



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:25 PM GMT

PDB ID : 1TJV  
Title : Crystal Structure of T161D Duck Delta 2 Crystallin Mutant  
Authors : Sampaleanu, L.M.; Coddling, P.W.; Lobsanov, Y.D.; Tsai, M.; Smith, G.D.; Horvatin, C.; Howell, P.L.  
Deposited on : 2004-06-07  
Resolution : 2.00 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

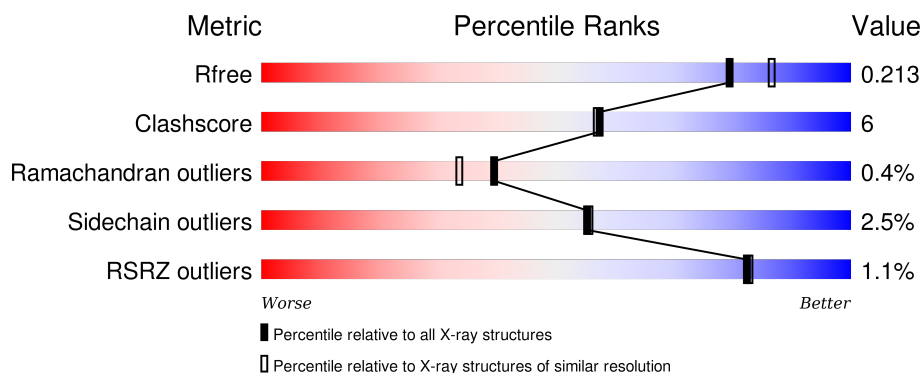
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	474	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 13%, green 81%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>81%</span> <span>13%</span> <span>• 5%</span> </div> </div>
1	B	474	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, yellow 12%, green 82%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span></span> <span>82%</span> <span>12%</span> <span>• 5%</span> </div> </div>
1	C	474	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 1%, yellow 12%, green 82%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>2%</span> <span>82%</span> <span>12%</span> <span>• 5%</span> </div> </div>
1	D	474	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 83%, yellow 10%, grey 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span></span> <span>83%</span> <span>10%</span> <span>• 5%</span> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14813 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 Delta crystallin II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	449	Total	C	N	O	S	0	0	0
			3470	2197	585	676	12			
1	B	449	Total	C	N	O	S	0	0	0
			3468	2195	587	674	12			
1	C	450	Total	C	N	O	S	0	0	0
			3469	2196	587	674	12			
1	D	450	Total	C	N	O	S	0	0	0
			3481	2203	588	678	12			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	161	ASP	THR	ENGINEERED	UNP P24058
B	161	ASP	THR	ENGINEERED	UNP P24058
C	161	ASP	THR	ENGINEERED	UNP P24058
D	161	ASP	THR	ENGINEERED	UNP P24058
A	469	HIS	-	EXPRESSION TAG	UNP P24058
A	470	HIS	-	EXPRESSION TAG	UNP P24058
A	471	HIS	-	EXPRESSION TAG	UNP P24058
A	472	HIS	-	EXPRESSION TAG	UNP P24058
A	473	HIS	-	EXPRESSION TAG	UNP P24058
A	474	HIS	-	EXPRESSION TAG	UNP P24058
B	469	HIS	-	EXPRESSION TAG	UNP P24058
B	470	HIS	-	EXPRESSION TAG	UNP P24058
B	471	HIS	-	EXPRESSION TAG	UNP P24058
B	472	HIS	-	EXPRESSION TAG	UNP P24058
B	473	HIS	-	EXPRESSION TAG	UNP P24058
B	474	HIS	-	EXPRESSION TAG	UNP P24058
C	469	HIS	-	EXPRESSION TAG	UNP P24058
C	470	HIS	-	EXPRESSION TAG	UNP P24058
C	471	HIS	-	EXPRESSION TAG	UNP P24058
C	472	HIS	-	EXPRESSION TAG	UNP P24058
C	473	HIS	-	EXPRESSION TAG	UNP P24058

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Chain	Residue	Modelled	Actual	Comment	Reference
C	474	HIS	-	EXPRESSION TAG	UNP P24058
D	469	HIS	-	EXPRESSION TAG	UNP P24058
D	470	HIS	-	EXPRESSION TAG	UNP P24058
D	471	HIS	-	EXPRESSION TAG	UNP P24058
D	472	HIS	-	EXPRESSION TAG	UNP P24058
D	473	HIS	-	EXPRESSION TAG	UNP P24058
D	474	HIS	-	EXPRESSION TAG	UNP P24058

