



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:23 AM UTC

PDB ID : 2D3E / pdb\_00002d3e  
Title : Crystal structure of the C-Terminal fragment of rabbit skeletal alpha-tropomyosin  
Authors : Nitandai, Y.; Maeda, K.; Oda, N.; Minakata, S.; Maeda, Y.  
Deposited on : 2005-09-27  
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](mailto: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>.

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

MolProbity : 4-5-2 with Phenix2.0  
Xtrriage (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

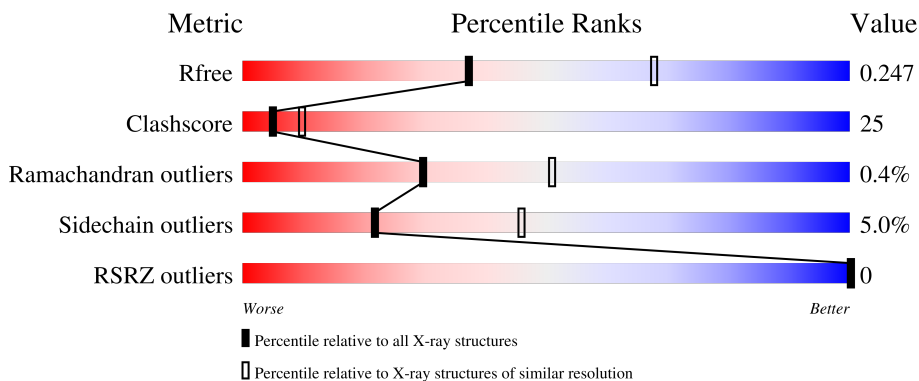
# 1 Overall quality at a glance

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  $\geq 3$ , 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	134	
1	B	134	
1	C	134	
1	D	134	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4288 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 General control protein GCN4 and Tropomyosin 1 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	130	1034	644	172	216	2	0	0	0
1	B	130	1037	644	172	219	2	0	0	0
1	C	130	1034	644	172	216	2	0	0	0
1	D	130	1037	644	172	219	2	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	151	MET	-	initiating methionine	UNP P58772
B	151	MET	-	initiating methionine	UNP P58772
C	151	MET	-	initiating methionine	UNP P58772
D	151	MET	-	initiating methionine	UNP P58772

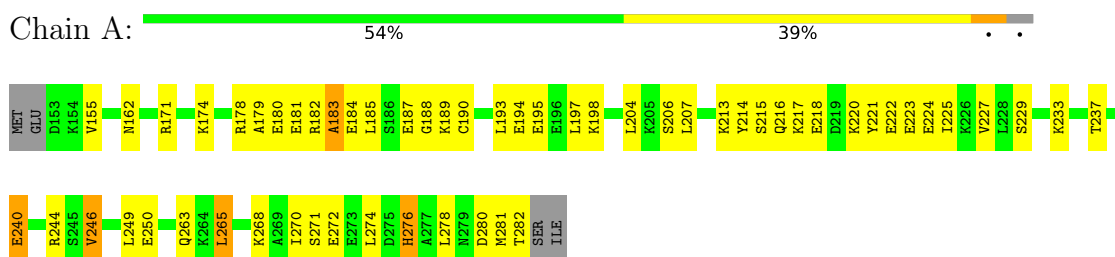
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	35	Total	O	0	0
			35	35		
2	B	37	Total	O	0	0
			37	37		
2	C	41	Total	O	0	0
			41	41		
2	D	33	Total	O	0	0
			33	33		

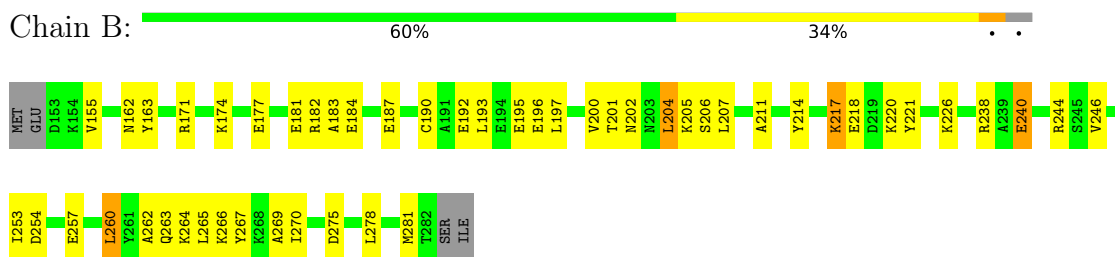
### 3 Residue-property plots

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.

