



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 03:16 PM GMT

PDB ID : 2EFS
Title : Crystal structure of the C-terminal tropomyosin fragment with N- and C-terminal extensions of the leucine zipper at 2.0 angstroms resolution
Authors : Minakata, S.; Nitandai, Y.; Maeda, K.; Oda, N.; Wakabayashi, K.; Maeda, Y.
Deposited on : 2007-02-23
Resolution : 2.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

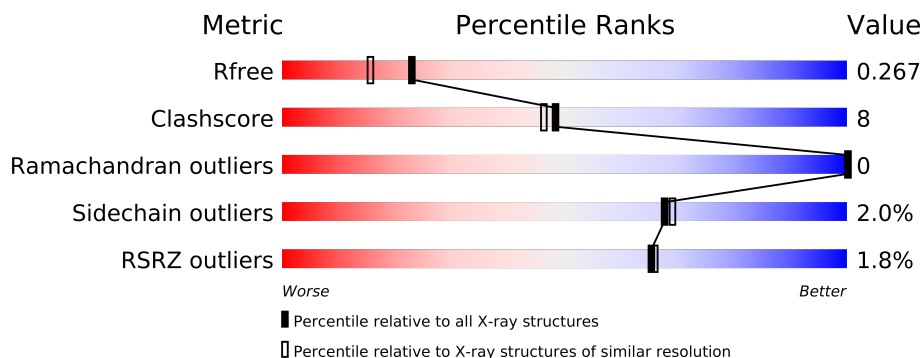
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.00 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}	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	155	
1	B	155	
1	C	155	
1	D	155	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5457 atoms, of which 0 are hydrogen and 0 are deuterium.

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 General control protein GCN4 and Tropomyosin 1 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	154	Total	C	N	O	S	0	0	0
			1274	793	211	268	2			
1	B	154	Total	C	N	O	S	0	0	0
			1274	793	211	268	2			
1	C	154	Total	C	N	O	S	0	0	0
			1274	793	211	268	2			
1	D	154	Total	C	N	O	S	0	0	0
			1274	793	211	268	2			

There are 116 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	190	SER	CYS	ENGINEERED	UNP P58772
A	274	MET	-	EXPRESSION TAG	UNP P58772
A	275	LYS	-	EXPRESSION TAG	UNP P58772
A	276	GLN	-	EXPRESSION TAG	UNP P58772
A	277	LEU	-	EXPRESSION TAG	UNP P58772
A	278	GLU	-	EXPRESSION TAG	UNP P58772
A	279	ASP	-	EXPRESSION TAG	UNP P58772
A	280	LYS	-	EXPRESSION TAG	UNP P58772
A	281	VAL	-	EXPRESSION TAG	UNP P58772
A	282	GLU	-	EXPRESSION TAG	UNP P58772
A	283	GLU	-	EXPRESSION TAG	UNP P58772
A	284	LEU	-	EXPRESSION TAG	UNP P58772
A	285	LEU	-	EXPRESSION TAG	UNP P58772
A	286	SER	-	EXPRESSION TAG	UNP P58772
A	287	LYS	-	EXPRESSION TAG	UNP P58772
A	288	ASN	-	EXPRESSION TAG	UNP P58772
A	289	TYR	-	EXPRESSION TAG	UNP P58772
A	290	HIS	-	EXPRESSION TAG	UNP P58772
A	291	LEU	-	EXPRESSION TAG	UNP P58772
A	292	GLU	-	EXPRESSION TAG	UNP P58772
A	293	ASN	-	EXPRESSION TAG	UNP P58772