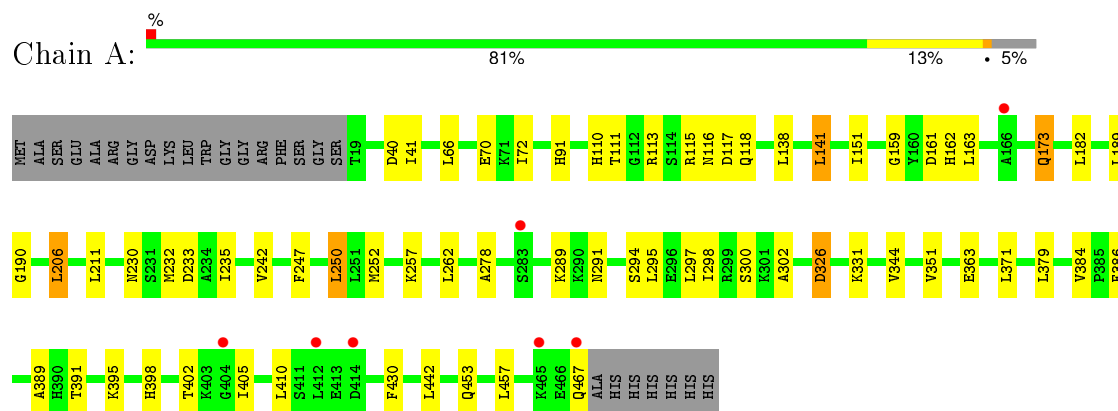
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	247	Total O 247 247	0	0
2	B	214	Total O 214 214	0	0
2	C	229	Total O 229 229	0	0
2	D	235	Total O 235 235	0	0

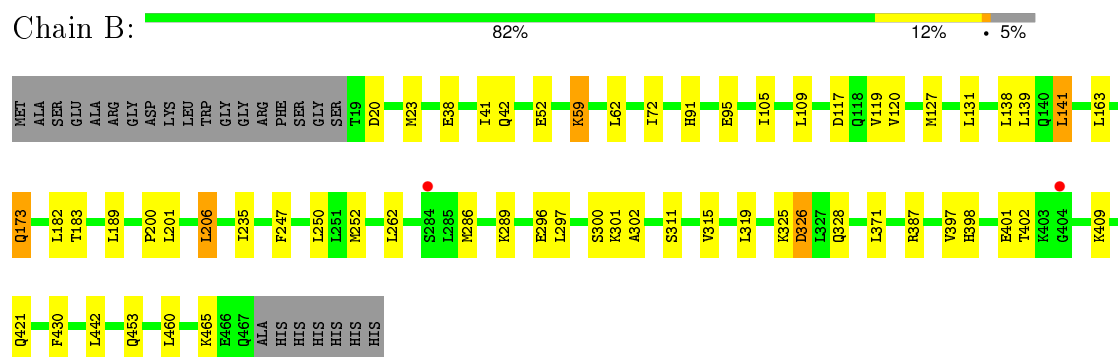
### 3 Residue-property plots

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.

