR	N-CA-C	5.77	123.10	110.80
1	B	110	GLN	N-CA-C	-5.69	100.13	109.40
1	A	163	LEU	N-CA-C	5.66	117.45	111.28
1	A	102	SER	N-CA-C	-5.61	100.74	109.72
1	B	189	ILE	N-CA-C	5.58	116.70	108.45
1	A	172	THR	N-CA-C	-5.35	102.36	110.28
1	A	149	TRP	N-CA-C	-5.14	105.90	113.61
1	A	220	GLN	N-CA-C	5.04	118.22	110.42

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	1313	0	1255	52	0
1	B	1313	0	1255	69	0
2	A	8	0	0	0	0
2	B	8	0	0	2	0
All	All	2642	0	2510	120	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 23.

All (120) 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:223:THR:HG22	1:A:226:GLN:H	1.37	0.87
1:A:223:THR:CG2	1:A:225:GLN:H	1.96	0.78
1:A:171:GLY:HA3	1:A:243:VAL:HG23	1.65	0.77
1:B:149:TRP:CZ3	1:B:175:ILE:HD12	2.21	0.76
1:B:113:ASN:HB3	1:B:121:GLU:HB3	1.69	0.74
1:A:223:THR:HG23	1:A:225:GLN:H	1.53	0.71
1:A:110:GLN:HG2	1:A:123:ARG:HB3	1.72	0.70
1:B:207:LYS:HD3	1:B:213:TYR:CE2	2.27	0.69
1:B:223:THR:HG22	1:B:226:GLN:H	1.59	0.67
1:B:223:THR:HB	1:B:226:GLN:HG3	1.78	0.65
1:A:171:GLY:HA3	1:A:243:VAL:CG2	2.25	0.65
1:A:174:LEU:HD12	1:A:174:LEU:C	2.23	0.64
1:B:100:ASP:HA	1:B:132:TYR:H	1.63	0.64
1:B:223:THR:HG23	1:B:225:GLN:OE1	1.97	0.64
1:A:147:GLU:O	1:A:148:GLU:HB2	1.97	0.63
1:B:223:THR:CG2	1:B:225:GLN:H	2.12	0.61
1:B:154:LEU:HD11	1:B:158:ASP:HB3	1.82	0.61
1:A:216:THR:HG22	1:A:216:THR:O	1.98	0.60
1:A:223:THR:HG23	1:A:225:GLN:OE1	2.01	0.60
1:A:86:LEU:C	1:A:141:VAL:HG23	2.27	0.59
1:B:110:GLN:HG2	1:B:123:ARG:HB3	1.84	0.59
1:B:216:THR:O	1:B:218:ARG:HG3	2.02	0.59
1:A:143:SER:HB3	1:A:145:GLN:OE1	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:193:ASP:HB2	1:B:199:HIS:HE1	1.68	0.59
1:B:232:SER:O	1:B:241:ARG:HD2	2.03	0.58
1:A:96:ARG:HG2	1:A:96:ARG:HH11	1.68	0.58
1:B:223:THR:HG22	1:B:225:GLN:H	1.69	0.58
1:B:198:ASP:C	1:B:198:ASP:OD1	2.47	0.56
1:B:216:THR:OG1	1:B:217:THR:N	2.37	0.56
1:A:149:TRP:HB3	1:A:245:PRO:HB3	1.87	0.56
1:A:95:ALA:HB2	1:A:102:SER:HB2	1.87	0.56
1:A:223:THR:HG22	1:A:225:GLN:N	2.21	0.55
1:B:217:THR:O	1:B:218:ARG:HB2	2.05	0.55
1:A:94:GLU:O	1:A:95:ALA:C	2.50	0.55
1:B:120:TRP:O	1:B:132:TYR:HA	2.08	0.54
1:B:223:THR:HG22	1:B:225:GLN:N	2.22	0.54
1:A:223:THR:HB	1:A:226:GLN:HG3	1.90	0.54
1:A:223:THR:HG22	1:A:225:GLN:H	1.73	0.54
1:B:189:ILE:O	1:B:200:VAL:HA	2.08	0.54
1:B:100:ASP:HB3	1:B:132:TYR:HB2	1.89	0.54
1:B:144:ILE:HD13	1:B:145:GLN:OE1	2.07	0.54
1:A:141:VAL:HG12	1:A:141:VAL:O	2.07	0.53
1:A:201:LYS:HD2	1:A:239:SER:OG	2.09	0.53
1:A:90:LEU:HD21	1:A:146:ALA:HB2	1.90	0.53
1:A:172:THR:HA	1:A:244:VAL:O	2.09	0.53
1:B:216:THR:O	1:B:218:ARG:N	2.42	0.52
1:A:215:ILE:O	1:A:217:THR:N	2.43	0.52
1:B:161:ARG:NH1	1:B:161:ARG:HB3	2.25	0.51
1:A:223:THR:CG2	1:A:226:GLN:HG3	2.41	0.50
1:B:95:ALA:HB1	1:B:100:ASP:OD1	2.12	0.50
1:B:100:ASP:CB	1:B:132:TYR:HB2	2.41	0.50
1:A:223:THR:HG22	1:A:226:GLN:N	2.18	0.50
1:A:218:ARG:NH2	1:B:100:ASP:OD2	2.38	0.50
1:B:95:ALA:HB1	1:B:100:ASP:O	2.12	0.49
1:A:147:GLU:HG2	1:A:149:TRP:CE2	2.47	0.49
1:B:224:LEU:O	1:B:228:VAL:HG23	2.12	0.49