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Chain	Residue	Modelled	Actual	Comment	Reference
A	294	GLU	-	EXPRESSION TAG	UNP P58772
A	295	VAL	-	EXPRESSION TAG	UNP P58772
A	296	ALA	-	EXPRESSION TAG	UNP P58772
A	297	ARG	-	EXPRESSION TAG	UNP P58772
A	298	LEU	-	EXPRESSION TAG	UNP P58772
A	299	LYS	-	EXPRESSION TAG	UNP P58772
A	300	LYS	-	EXPRESSION TAG	UNP P58772
A	301	LEU	-	EXPRESSION TAG	UNP P58772
B	190	SER	CYS	ENGINEERED	UNP P58772
B	274	MET	-	EXPRESSION TAG	UNP P58772
B	275	LYS	-	EXPRESSION TAG	UNP P58772
B	276	GLN	-	EXPRESSION TAG	UNP P58772
B	277	LEU	-	EXPRESSION TAG	UNP P58772
B	278	GLU	-	EXPRESSION TAG	UNP P58772
B	279	ASP	-	EXPRESSION TAG	UNP P58772
B	280	LYS	-	EXPRESSION TAG	UNP P58772
B	281	VAL	-	EXPRESSION TAG	UNP P58772
B	282	GLU	-	EXPRESSION TAG	UNP P58772
B	283	GLU	-	EXPRESSION TAG	UNP P58772
B	284	LEU	-	EXPRESSION TAG	UNP P58772
B	285	LEU	-	EXPRESSION TAG	UNP P58772
B	286	SER	-	EXPRESSION TAG	UNP P58772
B	287	LYS	-	EXPRESSION TAG	UNP P58772
B	288	ASN	-	EXPRESSION TAG	UNP P58772
B	289	TYR	-	EXPRESSION TAG	UNP P58772
B	290	HIS	-	EXPRESSION TAG	UNP P58772
B	291	LEU	-	EXPRESSION TAG	UNP P58772
B	292	GLU	-	EXPRESSION TAG	UNP P58772
B	293	ASN	-	EXPRESSION TAG	UNP P58772
B	294	GLU	-	EXPRESSION TAG	UNP P58772
B	295	VAL	-	EXPRESSION TAG	UNP P58772
B	296	ALA	-	EXPRESSION TAG	UNP P58772
B	297	ARG	-	EXPRESSION TAG	UNP P58772
B	298	LEU	-	EXPRESSION TAG	UNP P58772
B	299	LYS	-	EXPRESSION TAG	UNP P58772
B	300	LYS	-	EXPRESSION TAG	UNP P58772
B	301	LEU	-	EXPRESSION TAG	UNP P58772
C	190	SER	CYS	ENGINEERED	UNP P58772
C	274	MET	-	EXPRESSION TAG	UNP P58772
C	275	LYS	-	EXPRESSION TAG	UNP P58772
C	276	GLN	-	EXPRESSION TAG	UNP P58772
C	277	LEU	-	EXPRESSION TAG	UNP P58772

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Chain	Residue	Modelled	Actual	Comment	Reference
C	278	GLU	-	EXPRESSION TAG	UNP P58772
C	279	ASP	-	EXPRESSION TAG	UNP P58772
C	280	LYS	-	EXPRESSION TAG	UNP P58772
C	281	VAL	-	EXPRESSION TAG	UNP P58772
C	282	GLU	-	EXPRESSION TAG	UNP P58772
C	283	GLU	-	EXPRESSION TAG	UNP P58772
C	284	LEU	-	EXPRESSION TAG	UNP P58772
C	285	LEU	-	EXPRESSION TAG	UNP P58772
C	286	SER	-	EXPRESSION TAG	UNP P58772
C	287	LYS	-	EXPRESSION TAG	UNP P58772
C	288	ASN	-	EXPRESSION TAG	UNP P58772
C	289	TYR	-	EXPRESSION TAG	UNP P58772
C	290	HIS	-	EXPRESSION TAG	UNP P58772
C	291	LEU	-	EXPRESSION TAG	UNP P58772
C	292	GLU	-	EXPRESSION TAG	UNP P58772
C	293	ASN	-	EXPRESSION TAG	UNP P58772
C	294	GLU	-	EXPRESSION TAG	UNP P58772
C	295	VAL	-	EXPRESSION TAG	UNP P58772
C	296	ALA	-	EXPRESSION TAG	UNP P58772
C	297	ARG	-	EXPRESSION TAG	UNP P58772
C	298	LEU	-	EXPRESSION TAG	UNP P58772
C	299	LYS	-	EXPRESSION TAG	UNP P58772
C	300	LYS	-	EXPRESSION TAG	UNP P58772
C	301	LEU	-	EXPRESSION TAG	UNP P58772
D	190	SER	CYS	ENGINEERED	UNP P58772
D	274	MET	-	EXPRESSION TAG	UNP P58772
D	275	LYS	-	EXPRESSION TAG	UNP P58772
D	276	GLN	-	EXPRESSION TAG	UNP P58772
D	277	LEU	-	EXPRESSION TAG	UNP P58772
D	278	GLU	-	EXPRESSION TAG	UNP P58772
D	279	ASP	-	EXPRESSION TAG	UNP P58772
D	280	LYS	-	EXPRESSION TAG	UNP P58772
D	281	VAL	-	EXPRESSION TAG	UNP P58772
D	282	GLU	-	EXPRESSION TAG	UNP P58772
D	283	GLU	-	EXPRESSION TAG	UNP P58772
D	284	LEU	-	EXPRESSION TAG	UNP P58772
D	285	LEU	-	EXPRESSION TAG	UNP P58772
D	286	SER	-	EXPRESSION TAG	UNP P58772
D	287	LYS	-	EXPRESSION TAG	UNP P58772
D	288	ASN	-	EXPRESSION TAG	UNP P58772
D	289	TYR	-	EXPRESSION TAG	UNP P58772
D	290	HIS	-	EXPRESSION TAG	UNP P58772