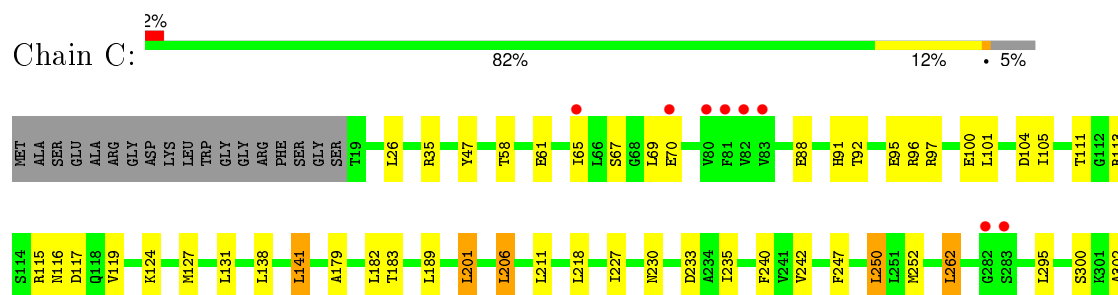
#### • Molecule 1: Delta crystallin II



#### • Molecule 1: Delta crystallin II

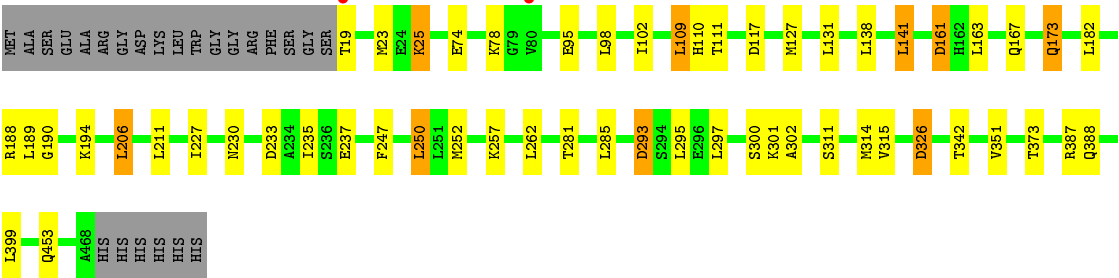
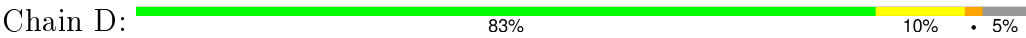


#### • Molecule 1: Delta crystallin II





● Molecule 1: Delta crystallin II



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.13 Å 98.48 Å 103.47 Å 90.00° 100.10° 90.00°	Depositor
Resolution (Å)	43.38 – 2.00 43.38 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.38-2.00) 99.9 (43.38-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	22.99 (at 2.00 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.174 , 0.213 0.174 , 0.213	Depositor DCC
$R_{free}$ test set	12399 reflections (9.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	16.2	Xtriage
Anisotropy	0.225	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 55.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 124250 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	14813	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.43	0/3515	0.61	1/4743 (0.0%)
1	B	0.41	0/3513	0.60	0/4741
1	C	0.40	0/3514	0.60	0/4743
1	D	0.41	0/3526	0.60	0/4757
All	All	0.41	0/14068	0.60	1/18984 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	159	GLY	N-CA-C	-5.26	99.94	113.10

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	3470	0	3562	51	0
1	B	3468	0	3558	45	0
1	C	3469	0	3554	51	0
1	D	3481	0	3576	47	0
2	A	247	0	0	2	0
2	B	214	0	0	3	0
2	C	229	0	0	4	0
2	D	235	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	14813	0	14250	170	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 6.