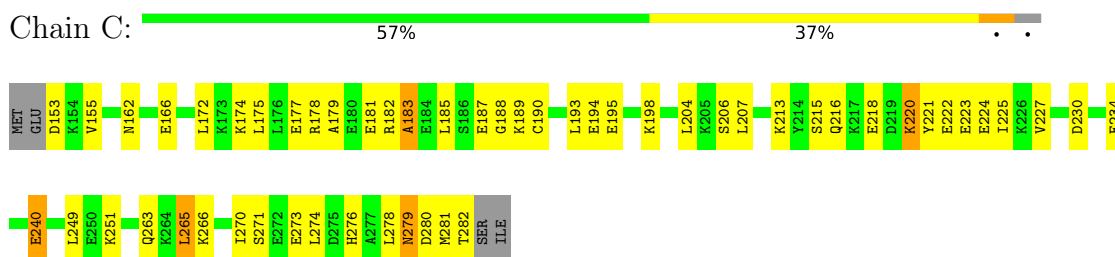
- Molecule 1: General control protein GCN4 and Tropomyosin 1 alpha chain



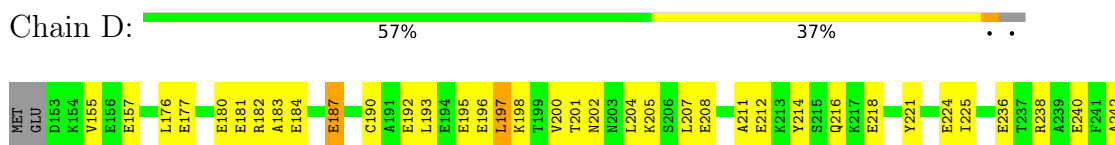
- Molecule 1: General control protein GCN4 and Tropomyosin 1 alpha chain



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



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





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	39.08Å 46.86Å 95.26Å 101.49° 101.66° 90.13°	Depositor
Resolution (Å)	45.67 – 2.60 45.67 – 2.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (45.67-2.60) 96.4 (45.67-2.60)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	9.32 (at 2.61Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.249 , 0.309 0.251 , 0.247	Depositor DCC
$R_{free}$ test set	989 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	53.1	Xtrriage
Anisotropy	0.680	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 98.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.460 for h,-k,-h-l 0.000 for -h,k,-k-l 0.000 for -h,-k,h+k+l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4288	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	90.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.54% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.38	0/1041	0.83	2/1393 (0.1%)
1	B	0.35	0/1044	0.85	2/1396 (0.1%)
1	C	0.37	0/1041	0.84	3/1393 (0.2%)
1	D	0.34	0/1044	0.87	2/1396 (0.1%)
All	All	0.36	0/4170	0.84	9/5578 (0.2%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	220	LYS	N-CA-C	-6.03	105.58	113.12
1	C	198	LYS	N-CA-C	-5.70	106.46	113.41
1	A	198	LYS	N-CA-C	-5.65	106.52	113.41
1	B	244	ARG	N-CA-C	-5.62	105.06	111.07
1	C	216	GLN	N-CA-C	-5.54	106.36	113.01
1	D	244	ARG	N-CA-C	-5.50	105.19	111.07
1	D	187	GLU	N-CA-C	-5.46	106.62	113.72
1	A	216	GLN	N-CA-C	-5.29	106.89	113.19
1	B	187	GLU	N-CA-C	-5.16	107.01	113.72

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	1034	0	1022	70	0
1	B	1037	0	1016	57	0
1	C	1034	0	1022	71	0
1	D	1037	0	1016	66	0
2	A	35	0	0	1	0
2	B	37	0	0	5	0
2	C	41	0	0	3	0
2	D	33	0	0	0	0
All	All	4288	0	4076	205	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 25.