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Chain	Residue	Modelled	Actual	Comment	Reference
D	291	LEU	-	EXPRESSION TAG	UNP P58772
D	292	GLU	-	EXPRESSION TAG	UNP P58772
D	293	ASN	-	EXPRESSION TAG	UNP P58772
D	294	GLU	-	EXPRESSION TAG	UNP P58772
D	295	VAL	-	EXPRESSION TAG	UNP P58772
D	296	ALA	-	EXPRESSION TAG	UNP P58772
D	297	ARG	-	EXPRESSION TAG	UNP P58772
D	298	LEU	-	EXPRESSION TAG	UNP P58772
D	299	LYS	-	EXPRESSION TAG	UNP P58772
D	300	LYS	-	EXPRESSION TAG	UNP P58772
D	301	LEU	-	EXPRESSION TAG	UNP P58772

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	90	Total 90	O 90	0	0
2	B	105	Total 105	O 105	0	0
2	C	71	Total 71	O 71	0	0
2	D	95	Total 95	O 95	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: General control protein GCN4 and Tropomyosin 1 alpha chain

Chain A: 



- Molecule 1: General control protein GCN4 and Tropomyosin 1 alpha chain

Chain B: 



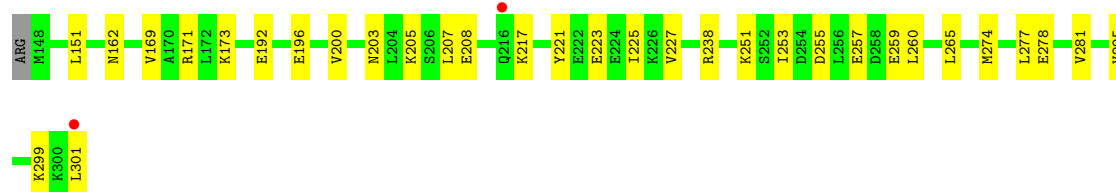
- Molecule 1: General control protein GCN4 and Tropomyosin 1 alpha chain

Chain C: 



- Molecule 1: General control protein GCN4 and Tropomyosin 1 alpha chain

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	43.20Å 63.72Å 73.26Å 66.64° 89.99° 90.30°	Depositor
Resolution (Å)	19.92 – 2.00 29.25 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.2 (19.92-2.00) 96.4 (29.25-2.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.249 , 0.304 0.233 , 0.267	Depositor DCC
R_{free} test set	2399 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	33.0	Xtriage
Anisotropy	0.203	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.3	EDS
Estimated twinning fraction	0.486 for h,-k,-l 0.000 for -h,k,k-l 0.000 for -h,-k,-k+l	Xtriage
L-test for twinning	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Outliers	0 of 47224 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5457	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.07% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.41	0/1282	0.50	0/1706
1	B	0.41	0/1282	0.51	0/1706
1	C	0.42	0/1282	0.51	0/1706
1	D	0.41	0/1282	0.51	0/1706
All	All	0.41	0/5128	0.51	0/6824