All (170) 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:B:183:THR:HG21	1:B:460:LEU:HD13	1.51	0.91
1:B:409:LYS:HA	1:B:409:LYS:HE2	1.53	0.89
1:D:387:ARG:C	1:D:387:ARG:HD3	1.99	0.82
1:D:173:GLN:HE22	1:D:453:GLN:HE22	1.27	0.82
1:A:278:ALA:HA	1:D:19:THR:HB	1.64	0.78
1:A:442:LEU:HD13	1:C:227:ILE:HD11	1.64	0.78
1:A:379:LEU:HB3	1:A:384:VAL:HG21	1.66	0.77
1:C:88:GLU:OE2	1:C:92:THR:HG21	1.87	0.75
1:B:173:GLN:HE22	1:B:453:GLN:HE22	1.35	0.74
1:C:67:SER:O	1:C:70:GLU:HG2	1.89	0.73
1:D:230:ASN:HD22	1:D:233:ASP:H	1.36	0.72
1:B:91:HIS:O	1:B:95:GLU:HG3	1.91	0.71
1:A:173:GLN:HE22	1:A:453:GLN:HE22	1.36	0.71
1:A:384:VAL:HG23	1:A:389:ALA:HB2	1.73	0.70
1:C:111:THR:HG22	1:C:211:LEU:HD11	1.74	0.70
1:A:151:ILE:HB	2:A:619:HOH:O	1.92	0.70
1:A:230:ASN:HD22	1:A:233:ASP:H	1.38	0.69
1:D:127:MET:O	1:D:131:LEU:HD13	1.92	0.69
1:A:110:HIS:HD2	1:A:113:ARG:HE	1.42	0.68
1:C:230:ASN:HD22	1:C:233:ASP:H	1.41	0.68
1:B:297:LEU:HD21	1:C:26:LEU:HD13	1.77	0.68
1:B:20:ASP:HB3	1:B:23:MET:HB2	1.76	0.67
1:C:386:PHE:HD2	1:C:387:ARG:HD2	1.60	0.65
1:B:173:GLN:HA	1:B:173:GLN:HE21	1.63	0.64
1:B:117:ASP:HB3	1:B:235:ILE:HD11	1.81	0.62
1:B:141:LEU:HD13	1:B:182:LEU:HD13	1.82	0.61
1:C:91:HIS:O	1:C:95:GLU:HG3	2.00	0.61
1:A:111:THR:HG22	1:A:211:LEU:HD11	1.81	0.60
1:A:379:LEU:HB3	1:A:384:VAL:CG2	2.32	0.60
1:A:115:ARG:NH2	1:A:118:GLN:HE22	1.99	0.60
1:C:141:LEU:HD13	1:C:182:LEU:HD13	1.83	0.59
1:C:442:LEU:HD23	2:C:510:HOH:O	2.03	0.59
1:C:127:MET:O	1:C:131:LEU:HD13	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:387:ARG:C	1:D:387:ARG:CD	2.71	0.58
1:C:97:ARG:CZ	1:C:101:LEU:HD21	2.33	0.58
1:A:379:LEU:O	1:A:384:VAL:HG22	2.04	0.58
1:A:115:ARG:HH21	1:A:118:GLN:HE22	1.52	0.58
1:A:297:LEU:HD12	1:D:23:MET:CE	2.34	0.57
1:C:386:PHE:CD2	1:C:387:ARG:HD2	2.39	0.57
1:C:387:ARG:HD3	2:C:518:HOH:O	2.04	0.56
1:A:384:VAL:CG2	1:A:389:ALA:HB2	2.35	0.56
1:A:379:LEU:C	1:A:384:VAL:HG22	2.27	0.55
1:C:396:ALA:HA	1:C:418:ILE:CD1	2.36	0.55
1:A:297:LEU:HD12	1:D:23:MET:SD	2.47	0.55
1:D:230:ASN:ND2	1:D:233:ASP:H	2.05	0.55
1:D:387:ARG:HD3	1:D:387:ARG:O	2.07	0.54
1:C:117:ASP:HB3	1:C:235:ILE:HD11	1.90	0.54
1:D:173:GLN:HA	1:D:173:GLN:HE21	1.71	0.54
1:D:25:LYS:N	1:D:25:LYS:HE3	2.23	0.53
1:B:397:VAL:O	1:B:401:GLU:HG3	2.08	0.53
1:A:391:THR:HG22	1:A:395:LYS:HE3	1.90	0.53
1:D:295:LEU:HD21	1:D:351:VAL:HG11	1.90	0.53
1:C:179:ALA:O	1:C:183:THR:HG23	2.08	0.53
1:B:297:LEU:HD21	1:C:26:LEU:CD1	2.38	0.53
1:B:398:HIS:O	1:B:402:THR:HG23	2.09	0.53
1:D:342:THR:HG22	2:D:651:HOH:O	2.08	0.52
1:B:183:THR:CG2	1:B:460:LEU:HD13	2.32	0.52
1:D:161:ASP:HB3	2:D:709:HOH:O	2.09	0.52
1:B:206:LEU:HB3	1:D:167:GLN:HE21	1.74	0.52
1:A:297:LEU:HD12	1:D:23:MET:HE1	1.92	0.52
1:C:206:LEU:HD23	1:C:206:LEU:C	2.30	0.51
1:D:111:THR:HG22	1:D:211:LEU:HD11	1.91	0.51
1:A:173:GLN:HE21	1:A:173:GLN:HA	1.75	0.51
1:A:141:LEU:HD13	1:A:182:LEU:HD13	1.93	0.51
1:D:95:GLU:OE1	1:D:110:HIS:HE1	1.94	0.50
1:C:96:ARG:HG2	1:C:96:ARG:HH11	1.77	0.50
1:C:35:ARG:HD2	2:C:604:HOH:O	2.11	0.50
1:D:252:MET:HB3	1:D:302:ALA:HA	1.94	0.50