All (205) 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:197:LEU:HD21	1:B:197:LEU:HB2	1.43	1.01
1:A:207:LEU:HD22	1:B:207:LEU:HD13	1.49	0.91
1:D:236:GLU:O	1:D:240:GLU:HG2	1.71	0.91
1:A:278:LEU:HD12	1:D:278:LEU:HD22	1.58	0.83
1:D:212:GLU:HG2	1:D:216:GLN:HE21	1.51	0.76
1:A:182:ARG:HE	1:B:183:ALA:HB1	1.51	0.75
1:B:211:ALA:HA	1:B:214:TYR:HD2	1.53	0.73
1:B:281:MET:HG2	1:D:270:ILE:HG13	1.70	0.73
1:C:278:LEU:HA	1:C:281:MET:HE3	1.70	0.72
1:A:204:LEU:HD23	1:B:204:LEU:HB2	1.71	0.72
1:B:171:ARG:O	1:B:174:LYS:HG2	1.91	0.71
1:A:278:LEU:HD21	1:D:274:LEU:HB3	1.73	0.71
1:C:207:LEU:HD22	1:D:207:LEU:HG	1.71	0.70
1:A:281:MET:SD	1:C:270:ILE:HD11	2.33	0.69
1:A:229:SER:O	1:A:233:LYS:HD3	1.92	0.69
1:D:211:ALA:O	1:D:214:TYR:HB2	1.93	0.68
1:D:177:GLU:O	1:D:181:GLU:HG2	1.93	0.68
1:D:184:GLU:O	1:D:187:GLU:HG2	1.94	0.68
1:D:205:LYS:O	1:D:205:LYS:HD3	1.94	0.67
1:D:253:ILE:O	1:D:257:GLU:HG3	1.94	0.67
1:C:204:LEU:HD23	1:D:204:LEU:HD23	1.76	0.67
1:D:211:ALA:HA	1:D:214:TYR:HD2	1.60	0.67
1:B:270:ILE:HG13	1:D:281:MET:HG2	1.78	0.66
1:B:211:ALA:HA	1:B:214:TYR:CD2	2.31	0.66
1:C:178:ARG:HG3	1:C:178:ARG:HH11	1.60	0.66
1:C:174:LYS:HE3	1:C:178:ARG:NH2	2.12	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:MET:HB3	1:D:267:TYR:OH	1.97	0.65
1:A:197:LEU:CD2	1:B:197:LEU:HB2	2.23	0.65
1:C:183:ALA:HB2	1:D:183:ALA:HB2	1.79	0.64
1:A:183:ALA:HB2	1:B:183:ALA:HB2	1.80	0.64
1:B:171:ARG:HG3	1:B:174:LYS:HE2	1.81	0.63
1:A:270:ILE:HD11	1:C:281:MET:SD	2.39	0.63
1:A:171:ARG:HH21	1:A:171:ARG:HG3	1.64	0.62
1:C:263:GLN:HG3	2:C:72:HOH:O	1.99	0.62
1:D:184:GLU:HA	1:D:187:GLU:OE1	2.00	0.61
1:C:194:GLU:HB3	1:D:193:LEU:HD13	1.83	0.60
1:C:189:LYS:HZ3	1:C:193:LEU:HD21	1.66	0.60
1:A:182:ARG:HE	1:B:183:ALA:CB	2.14	0.60
1:C:153:ASP:HA	2:C:81:HOH:O	2.01	0.60
1:D:246:VAL:O	1:D:250:GLU:HG3	2.01	0.60
1:B:202:ASN:C	1:B:204:LEU:H	2.09	0.59
1:D:202:ASN:C	1:D:204:LEU:H	2.10	0.59
1:B:211:ALA:O	1:B:214:TYR:HB2	2.02	0.58
1:B:278:LEU:HD22	1:C:278:LEU:HD22	1.85	0.58
1:A:246:VAL:O	1:A:250:GLU:HG3	2.03	0.58
1:C:249:LEU:HD11	1:D:246:VAL:HG13	1.85	0.58
1:A:223:GLU:HG3	1:A:224:GLU:N	2.18	0.58
1:A:193:LEU:O	1:A:197:LEU:HD23	2.05	0.57
1:A:271:SER:HA	1:D:281:MET:HE1	1.85	0.57
1:A:178:ARG:CB	1:A:178:ARG:HH11	2.18	0.57
1:A:194:GLU:HB2	1:B:193:LEU:HD13	1.86	0.57
1:A:207:LEU:C	1:A:207:LEU:HD23	2.30	0.56
