Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 08:29 AM BST

PDB ID	:	6IAK
Title	:	The crystal structure of the chicken CREB3 bZIP
Authors	:	Sabaratnam, K.; Renner, M.
Deposited on	:	2018-11-26
Resolution	:	3.95 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 following versions of software and data (see references ⓘ) were used in the production of this report:

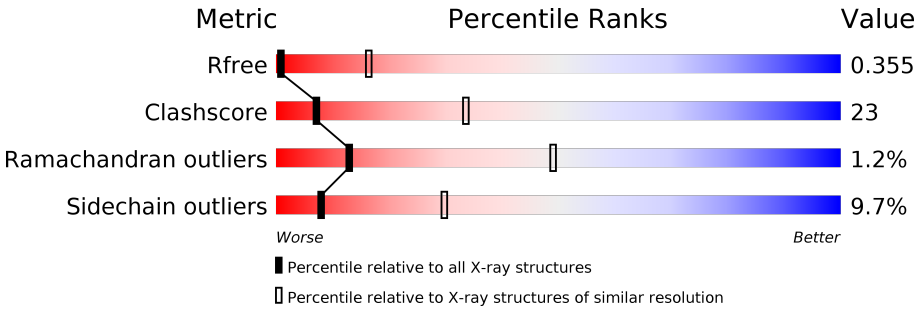
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

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 3.95 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	1025 (4.22-3.70)
Clashscore	141614	1085 (4.22-3.70)
Ramachandran outliers	138981	1047 (4.22-3.70)
Sidechain outliers	138945	1039 (4.22-3.70)

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 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 <=5%

Mol	Chain	Length	Quality of chain
1	A	441	7% 7% • 85%
1	B	441	7% 6% • 86%
1	C	441	7% 7% • 86%
1	D	441	7% 6% 86%
1	E	441	9% 6% 85%
1	F	441	9% 6% • 84%
1	G	441	6% • 92%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 3563 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 Uncharacterized protein.

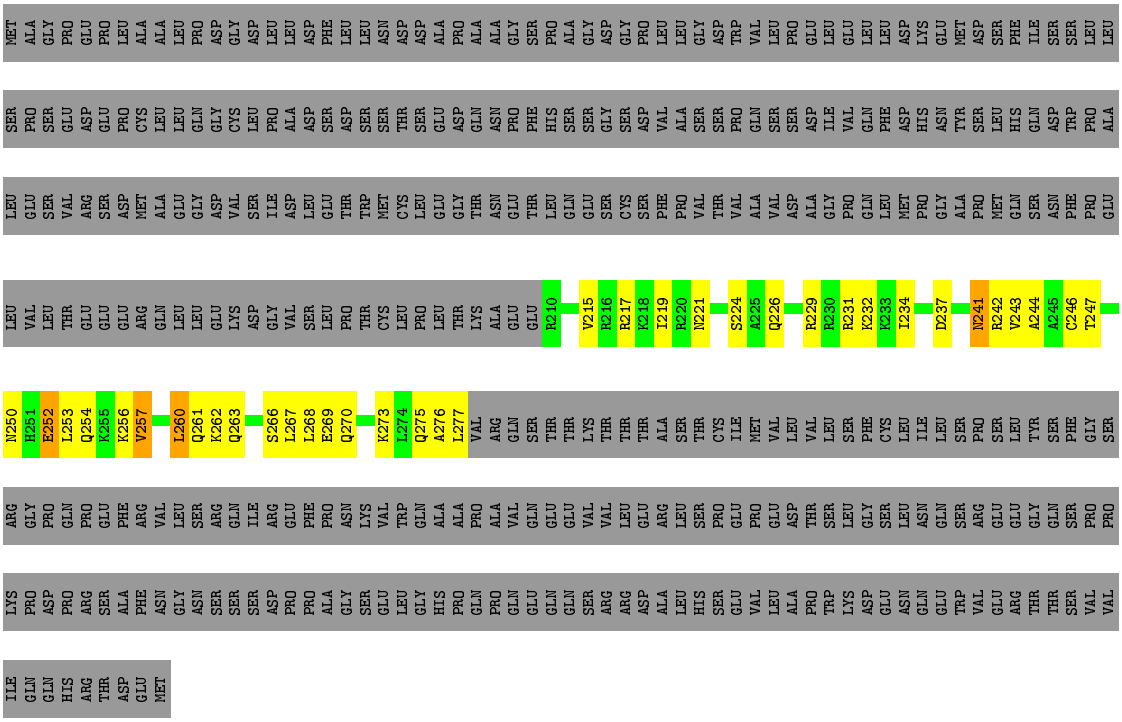
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	68	Total	C	N	O	S	0	0	0
			567	349	120	96	2			
1	B	62	Total	C	N	O	S	0	0	0
			511	313	106	90	2			
1	C	61	Total	C	N	O	S	0	0	0
			509	310	108	89	2			
1	D	62	Total	C	N	O	S	0	0	0
			518	317	110	89	2			
1	E	67	Total	C	N	O	S	0	0	0
			563	345	119	97	2			
1	F	71	Total	C	N	O	S	0	0	0
			593	363	124	104	2			
1	G	37	Total	C	N	O	S	0	0	0
			302	180	65	56	1			

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. 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. 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.