1:A:95:ALA:CB	1:A:102:SER:HB2	2.42	0.49
1:A:223:THR:CG2	1:A:225:GLN:N	2.69	0.49
1:B:181:THR:HB	1:B:184:ALA:HB3	1.94	0.49
1:A:97:THR:OG1	1:A:100:ASP:OD1	2.30	0.48
1:A:156:ARG:O	1:A:160:GLU:HG3	2.13	0.48
1:B:154:LEU:HD11	1:B:158:ASP:CB	2.42	0.48
1:A:96:ARG:HG2	1:A:96:ARG:NH1	2.28	0.48
1:A:193:ASP:HB2	1:A:199:HIS:CE1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:LEU:HD22	1:B:172:THR:HG22	1.95	0.48
1:B:163:LEU:HD21	1:B:174:LEU:HD23	1.96	0.48
1:B:191:ASP:OD1	1:B:240:SER:HB2	2.14	0.48
1:B:156:ARG:NH1	1:B:160:GLU:OE2	2.47	0.47
1:B:104:HIS:NE2	1:B:107:GLU:HB2	2.29	0.47
1:B:172:THR:HA	1:B:244:VAL:O	2.15	0.47
1:B:85:THR:HG22	1:B:86:LEU:N	2.29	0.47
1:A:223:THR:CG2	1:A:225:GLN:HB2	2.45	0.46
1:B:112:LEU:HB2	1:B:121:GLU:HG3	1.97	0.46
1:B:113:ASN:ND2	1:B:116:GLU:OE2	2.48	0.46
1:B:161:ARG:HB3	1:B:161:ARG:HH11	1.80	0.46
1:A:216:THR:O	1:A:216:THR:CG2	2.62	0.46
1:B:227:LEU:C	1:B:227:LEU:HD23	2.41	0.46
1:A:116:GLU:OE2	1:A:132:TYR:OH	2.30	0.45
1:A:227:LEU:C	1:A:227:LEU:HD23	2.41	0.45
1:A:159:ALA:HB1	1:A:174:LEU:HD21	1.98	0.45
1:B:216:THR:O	1:B:217:THR:C	2.60	0.45
1:B:190:ARG:NH1	1:B:198:ASP:OD2	2.50	0.44
1:A:113:ASN:HB3	1:A:121:GLU:HB3	1.99	0.44
1:A:223:THR:HG21	1:A:225:GLN:HB2	2.00	0.44
1:B:222:GLU:HG3	2:B:14:HOH:O	2.17	0.44
1:B:234:ARG:HH11	1:B:234:ARG:HB3	1.82	0.44
1:A:244:VAL:HA	1:A:245:PRO:HD3	1.78	0.44
1:B:100:ASP:OD1	1:B:100:ASP:C	2.60	0.44
1:B:215:ILE:HG22	1:B:216:THR:HG23	1.99	0.44
1:A:142:ASP:OD1	1:A:142:ASP:N	2.50	0.44
1:B:215:ILE:O	1:B:216:THR:C	2.61	0.44
1:B:173:PHE:O	1:B:245:PRO:HA	2.18	0.43
1:A:141:VAL:O	1:A:141:VAL:CG1	2.66	0.43
1:B:174:LEU:C	1:B:174:LEU:HD12	2.42	0.43
1:A:189:ILE:O	1:A:200:VAL:HG13	2.19	0.43
1:A:170:ARG:O	1:A:243:VAL:HG21	2.19	0.43
1:A:150:TYR:C	1:A:151:PHE:HD1	2.27	0.43
1:B:113:ASN:HA	2:B:6:HOH:O	2.19	0.43
1:B:93:TYR:HB3	1:B:103:PHE:CE2	2.53	0.43
1:B:181:THR:HG22	1:B:182:LYS:N	2.33	0.43
1:B:142:ASP:CG	1:B:143:SER:N	2.77	0.42
1:B:116:GLU:H	1:B:116:GLU:HG2	1.51	0.42
1:B:93:TYR:HB3	1:B:103:PHE:CZ	2.55	0.42
1:A:158:ASP:OD1	1:A:161:ARG:NH2	2.50	0.42
1:A:107:GLU:OE2	1:A:126:THR:HB	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:LEU:O	1:B:176:ARG:HD2	2.20	0.42
1:B:140:PRO:HD2	1:B:143:SER:OG	2.20	0.42
1:B:113:ASN:CB	1:B:121:GLU:HB3	2.43	0.41
1:B:143:SER:OG	1:B:145:GLN:NE2	2.52	0.41
1:B:216:THR:C	1:B:218:ARG:N	2.77	0.41
1:B:223:THR:HG23	1:B:225:GLN:H	1.84	0.41
1:A:208:LEU:O	1:A:209:ASP:C	2.63	0.41
1:B:164:LEU:HD22	1:B:198:ASP:OD1	2.20	0.41
1:B:104:HIS:CD2	1:B:107:GLU:HB2	2.55	0.41
1:B:125:LEU:HD23	1:B:125:LEU:HA	1.90	0.41
1:A:116:GLU:H	1:A:116:GLU:HG2	1.70	0.41
1:B:142:ASP:OD1	1:B:143:SER:N	2.54	0.41
1:A:156:ARG:NH1	1:A:160:GLU:OE2	2.54	0.40
1:B:110:GLN:OE1	1:B:123:ARG:HD2	2.21	0.40
1:B:114:SER:HA	1:B:120:TRP:CD1	2.56	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	159/165 (96%)	141 (89%)	14 (9%)	4 (2%)	4	8
1	B	159/165 (96%)	147 (92%)	9 (6%)	3 (2%)	6	13
All	All	318/330 (96%)	288 (91%)	23 (7%)	7 (2%)	5	10