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1274	0	1293	26	0
1	B	1274	0	1293	27	0
1	C	1274	0	1293	26	0
1	D	1274	0	1293	30	0
2	A	90	0	0	4	0
2	B	105	0	0	3	0
2	C	71	0	0	1	0
2	D	95	0	0	2	0
All	All	5457	0	5172	80	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (80) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:149:LYS:HA	1:C:152:GLU:OE1	1.86	0.76
1:C:152:GLU:HG3	1:D:151:LEU:HD21	1.67	0.76
1:A:162:ASN:HB3	2:A:374:HOH:O	1.86	0.75
1:A:238:ARG:HD2	1:B:243:GLU:OE2	1.91	0.71
1:C:151:LEU:HB2	1:D:151:LEU:HD13	1.74	0.70
1:C:203:ASN:O	1:C:207:LEU:HD13	1.92	0.70
1:A:298:LEU:HD11	1:B:299:LYS:HZ2	1.58	0.67
1:A:203:ASN:O	1:A:207:LEU:HD13	1.96	0.66
1:A:148:MET:HG3	1:B:151:LEU:HD12	1.78	0.65
1:D:253:ILE:O	1:D:257:GLU:HG2	1.99	0.63
1:C:200:VAL:HG12	1:D:200:VAL:HG12	1.82	0.60
1:C:259:GLU:HG2	1:D:260:LEU:HD21	1.83	0.60
1:A:298:LEU:HD11	1:B:299:LYS:NZ	2.19	0.58
1:D:295:VAL:HG13	1:D:299:LYS:HZ2	1.68	0.58
1:D:295:VAL:HG13	1:D:299:LYS:NZ	2.18	0.58
1:C:298:LEU:HD11	1:D:299:LYS:NZ	2.18	0.58
1:A:222:GLU:HG2	2:A:383:HOH:O	2.03	0.58
1:C:153:ASP:O	1:C:157:GLU:HG3	2.04	0.58
1:B:300:LYS:HG3	1:B:300:LYS:O	2.05	0.56
1:A:214:TYR:CD1	1:B:214:TYR:HB3	2.41	0.56
1:B:204:LEU:O	1:B:208:GLU:HG3	2.06	0.55
1:D:221:TYR:O	1:D:225:ILE:HG12	2.06	0.55
1:D:203:ASN:O	1:D:207:LEU:HD13	2.05	0.55
1:B:189:LYS:HG2	2:B:369:HOH:O	2.07	0.55
1:C:228:LEU:HD21	2:C:351:HOH:O	2.07	0.54
1:A:200:VAL:HG12	1:B:200:VAL:HG12	1.88	0.53
1:A:207:LEU:HD22	1:B:204:LEU:HD12	1.91	0.53
1:D:223:GLU:O	1:D:227:VAL:HG23	2.10	0.52
1:C:162:ASN:HD21	1:D:162:ASN:HB2	1.75	0.52
1:B:253:ILE:O	1:B:257:GLU:HG2	2.10	0.52
1:D:217:LYS:HD3	1:D:221:TYR:CE2	2.45	0.51
1:A:209:ALA:O	1:A:212:GLU:HB3	2.10	0.51
1:C:253:ILE:O	1:C:257:GLU:HG3	2.11	0.51
1:D:238:ARG:HD2	2:D:367:HOH:O	2.11	0.49
1:C:152:GLU:HG3	1:D:151:LEU:CD2	2.40	0.49
1:B:183:ALA:O	1:B:187:GLU:HG3	2.11	0.49
1:C:200:VAL:CG1	1:D:200:VAL:HG12	2.43	0.48
1:B:266:LYS:HB3	1:B:266:LYS:NZ	2.28	0.48
1:A:158:LEU:O	1:A:162:ASN:HB2	2.12	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:227:VAL:HG12	1:C:231:LYS:HE3	1.95	0.48
1:C:277:LEU:HD22	1:D:281:VAL:HG21	1.96	0.48
1:A:162:ASN:O	1:A:166:GLU:HG2	2.13	0.48
1:D:171:ARG:NE	2:D:381:HOH:O	2.30	0.47
1:B:239:ALA:O	1:B:243:GLU:HG3	2.14	0.47
1:C:260:LEU:HD21	1:D:259:GLU:CD	2.35	0.47
1:C:160:SER:O	1:C:163:TYR:HB3	2.16	0.46
1:A:162:ASN:ND2	1:B:161:LYS:HD3	2.30	0.46
1:A:151:LEU:HD11	1:B:155:VAL:HG21	1.98	0.46
1:C:261:TYR:CE2	1:C:265:LEU:HD11	2.50	0.46
1:C:152:GLU:CG	1:D:151:LEU:HD21	2.42	0.46
1:C:151:LEU:CB	1:D:151:LEU:HD13	2.45	0.46
1:A:268:LYS:HG3	2:A:343:HOH:O	2.16	0.45
1:B:149:LYS:O	1:B:152:GLU:HB2	2.15	0.45
1:D:217:LYS:HD3	1:D:221:TYR:HE2	1.81	0.45
1:A:149:LYS:O	1:A:153:ASP:OD1	2.35	0.45
1:A:218:GLU:CD	1:B:217:LYS:HZ1	2.20	0.45
1:C:272:GLU:O	1:C:276:GLN:HG3	2.17	0.45
1:A:204:LEU:C	1:A:204:LEU:HD23	2.37	0.45
1:A:162:ASN:HD22	1:B:161:LYS:HD3	1.82	0.45
1:B:171:ARG:NE	2:B:390:HOH:O	2.42	0.45
1:D:192:GLU:O	1:D:196:GLU:HG3	2.17	0.44
1:A:253:ILE:O	1:A:257:GLU:HG3	2.18	0.44
1:C:189:LYS:HE3	1:C:193:LEU:HD11	2.00	0.44
1:A:215:SER:O	1:A:218:GLU:HB3	2.17	0.44
1:C:298:LEU:HD11	1:D:299:LYS:HZ2	1.82	0.44
1:C:260:LEU:HD21	1:D:259:GLU:OE1	2.18	0.43
1:B:159:LEU:HA	1:B:159:LEU:HD23	1.88	0.43
1:D:205:LYS:HA	1:D:208:GLU:OE1	2.19	0.43
1:A:259:GLU:HG2	1:B:260:LEU:HD21	2.01	0.43
1:B:180:GLU:O	1:B:184:GLU:HG3	2.19	0.42
1:C:278:GLU:HG3	1:D:277:LEU:HD21	2.01	0.42
1:A:182:ARG:NH1	2:A:304:HOH:O	2.25	0.42
1:C:204:LEU:HD23	1:C:204:LEU:C	2.39	0.42
1:B:214:TYR:O	1:B:218:GLU:HB2	2.20	0.41
1:B:189:LYS:HE2	2:B:369:HOH:O	2.20	0.41
1:A:207:LEU:HD22	1:B:204:LEU:CD1	2.51	0.40
1:A:148:MET:HG3	1:B:151:LEU:CD1	2.46	0.40
1:D:251:LYS:HD3	1:D:255:ASP:OD2	2.20	0.40
1:D:169:VAL:CG1	1:D:173:LYS:HE3	2.51	0.40
1:D:274:MET:O	1:D:278:GLU:HG3	2.21	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	152/155 (98%)	152 (100%)	0	0	100	100
1	B	152/155 (98%)	151 (99%)	1 (1%)	0	100	100
1	C	152/155 (98%)	150 (99%)	2 (1%)	0	100	100
1	D	152/155 (98%)	151 (99%)	1 (1%)	0	100	100
All	All	608/620 (98%)	604 (99%)	4 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	141/142 (99%)	139 (99%)	2 (1%)	78	81
1	B	141/142 (99%)	139 (99%)	2 (1%)	78	81
1	C	141/142 (99%)	136 (96%)	5 (4%)	48	43
1	D	141/142 (99%)	139 (99%)	2 (1%)	78	81
All	All	564/568 (99%)	553 (98%)	11 (2%)	68	69