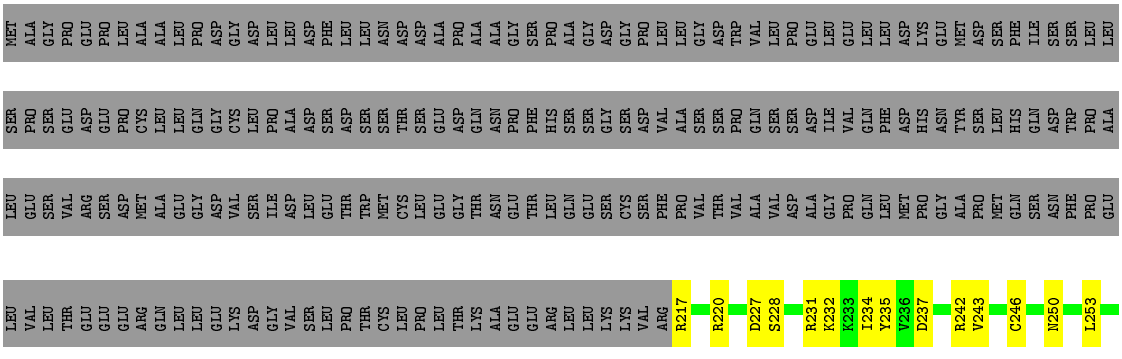
- Molecule 1: Uncharacterized protein

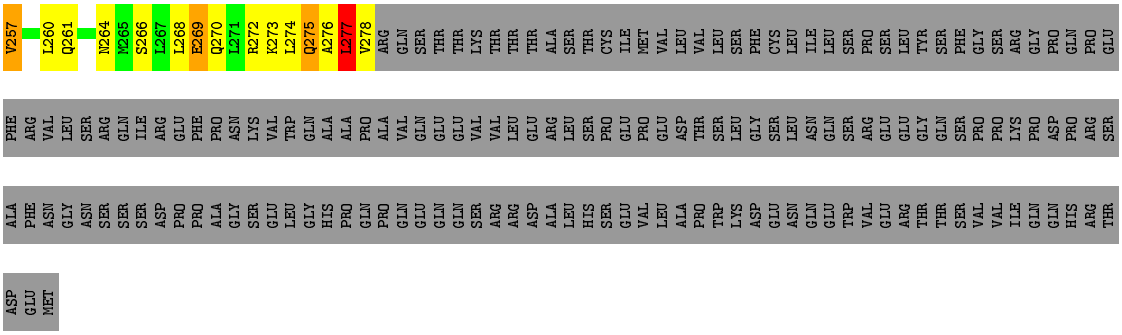
Chain A:



- Molecule 1: Uncharacterized protein

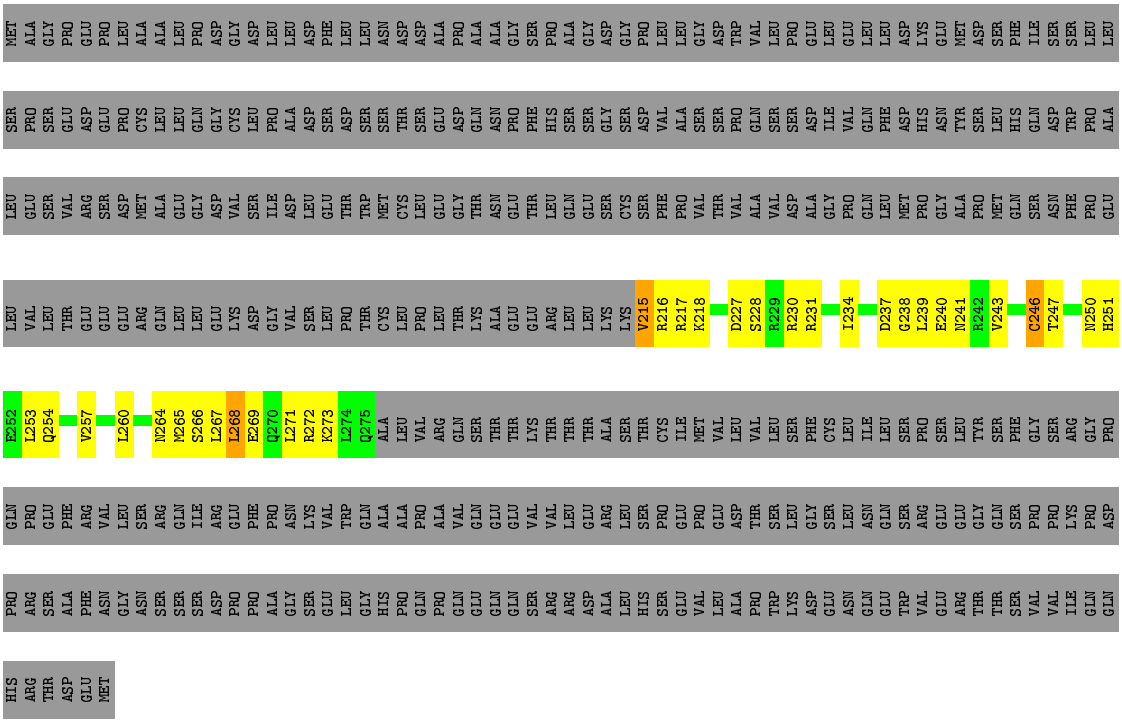
Chain B:





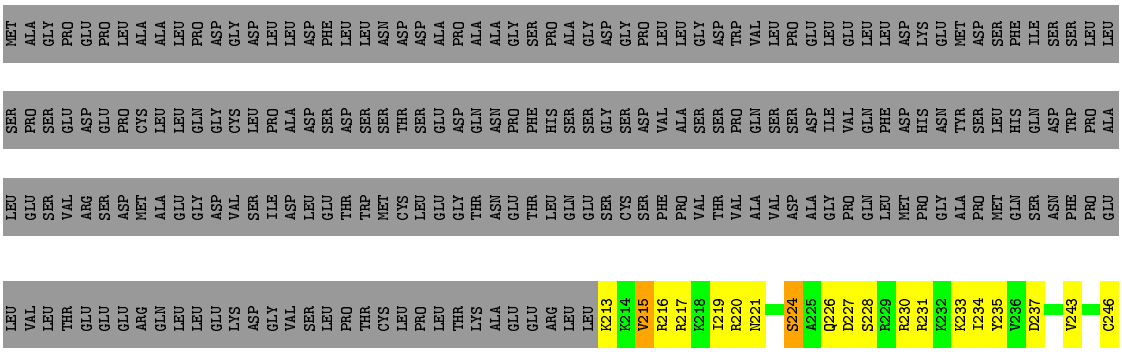
- Molecule 1: Uncharacterized protein

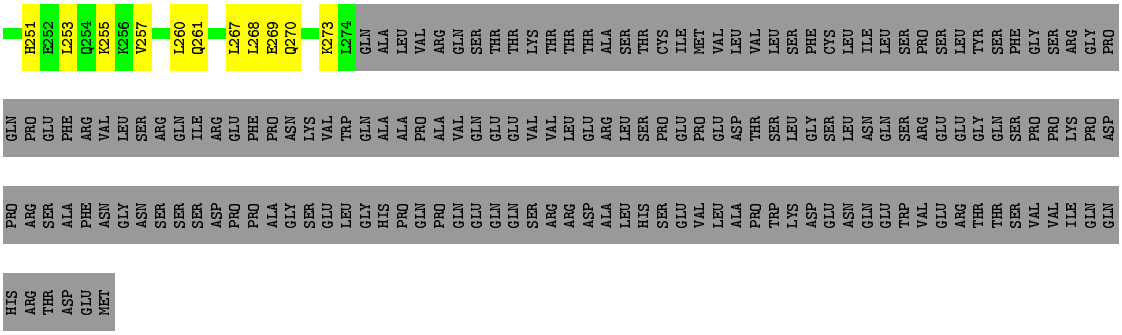
Chain C: 7% 7% 86%



- Molecule 1: Uncharacterized protein

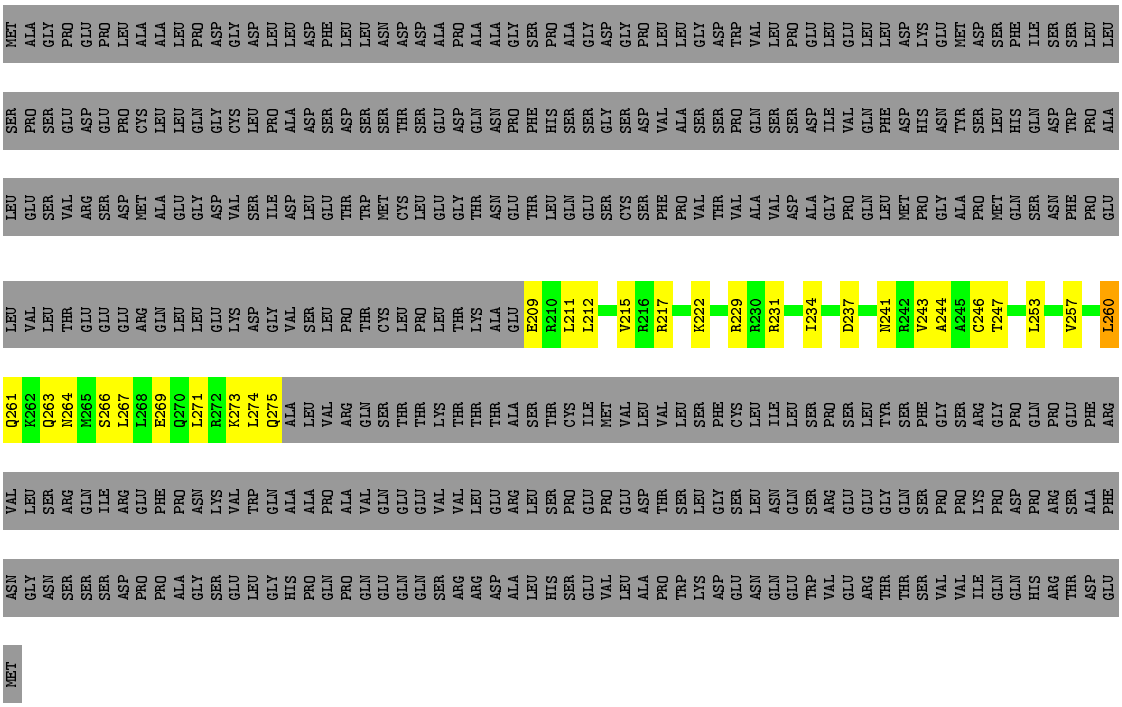
Chain D: 7% 6% 86%





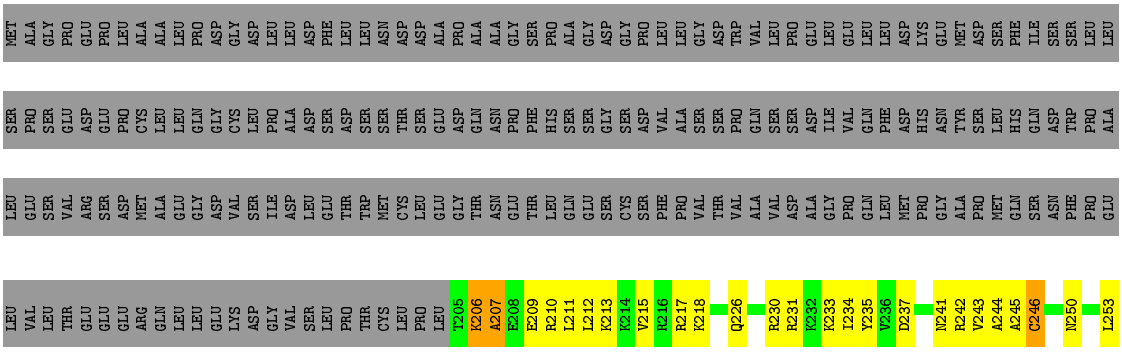
- Molecule 1: Uncharacterized protein

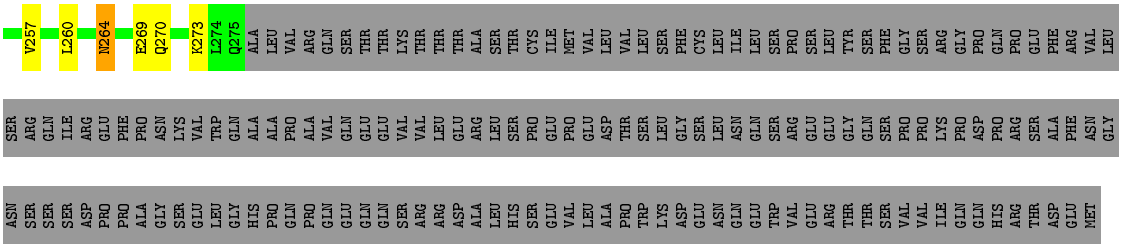
Chain E:



- Molecule 1: Uncharacterized protein

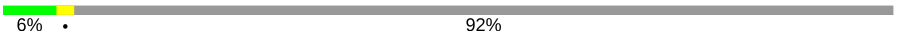
Chain F:

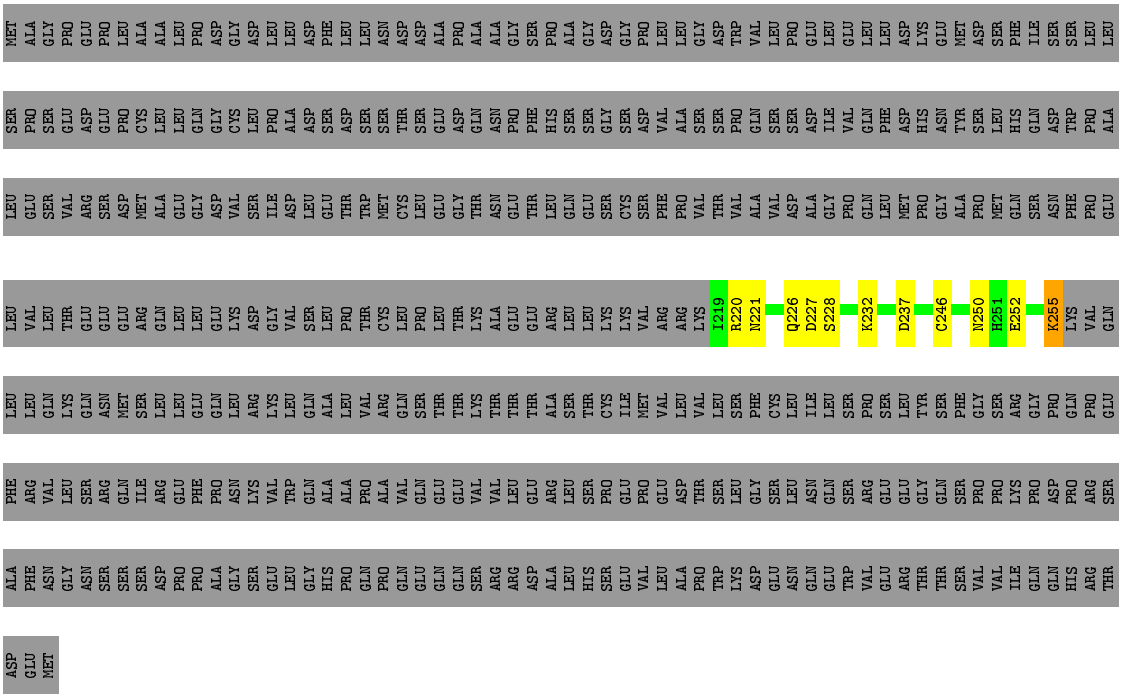




• Molecule 1: Uncharacterized protein

Chain G:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	137.90Å 167.06Å 115.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.64 – 3.95 51.64 – 3.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (51.64-3.95) 99.9 (51.64-3.95)	Depositor EDS
R_{merge}	0.30	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.92 (at 4.00Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.261 , 0.300 0.312 , 0.355	Depositor DCC
R_{free} test set	605 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	126.6	Xtriage
Anisotropy	0.469	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 168.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	3563	wwPDB-VP
Average B, all atoms (Å^2)	158.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.96% 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, 0.333 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	0.74	0/568	1.10	3/751 (0.4%)
1	B	0.76	0/512	1.07	2/679 (0.3%)
1	C	0.78	0/510	1.07	2/675 (0.3%)
1	D	0.81	0/519	0.99	0/685
1	E	0.75	0/564	1.05	3/745 (0.4%)
1	F	0.77	0/594	1.03	1/785 (0.1%)
1	G	0.70	0/303	1.07	1/402 (0.2%)
All	All	0.76	0/3570	1.06	12/4722 (0.3%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	276	ALA	N-CA-CB	-6.40	101.14	110.10
1	A	252	GLU	N-CA-CB	6.15	121.67	110.60
1	F	207	ALA	C-N-CA	5.93	136.52	121.70
1	E	212	LEU	N-CA-C	-5.62	95.81	111.00
1	E	266	SER	N-CA-CB	-5.54	102.19	110.50
1	C	268	LEU	CB-CA-C	-5.51	99.73	110.20
1	G	255	LYS	N-CA-CB	5.43	120.38	110.60
1	A	276	ALA	C-N-CA	5.38	135.16	121.70
1	B	277	LEU	CA-CB-CG	5.27	127.42	115.30
1	C	215	VAL	C-N-CA	5.19	134.67	121.70
1	E	217	ARG	N-CA-CB	5.10	119.78	110.60
1	A	241	ASN	CA-CB-CG	-5.09	102.19	113.40

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	567	0	626	45	0
1	B	511	0	552	34	0
1	C	509	0	550	40	0
1	D	518	0	568	38	0
1	E	563	0	616	18	0
1	F	593	0	647	38	0
1	G	302	0	309	3	0
All	All	3563	0	3868	169	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 (169) 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:F:211:LEU:CD1	1:F:212:LEU:HD12	1.12	1.56
1:F:211:LEU:CD1	1:F:212:LEU:CD1	1.74	1.55
1:F:211:LEU:HD11	1:F:212:LEU:CD1	1.34	1.39
1:C:253:LEU:HG	1:D:253:LEU:HD22	1.28	1.15
1:F:211:LEU:HD11	1:F:212:LEU:HD13	1.32	1.12
1:A:253:LEU:HG	1:B:253:LEU:HD22	1.14	1.11
1:F:211:LEU:HD12	1:F:212:LEU:CD1	1.56	1.10
1:F:211:LEU:HD11	1:F:212:LEU:HD11	1.42	1.01
1:A:242:ARG:NH2	1:D:237:ASP:OD2	1.92	1.01
1:B:272:ARG:HA	1:B:275:GLN:HB3	1.43	0.98
1:C:268:LEU:O	1:C:268:LEU:HD23	1.64	0.96
1:D:217:ARG:O	1:D:220:ARG:HG2	1.65	0.95
1:A:215:VAL:O	1:A:219:ILE:CD1	2.18	0.91
1:F:211:LEU:HD12	1:F:212:LEU:N	1.86	0.90
1:F:211:LEU:CD1	1:F:212:LEU:HD13	1.94	0.86
1:F:206:LYS:NZ	1:F:210:ARG:HG3	1.90	0.86
1:E:211:LEU:O	1:E:211:LEU:HD23	1.76	0.85
1:C:267:LEU:HD13	1:D:268:LEU:HD21	1.59	0.85
1:A:253:LEU:CG	1:B:253:LEU:HD22	2.06	0.80
1:F:226:GLN:HE22	1:F:230:ARG:HD3	1.46	0.80