1:A:278:LEU:HA	1:A:281:MET:HE2	1.87	0.56
1:B:202:ASN:HA	1:B:205:LYS:HG2	1.87	0.56
1:C:204:LEU:C	1:C:204:LEU:HD13	2.29	0.56
1:C:207:LEU:HD13	1:D:207:LEU:HB3	1.86	0.56
1:A:221:TYR:O	1:A:225:ILE:HG13	2.06	0.56
1:B:270:ILE:CD1	1:D:281:MET:HG2	2.36	0.56
1:A:204:LEU:C	1:A:204:LEU:HD13	2.31	0.56
1:C:266:LYS:HE3	1:D:267:TYR:CE2	2.41	0.56
1:A:197:LEU:HD22	1:B:197:LEU:HD13	1.87	0.56
1:A:180:GLU:O	1:A:184:GLU:HG2	2.06	0.55
1:C:174:LYS:HE3	1:C:178:ARG:HH21	1.71	0.55
1:B:204:LEU:C	1:B:204:LEU:HD13	2.32	0.55
1:B:262:ALA:O	1:B:265:LEU:HB2	2.06	0.55
1:A:178:ARG:HH11	1:A:178:ARG:HB2	1.72	0.55
1:A:281:MET:HG2	1:C:270:ILE:HD11	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:GLU:HA	1:B:240:GLU:OE1	2.06	0.54
1:C:204:LEU:CD2	1:D:204:LEU:HD23	2.37	0.54
1:C:279:ASN:C	1:C:279:ASN:HD22	2.15	0.54
1:D:205:LYS:HD3	1:D:205:LYS:C	2.33	0.54
1:C:278:LEU:HA	1:C:281:MET:CE	2.35	0.54
1:B:177:GLU:O	1:B:181:GLU:HG2	2.07	0.54
1:D:214:TYR:O	1:D:218:GLU:HG3	2.08	0.53
1:D:212:GLU:HG2	1:D:216:GLN:NE2	2.21	0.53
1:D:204:LEU:HD13	1:D:204:LEU:O	2.08	0.53
1:B:281:MET:HG2	1:D:270:ILE:CG1	2.39	0.53
1:A:207:LEU:HD23	1:A:207:LEU:O	2.10	0.52
1:B:226:LYS:HE2	2:B:98:HOH:O	2.10	0.52
1:D:204:LEU:HD13	1:D:204:LEU:C	2.34	0.52
1:C:276:HIS:HA	1:C:279:ASN:HD21	1.75	0.51
1:A:249:LEU:HD11	1:B:246:VAL:HG13	1.92	0.51
1:C:189:LYS:HB3	1:D:190:CYS:SG	2.51	0.51
1:C:185:LEU:C	1:C:185:LEU:HD23	2.35	0.51
1:A:185:LEU:C	1:A:185:LEU:HD23	2.36	0.51
1:C:207:LEU:HD22	1:D:207:LEU:CG	2.40	0.51
1:D:262:ALA:O	1:D:265:LEU:HB2	2.11	0.51
1:C:220:LYS:O	1:C:220:LYS:HD3	2.11	0.50
1:C:278:LEU:C	1:C:280:ASP:H	2.18	0.50
1:A:178:ARG:HB2	1:A:178:ARG:NH1	2.26	0.50
1:B:267:TYR:OH	1:C:281:MET:HB3	2.11	0.50
1:C:218:GLU:O	1:C:222:GLU:HG2	2.12	0.50
1:A:155:VAL:HG22	1:B:155:VAL:HG22	1.92	0.50
1:A:204:LEU:CD2	1:B:204:LEU:HD23	2.42	0.50
1:A:278:LEU:C	1:A:280:ASP:H	2.18	0.50
1:B:270:ILE:CG1	1:D:281:MET:HG2	2.41	0.50
1:C:178:ARG:HG3	1:C:178:ARG:NH1	2.26	0.50
1:C:182:ARG:HH11	1:D:183:ALA:CB	2.24	0.50
1:C:182:ARG:HH11	1:D:183:ALA:HB3	1.77	0.50
1:B:197:LEU:O	1:B:201:THR:N	2.45	0.50
1:D:197:LEU:O	1:D:201:THR:N	2.44	0.50
1:B:281:MET:HG2	1:D:270:ILE:CD1	2.43	0.49
1:A:263:GLN:HG3	2:B:74:HOH:O	2.13	0.49
1:C:207:LEU:HB3	1:D:207:LEU:HD23	1.95	0.49
1:A:233:LYS:HD2	1:A:233:LYS:N	2.28	0.49