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	216	THR
1	B	143	SER
1	B	216	THR

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Mol	Chain	Res	Type
1	A	97	THR
1	B	217	THR
1	A	96	ARG
1	A	152	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	138/141 (98%)	129 (94%)	9 (6%)	15	35
1	B	138/141 (98%)	127 (92%)	11 (8%)	11	25
All	All	276/282 (98%)	256 (93%)	20 (7%)	13	30

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	GLU
1	A	121	GLU
1	A	130	THR
1	A	181	THR
1	A	182	LYS
1	A	189	ILE
1	A	190	ARG
1	A	223	THR
1	A	244	VAL
1	B	116	GLU
1	B	142	ASP
1	B	144	ILE
1	B	172	THR
1	B	186	SER
1	B	190	ARG
1	B	198	ASP
1	B	209	ASP
1	B	223	THR
1	B	232	SER

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Mol	Chain	Res	Type
1	B	234	ARG

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	104	HIS
1	A	110	GLN
1	B	104	HIS
1	B	199	HIS

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 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	161/165 (97%)	0.26	11 (6%) 23 19	9, 29, 54, 65	0
1	B	161/165 (97%)	0.49	6 (3%) 45 39	15, 37, 63, 67	0
All	All	322/330 (97%)	0.38	17 (5%) 32 26	9, 34, 60, 67	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	217	THR	8.2
1	A	216	THR	7.7
1	A	97	THR	6.5
1	B	216	THR	6.2
1	B	97	THR	5.6
1	B	217	THR	5.6
1	B	197	GLY	3.9
1	B	198	ASP	3.1
1	A	96	ARG	3.1
1	A	98	GLU	3.1
1	B	96	ARG	2.8
1	A	193	ASP	2.7
1	A	94	GLU	2.5
1	A	218	ARG	2.5
1	A	195	MET	2.3
1	A	197	GLY	2.3
1	A	95	ALA	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.