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:267:LEU:HD13	1:D:268:LEU:CD2	2.12	0.79
1:D:217:ARG:HH11	1:D:220:ARG:NH2	1.79	0.79
1:C:267:LEU:CD1	1:D:268:LEU:HD21	2.12	0.79
1:A:254:GLN:HG3	1:B:253:LEU:HD11	1.66	0.77
1:C:257:VAL:HG22	1:D:257:VAL:HG22	1.66	0.77
1:A:215:VAL:O	1:A:219:ILE:HD11	1.83	0.77
1:D:213:LYS:NZ	1:D:217:ARG:HD3	1.99	0.77
1:F:211:LEU:HD13	1:F:212:LEU:HD12	1.59	0.74
1:F:211:LEU:CD1	1:F:212:LEU:HD11	2.03	0.74
1:A:269:GLU:O	1:A:273:LYS:HB2	1.89	0.73
1:D:213:LYS:HZ3	1:D:217:ARG:HD3	1.52	0.72
1:C:266:SER:HA	1:C:269:GLU:HG2	1.72	0.71
1:C:237:ASP:O	1:C:241:ASN:ND2	2.24	0.70
1:A:268:LEU:O	1:A:268:LEU:HD23	1.92	0.70
1:B:268:LEU:HD23	1:B:268:LEU:O	1.92	0.69
1:E:260:LEU:HD13	1:F:260:LEU:HB3	1.74	0.69
1:B:217:ARG:HB3	1:B:220:ARG:HG2	1.72	0.69
1:A:257:VAL:HG22	1:B:257:VAL:HG22	1.75	0.68
1:F:211:LEU:HD12	1:F:212:LEU:HD12	0.68	0.68
1:A:256:LYS:HE3	1:A:260:LEU:HG	1.77	0.67
1:C:253:LEU:CG	1:D:253:LEU:HD22	2.17	0.67
1:C:268:LEU:C	1:C:268:LEU:HD23	2.14	0.66
1:C:269:GLU:O	1:C:273:LYS:HG3	1.96	0.66
1:A:226:GLN:HG2	1:A:229:ARG:NH2	2.10	0.66
1:C:260:LEU:HD13	1:D:260:LEU:HB3	1.77	0.66
1:G:220:ARG:NH2	1:G:221:ASN:HD21	1.94	0.65
1:A:243:VAL:HG23	1:B:243:VAL:HG23	1.77	0.65
1:F:231:ARG:HA	1:F:234:ILE:HD12	1.79	0.64
1:D:217:ARG:NH1	1:D:220:ARG:NH2	2.44	0.64
1:E:211:LEU:HG	1:E:215:VAL:HG13	1.78	0.63
1:C:231:ARG:HA	1:C:234:ILE:HD12	1.80	0.63
1:A:231:ARG:HA	1:A:234:ILE:HD12	1.81	0.62
1:B:231:ARG:HA	1:B:234:ILE:HD12	1.81	0.62
1:B:266:SER:O	1:B:270:GLN:HG2	2.00	0.62
1:C:268:LEU:CD2	1:C:272:ARG:HG2	2.28	0.62
1:C:266:SER:HA	1:C:269:GLU:OE2	1.99	0.61
1:E:243:VAL:HG23	1:F:243:VAL:HG23	1.82	0.61
1:E:231:ARG:HA	1:E:234:ILE:HD12	1.82	0.61
1:E:253:LEU:HG	1:F:253:LEU:HG	1.82	0.60
1:A:266:SER:O	1:A:270:GLN:HG2	2.02	0.59
1:B:269:GLU:O	1:B:273:LYS:HB2	2.02	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:217:ARG:HD2	1:D:220:ARG:CZ	2.34	0.58
1:A:215:VAL:O	1:A:219:ILE:HD12	2.03	0.58
1:A:260:LEU:HD13	1:B:260:LEU:HB3	1.85	0.58
1:F:269:GLU:O	1:F:273:LYS:HB2	2.03	0.58
1:C:243:VAL:HG23	1:D:243:VAL:HG12	1.86	0.57
1:D:231:ARG:HA	1:D:234:ILE:HD12	1.87	0.57
1:E:269:GLU:O	1:E:273:LYS:HB2	2.05	0.57
1:D:269:GLU:O	1:D:273:LYS:HB2	2.05	0.57
1:F:211:LEU:O	1:F:215:VAL:HG23	2.05	0.57
1:F:206:LYS:HZ1	1:F:210:ARG:HG3	1.69	0.56
1:A:275:GLN:H	1:A:277:LEU:HD13	1.71	0.56