1:C:207:LEU:HD22	1:D:207:LEU:CD2	2.43	0.49
1:C:274:LEU:CD1	1:D:274:LEU:HD11	2.43	0.49
1:C:195:GLU:OE2	1:C:195:GLU:HA	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:266:LYS:O	1:D:269:ALA:HB3	2.13	0.48
1:C:221:TYR:O	1:C:225:ILE:HG13	2.14	0.48
1:A:281:MET:CG	1:C:270:ILE:HD11	2.44	0.48
1:D:242:ALA:O	1:D:246:VAL:HG23	2.14	0.48
1:B:217:LYS:HD2	1:B:221:TYR:CE2	2.49	0.48
1:C:162:ASN:O	1:C:166:GLU:HG3	2.14	0.48
1:B:281:MET:HE1	1:C:271:SER:HA	1.95	0.47
1:C:189:LYS:HZ2	1:C:193:LEU:HD11	1.79	0.47
1:A:274:LEU:HD12	1:D:278:LEU:HD13	1.97	0.47
1:B:205:LYS:HG3	1:B:206:SER:N	2.29	0.47
1:D:197:LEU:C	1:D:197:LEU:HD12	2.39	0.47
1:A:222:GLU:HA	1:A:222:GLU:OE1	2.14	0.47
1:C:155:VAL:HG22	1:D:155:VAL:HG22	1.95	0.47
1:C:234:GLU:CD	1:C:234:GLU:C	2.81	0.47
1:C:189:LYS:NZ	1:C:193:LEU:HD11	2.29	0.47
1:D:281:MET:O	1:D:282:THR:HB	2.14	0.47
1:A:223:GLU:HG3	1:A:224:GLU:H	1.78	0.47
1:A:162:ASN:HD21	1:B:162:ASN:HB2	1.80	0.47
1:D:200:VAL:O	1:D:200:VAL:HG12	2.16	0.46
1:C:274:LEU:HD11	1:D:274:LEU:HD11	1.98	0.46
1:B:207:LEU:O	1:B:211:ALA:HB2	2.16	0.46
1:C:189:LYS:NZ	1:C:193:LEU:HD21	2.31	0.46
1:C:207:LEU:HD23	1:C:207:LEU:O	2.15	0.46
1:C:174:LYS:O	1:C:178:ARG:HG2	2.16	0.46
1:A:207:LEU:CD2	1:B:207:LEU:HB3	2.46	0.45
1:C:204:LEU:HA	1:D:204:LEU:HD23	1.99	0.45
1:B:204:LEU:HD13	1:B:204:LEU:O	2.16	0.45
1:D:207:LEU:O	1:D:211:ALA:HB2	2.17	0.45
1:C:224:GLU:HA	1:C:227:VAL:HG12	1.98	0.45
1:A:233:LYS:O	1:A:237:THR:HG23	2.17	0.45
1:D:176:LEU:O	1:D:180:GLU:HG3	2.17	0.45
1:A:204:LEU:HD23	1:B:204:LEU:HD23	1.99	0.45
1:A:224:GLU:O	1:A:227:VAL:HG12	2.17	0.45
1:C:188:GLY:C	1:C:190:CYS:N	2.73	0.45
1:A:217:LYS:HE2	1:A:221:TYR:CZ	2.52	0.44
1:B:253:ILE:O	1:B:257:GLU:HG3	2.17	0.44
1:A:171:ARG:HH21	1:A:171:ARG:CG	2.27	0.44
1:A:188:GLY:C	1:A:190:CYS:N	2.73	0.44
1:A:265:LEU:HD12	1:A:265:LEU:HA	1.84	0.44
1:C:204:LEU:C	1:C:206:SER:N	2.73	0.44
1:C:240:GLU:OE1	1:D:238:ARG:NH2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:275:ASP:HA	1:C:278:LEU:HD21	1.99	0.44
1:D:197:LEU:HD12	1:D:198:LYS:N	2.33	0.44
1:B:200:VAL:HG12	1:B:200:VAL:O	2.18	0.44
1:A:224:GLU:HA	1:A:227:VAL:HG12	2.00	0.43
1:A:213:LYS:C	1:A:215:SER:H	2.26	0.43
1:A:270:ILE:HD11	1:C:281:MET:CG	2.47	0.43
1:B:220:LYS:HE2	2:B:99:HOH:O	2.17	0.43
1:A:189:LYS:HB3	1:B:190:CYS:SG	2.58	0.43
1:B:214:TYR:O	1:B:218:GLU:HG3	2.19	0.43