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	ASP
1	A	162	ASN
1	B	174	LYS
1	B	301	LEU
1	C	148	MET

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Mol	Chain	Res	Type
1	C	153	ASP
1	C	175	LEU
1	C	220	LYS
1	C	251	LYS
1	D	265	LEU
1	D	301	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	203	ASN
1	A	210	GLN
1	A	216	GLN
1	A	263	GLN
1	A	276	GLN
1	C	162	ASN
1	C	167	ASN
1	C	203	ASN
1	C	210	GLN
1	C	216	GLN
1	C	263	GLN
1	D	216	GLN
1	D	290	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates 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 ⓘ

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	154/155 (99%)	0.09	4 (2%) 53 53	30, 51, 74, 88	0
1	B	154/155 (99%)	-0.06	3 (1%) 64 64	29, 51, 72, 78	0
1	C	154/155 (99%)	0.16	2 (1%) 74 74	29, 51, 74, 86	0
1	D	154/155 (99%)	0.13	2 (1%) 74 74	31, 50, 71, 78	0
All	All	616/620 (99%)	0.08	11 (1%) 65 66	29, 51, 74, 88	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	163	TYR	4.0
1	A	155	VAL	3.6
1	D	301	LEU	3.0
1	B	214	TYR	2.9
1	A	148	MET	2.9
1	B	151	LEU	2.9
1	B	301	LEU	2.5
1	C	301	LEU	2.4
1	C	214	TYR	2.3
1	D	216	GLN	2.1
1	A	159	LEU	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.