1:E:260:LEU:HB3	1:F:260:LEU:HD13	1.87	0.56
1:B:272:ARG:CA	1:B:275:GLN:HB3	2.28	0.56
1:C:215:VAL:HG22	1:C:218:LYS:HB2	1.87	0.56
1:A:260:LEU:HB3	1:B:260:LEU:HD13	1.88	0.55
1:C:271:LEU:HD11	1:D:270:GLN:CD	2.26	0.55
1:C:266:SER:HA	1:C:269:GLU:CG	2.37	0.54
1:E:257:VAL:HG22	1:F:257:VAL:HG22	1.89	0.54
1:B:268:LEU:HD21	1:B:272:ARG:HD3	1.90	0.54
1:A:217:ARG:O	1:A:221:ASN:OD1	2.26	0.54
1:A:250:ASN:CG	1:B:250:ASN:HD22	2.12	0.54
1:A:250:ASN:OD1	1:B:250:ASN:HA	2.09	0.52
1:B:268:LEU:HD23	1:B:268:LEU:C	2.30	0.52
1:B:274:LEU:O	1:B:277:LEU:HB3	2.10	0.52
1:D:226:GLN:HE22	1:D:230:ARG:HD3	1.75	0.52
1:C:268:LEU:CD2	1:C:268:LEU:O	2.49	0.51
1:F:206:LYS:HZ2	1:F:210:ARG:HG3	1.72	0.51
1:E:222:LYS:HE3	1:E:229:ARG:HH22	1.74	0.51
1:A:250:ASN:O	1:A:253:LEU:HB3	2.11	0.51
1:C:254:GLN:HG3	1:D:253:LEU:HD11	1.92	0.51
1:B:231:ARG:HD3	1:F:235:TYR:OH	2.11	0.51
1:A:253:LEU:HD23	1:B:253:LEU:HD13	1.93	0.51
1:B:273:LYS:O	1:B:277:LEU:HB2	2.12	0.50
1:B:235:TYR:OH	1:F:231:ARG:HD3	2.12	0.50
1:C:268:LEU:HD23	1:C:272:ARG:HG2	1.92	0.50
1:E:257:VAL:HG12	1:E:261:GLN:HE21	1.77	0.50
1:B:257:VAL:HG12	1:B:261:GLN:HE21	1.75	0.50
1:C:238:GLY:HA2	1:C:241:ASN:HD21	1.76	0.50
1:C:246:CYS:SG	1:C:250:ASN:ND2	2.85	0.49
1:D:213:LYS:HZ2	1:D:217:ARG:HD3	1.78	0.49
1:F:209:GLU:O	1:F:213:LYS:HB2	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:ASP:OD2	1:F:242:ARG:NH2	2.46	0.48
1:A:232:LYS:HD2	1:B:232:LYS:HE3	1.95	0.48
1:D:220:ARG:HG3	1:D:221:ASN:N	2.28	0.48
1:F:211:LEU:HD12	1:F:212:LEU:CA	2.44	0.47
1:A:232:LYS:NZ	1:B:232:LYS:CE	2.78	0.47
1:A:277:LEU:HD12	1:A:277:LEU:H	1.80	0.47
1:C:260:LEU:HB3	1:D:260:LEU:HD13	1.97	0.47
1:C:266:SER:HA	1:C:269:GLU:CD	2.35	0.47
1:A:215:VAL:O	1:A:215:VAL:HG12	2.14	0.46
1:F:213:LYS:O	1:F:217:ARG:HG3	2.14	0.46
1:E:211:LEU:CD2	1:E:211:LEU:O	2.56	0.46
1:G:220:ARG:HG3	1:G:221:ASN:ND2	2.31	0.46
1:F:207:ALA:O	1:F:211:LEU:HG	2.16	0.46
1:C:266:SER:CA	1:C:269:GLU:HG2	2.43	0.46
1:A:232:LYS:CD	1:B:232:LYS:HE3	2.46	0.45
1:A:215:VAL:O	1:A:219:ILE:CG1	2.65	0.45
1:D:215:VAL:O	1:D:219:ILE:HG12	2.17	0.45
1:E:243:VAL:O	1:E:247:THR:HB	2.17	0.45
1:C:268:LEU:HD21	1:C:272:ARG:CG	2.46	0.45
1:C:271:LEU:HD11	1:D:270:GLN:OE1	2.16	0.45
1:F:241:ASN:O	1:F:244:ALA:HB3	2.17	0.45
1:A:232:LYS:HZ3	1:B:232:LYS:CE	2.30	0.45
1:D:217:ARG:HD2	1:D:220:ARG:NE	2.31	0.45