1:C:279:ASN:C	1:C:279:ASN:ND2	2.76	0.43
1:A:204:LEU:C	1:A:206:SER:N	2.73	0.43
1:D:251:LYS:O	1:D:254:ASP:HB2	2.19	0.43
1:B:192:GLU:O	1:B:195:GLU:HB3	2.18	0.43
1:A:185:LEU:C	1:A:187:GLU:N	2.75	0.43
1:C:179:ALA:C	1:C:181:GLU:N	2.76	0.43
1:D:182:ARG:C	1:D:184:GLU:H	2.27	0.42
1:A:179:ALA:C	1:A:181:GLU:N	2.75	0.42
1:B:266:LYS:O	1:B:269:ALA:HB3	2.19	0.42
1:A:195:GLU:HA	1:A:195:GLU:OE1	2.19	0.42
1:A:276:HIS:ND1	1:A:276:HIS:C	2.77	0.42
1:C:213:LYS:C	1:C:215:SER:H	2.27	0.42
1:B:195:GLU:O	1:B:196:GLU:C	2.62	0.42
1:C:182:ARG:HE	1:D:183:ALA:HB1	1.85	0.42
1:A:268:LYS:NZ	1:A:272:GLU:OE1	2.50	0.42
1:C:223:GLU:HG3	1:C:224:GLU:N	2.34	0.42
1:D:184:GLU:HA	1:D:187:GLU:CD	2.45	0.42
1:C:185:LEU:C	1:C:187:GLU:N	2.76	0.42
1:C:172:LEU:HD23	1:C:172:LEU:HA	1.94	0.42
1:D:195:GLU:O	1:D:196:GLU:C	2.62	0.42
1:A:183:ALA:O	1:A:187:GLU:HB2	2.20	0.42
1:A:278:LEU:C	1:A:280:ASP:N	2.78	0.42
1:B:182:ARG:C	1:B:184:GLU:H	2.28	0.42
1:C:251:LYS:HE3	2:C:125:HOH:O	2.20	0.42
1:A:240:GLU:OE1	1:B:238:ARG:NH2	2.53	0.42
1:C:278:LEU:C	1:C:280:ASP:N	2.78	0.42
1:A:206:SER:O	1:A:207:LEU:C	2.62	0.41
1:D:184:GLU:HA	1:D:187:GLU:HG2	2.02	0.41
1:D:221:TYR:O	1:D:225:ILE:HG13	2.20	0.41
1:C:207:LEU:CD2	1:D:207:LEU:HG	2.47	0.41
1:D:192:GLU:HG3	1:D:193:LEU:N	2.36	0.41
1:C:188:GLY:C	1:C:190:CYS:H	2.27	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:TYR:HD2	2:B:4:HOH:O	2.03	0.41
1:D:204:LEU:O	1:D:208:GLU:HG3	2.20	0.41
1:B:260:LEU:O	1:B:264:LYS:HG3	2.21	0.41
1:C:206:SER:O	1:C:207:LEU:C	2.63	0.41
1:A:214:TYR:O	1:A:218:GLU:HG3	2.21	0.41
1:B:278:LEU:CD2	1:C:278:LEU:HD22	2.48	0.41
1:A:215:SER:C	1:A:217:LYS:H	2.28	0.41
1:A:184:GLU:HA	1:A:187:GLU:HB3	2.03	0.41
1:B:267:TYR:HB2	2:B:74:HOH:O	2.21	0.41
1:B:202:ASN:O	1:B:205:LYS:HG2	2.20	0.41
1:A:188:GLY:C	1:A:190:CYS:H	2.28	0.40
1:A:220:LYS:HE2	2:A:138:HOH:O	2.20	0.40
1:A:174:LYS:HE3	1:A:178:ARG:HE	1.86	0.40
1:C:265:LEU:HD12	1:C:265:LEU:HA	1.95	0.40
1:C:175:LEU:C	1:C:177:GLU:N	2.79	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	128/134 (96%)	115 (90%)	12 (9%)	1 (1%)	16	34
1	B	128/134 (96%)	121 (94%)	7 (6%)	0	100	100
1	C	128/134 (96%)	115 (90%)	12 (9%)	1 (1%)	16	34
1	D	128/134 (96%)	121 (94%)	7 (6%)	0	100	100
All	All	512/536 (96%)	472 (92%)	38 (7%)	2 (0%)	30	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	183	ALA
1	C	183	ALA