1:F:226:GLN:NE2	1:F:230:ARG:HD3	2.24	0.45
1:E:271:LEU:HD11	1:F:270:GLN:HE21	1.82	0.45
1:A:257:VAL:HG12	1:A:261:GLN:HE21	1.81	0.45
1:B:260:LEU:HD23	1:B:260:LEU:HA	1.84	0.44
1:D:255:LYS:HA	1:D:255:LYS:HD2	1.74	0.44
1:A:231:ARG:HD3	1:D:235:TYR:OH	2.17	0.44
1:A:242:ARG:HH12	1:D:237:ASP:HB3	1.83	0.44
1:F:215:VAL:HA	1:F:218:LYS:HE2	1.99	0.44
1:A:268:LEU:HD23	1:A:268:LEU:C	2.38	0.44
1:A:232:LYS:HZ3	1:B:232:LYS:HE2	1.83	0.44
1:E:274:LEU:O	1:E:275:GLN:HG3	2.18	0.44
1:C:237:ASP:HA	1:C:240:GLU:HG3	1.99	0.43
1:D:226:GLN:NE2	1:D:230:ARG:HB2	2.33	0.43
1:A:241:ASN:O	1:A:244:ALA:HB3	2.17	0.43
1:C:268:LEU:HD21	1:C:272:ARG:HG2	1.97	0.43
1:D:243:VAL:HA	1:D:246:CYS:SG	2.58	0.43
1:C:247:THR:O	1:C:251:HIS:N	2.41	0.43
1:C:268:LEU:O	1:C:272:ARG:HG2	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:GLN:O	1:A:267:LEU:HD13	2.19	0.43
1:B:242:ARG:HH22	1:F:233:LYS:HE2	1.84	0.42
1:A:232:LYS:NZ	1:B:232:LYS:HE2	2.34	0.42
1:C:260:LEU:HD11	1:D:261:GLN:HG3	2.01	0.42
1:C:271:LEU:C	1:C:271:LEU:HD23	2.40	0.42
1:A:262:LYS:O	1:A:262:LYS:HG3	2.19	0.42
1:A:242:ARG:HH12	1:D:237:ASP:CB	2.32	0.42
1:G:227:ASP:O	1:G:228:SER:C	2.57	0.42
1:A:243:VAL:O	1:A:247:THR:HB	2.20	0.42
1:D:227:ASP:O	1:D:228:SER:C	2.58	0.42
1:C:269:GLU:O	1:C:273:LYS:CG	2.67	0.41
1:C:227:ASP:O	1:C:228:SER:C	2.59	0.41
1:E:222:LYS:HA	1:E:222:LYS:HD2	1.90	0.41
1:A:242:ARG:HH21	1:D:233:LYS:HD3	1.84	0.41
1:A:242:ARG:HH22	1:D:237:ASP:CG	2.11	0.41
1:C:268:LEU:CD2	1:C:272:ARG:CG	2.96	0.41
1:D:220:ARG:O	1:D:224:SER:HB3	2.21	0.41
1:C:253:LEU:HD23	1:D:253:LEU:HD13	2.02	0.41
1:E:241:ASN:O	1:E:244:ALA:HB3	2.21	0.41
1:F:245:ALA:O	1:F:246:CYS:C	2.59	0.41
1:F:206:LYS:HG2	1:F:209:GLU:HB3	2.02	0.41
1:A:215:VAL:O	1:A:219:ILE:HG13	2.21	0.41
1:B:227:ASP:O	1:B:228:SER:C	2.59	0.41
1:C:239:LEU:HA	1:C:239:LEU:HD23	1.87	0.40
1:E:263:GLN:HE21	1:F:264:ASN:ND2	2.18	0.40
1:A:215:VAL:O	1:A:215:VAL:CG1	2.70	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	66/441 (15%)	59 (89%)	7 (11%)	0	100 100
1	B	60/441 (14%)	55 (92%)	3 (5%)	2 (3%)	4 30
1	C	59/441 (13%)	57 (97%)	1 (2%)	1 (2%)	9 42
1	D	60/441 (14%)	56 (93%)	3 (5%)	1 (2%)	9 42
1	E	65/441 (15%)	55 (85%)	10 (15%)	0	100 100
1	F	69/441 (16%)	62 (90%)	6 (9%)	1 (1%)	11 45
1	G	35/441 (8%)	31 (89%)	4 (11%)	0	100 100
All	All	414/3087 (13%)	375 (91%)	34 (8%)	5 (1%)	13 49

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	206	LYS
1	B	275	GLN
1	C	216	ARG
1	B	277	LEU
1	D	215	VAL

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	62/392 (16%)	56 (90%)	6 (10%)	8 30
1	B	56/392 (14%)	50 (89%)	6 (11%)	6 27
1	C	56/392 (14%)	51 (91%)	5 (9%)	9 35
1	D	57/392 (14%)	53 (93%)	4 (7%)	15 43
1	E	62/392 (16%)	56 (90%)	6 (10%)	8 30
1	F	65/392 (17%)	61 (94%)	4 (6%)	18 46
1	G	32/392 (8%)	25 (78%)	7 (22%)	1 6
All	All	390/2744 (14%)	352 (90%)	38 (10%)	8 30

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	224	SER
1	A	237	ASP
1	A	246	CYS
1	A	252	GLU
1	A	257	VAL
1	A	260	LEU
1	B	246	CYS
1	B	257	VAL
1	B	264	ASN
1	B	269	GLU
1	B	277	LEU
1	B	278	VAL
1	C	217	ARG
1	C	230	ARG
1	C	246	CYS
1	C	264	ASN
1	C	265	MET
1	D	216	ARG
1	D	224	SER
1	D	251	HIS
1	D	267	LEU
1	E	209	GLU
1	E	237	ASP
1	E	246	CYS
1	E	260	LEU
1	E	264	ASN
1	E	267	LEU
1	F	237	ASP
1	F	246	CYS
1	F	250	ASN
1	F	264	ASN
1	G	226	GLN
1	G	232	LYS
1	G	237	ASP
1	G	246	CYS
1	G	250	ASN
1	G	252	GLU
1	G	255	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	241	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	251	HIS
1	A	261	GLN
1	A	264	ASN
1	B	226	GLN
1	B	250	ASN
1	B	251	HIS
1	B	261	GLN
1	B	275	GLN
1	C	226	GLN
1	C	241	ASN
1	C	249	GLN
1	C	250	ASN
1	D	226	GLN
1	D	264	ASN
1	E	226	GLN
1	E	258	GLN
1	E	261	GLN
1	E	263	GLN
1	E	270	GLN
1	F	226	GLN
1	F	258	GLN
1	F	270	GLN
1	F	275	GLN
1	G	221	ASN

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 monosaccharides 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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.