### 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	109/121 (90%)	103 (94%)	6 (6%)	19	41
1	B	109/121 (90%)	103 (94%)	6 (6%)	19	41
1	C	109/121 (90%)	103 (94%)	6 (6%)	19	41
1	D	109/121 (90%)	105 (96%)	4 (4%)	30	57
All	All	436/484 (90%)	414 (95%)	22 (5%)	22	46

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	240	GLU
1	A	244	ARG
1	A	246	VAL
1	A	265	LEU
1	A	276	HIS
1	A	282	THR
1	B	204	LEU
1	B	217	LYS
1	B	240	GLU
1	B	254	ASP
1	B	260	LEU
1	B	263	GLN
1	C	230	ASP
1	C	240	GLU
1	C	265	LEU
1	C	273	GLU
1	C	279	ASN
1	C	282	THR
1	D	157	GLU
1	D	197	LEU

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Mol	Chain	Res	Type
1	D	224	GLU
1	D	274	LEU

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	162	ASN
1	A	210	GLN
1	A	263	GLN
1	B	210	GLN
1	B	263	GLN
1	C	167	ASN
1	C	216	GLN
1	C	263	GLN
1	C	279	ASN
1	D	210	GLN
1	D	216	GLN

### 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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	130/134 (97%)	-0.79	0 100 100	45, 90, 135, 156	0
1	B	130/134 (97%)	-0.78	0 100 100	38, 83, 140, 152	0
1	C	130/134 (97%)	-0.83	0 100 100	43, 95, 134, 157	0
1	D	130/134 (97%)	-0.73	0 100 100	39, 82, 141, 149	0
All	All	520/536 (97%)	-0.78	0 100 100	38, 87, 140, 157	0

There are no RSRZ outliers to report.

### 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.