1:D:206:LEU:HD23	1:D:206:LEU:C	2.33	0.49
1:B:442:LEU:HD13	1:D:227:ILE:HD11	1.95	0.49
1:B:289:LYS:NZ	2:B:688:HOH:O	2.35	0.49
1:B:325:LYS:HG3	1:B:325:LYS:O	2.12	0.48
1:A:242:VAL:HG11	1:C:250:LEU:HG	1.96	0.48
1:B:371:LEU:HD13	1:B:430:PHE:HA	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:ASN:HB3	1:A:235:ILE:HD12	1.95	0.48
1:B:326:ASP:HA	1:C:300:SER:HB3	1.96	0.48
1:A:300:SER:HB3	1:D:326:ASP:HA	1.95	0.48
1:B:52:GLU:HG3	1:B:62:LEU:HD22	1.96	0.48
1:D:117:ASP:HB3	1:D:235:ILE:HD11	1.96	0.48
1:B:300:SER:HB3	1:C:326:ASP:HA	1.96	0.48
1:C:189:LEU:HD23	1:C:189:LEU:C	2.34	0.48
1:B:297:LEU:CD2	1:C:26:LEU:HD13	2.43	0.48
1:A:230:ASN:HD21	1:A:232:MET:HB2	1.79	0.47
1:A:66:LEU:O	1:A:70:GLU:HG3	2.14	0.47
1:A:295:LEU:HD21	1:A:351:VAL:HG11	1.96	0.47
1:C:124:LYS:HE2	1:C:240:PHE:CE2	2.48	0.47
1:C:396:ALA:HA	1:C:418:ILE:HD13	1.95	0.47
1:D:74:GLU:HG3	1:D:78:LYS:HE3	1.97	0.47
1:C:58:THR:OG1	1:C:61:GLU:HG3	2.15	0.47
1:A:117:ASP:HB3	1:A:235:ILE:HD11	1.97	0.47
1:A:331:LYS:HE3	2:A:529:HOH:O	2.15	0.47
1:D:109:LEU:HD12	1:D:109:LEU:HA	1.80	0.46
1:B:297:LEU:HD21	1:C:26:LEU:CB	2.45	0.46
1:D:188:ARG:NH2	1:D:250:LEU:HD13	2.31	0.46
1:B:289:LYS:HG3	1:B:289:LYS:O	2.16	0.46
1:C:262:LEU:HB3	1:C:295:LEU:CD1	2.46	0.46
1:A:40:ASP:OD2	1:A:91:HIS:HD2	1.97	0.46
1:C:252:MET:HB3	1:C:302:ALA:HA	1.97	0.46
1:D:387:ARG:HD3	1:D:388:GLN:N	2.29	0.46
1:C:96:ARG:O	1:C:100:GLU:HG3	2.16	0.46
1:A:161:ASP:HB3	2:B:688:HOH:O	2.15	0.45
1:A:189:LEU:O	1:A:189:LEU:HD23	2.17	0.45
1:A:41:ILE:HD11	1:A:72:ILE:HG22	1.98	0.45
1:B:286:MET:CB	1:B:289:LYS:HG2	2.46	0.45
1:D:297:LEU:O	1:D:301:LYS:HG2	2.16	0.45
1:C:116:ASN:HB3	1:C:235:ILE:HD12	1.98	0.45
1:B:189:LEU:HD23	1:B:189:LEU:O	2.17	0.45
1:B:252:MET:HB3	1:B:302:ALA:HA	1.99	0.45
1:A:298:ILE:HG12	1:A:344:VAL:HG13	1.98	0.45
1:C:115:ARG:O	1:C:119:VAL:HG23	2.17	0.45
1:B:139:LEU:HD13	1:B:465:LYS:HE2	1.98	0.45
1:B:41:ILE:HD11	1:B:72:ILE:HG22	1.99	0.45
1:C:442:LEU:HD22	1:C:442:LEU:N	2.32	0.44
1:C:124:LYS:HE2	1:C:240:PHE:CD2	2.52	0.44
1:C:390:HIS:CE1	1:D:285:LEU:HD13	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:HIS:CD2	1:A:113:ARG:HE	2.29	0.44
1:D:399:LEU:HD23	1:D:399:LEU:O	2.18	0.44
1:B:127:MET:O	1:B:131:LEU:HD13	2.17	0.44
1:C:61:GLU:HB3	1:C:105:ILE:HD11	2.00	0.44
1:C:262:LEU:HB3	1:C:295:LEU:HD11	2.00	0.44
1:A:326:ASP:HA	1:D:300:SER:HB3	2.00	0.44
1:A:371:LEU:HD13	1:A:430:PHE:HA	2.00	0.44
1:A:405:ILE:HD11	1:A:410:LEU:HD23	2.00	0.43
1:A:115:ARG:HH21	1:A:118:GLN:NE2	2.14	0.43
1:D:98:LEU:O	1:D:102:ILE:HG12	2.18	0.43
1:B:421:GLN:NE2	1:B:421:GLN:HA	2.33	0.43
1:B:319:LEU:O	1:D:257:LYS:HG3	2.18	0.43
1:B:286:MET:HB3	1:B:289:LYS:HG2	2.00	0.43
1:B:105:ILE:HG23	2:B:543:HOH:O	2.19	0.43
1:A:190:GLY:HA3	1:A:467:GLN:HE22	1.83	0.43
1:A:398:HIS:O	1:A:402:THR:HG23	2.18	0.43
1:A:206:LEU:HD23	1:A:206:LEU:C	2.39	0.43
1:A:162:HIS:HA	1:B:296:GLU:OE1	2.18	0.43
1:A:391:THR:O	1:A:395:LYS:HG3	2.18	0.43
1:D:262:LEU:HB3	1:D:295:LEU:CD1	2.49	0.43
1:A:252:MET:HB3	1:A:302:ALA:HA	2.01	0.43
1:C:97:ARG:NH2	1:C:101:LEU:HD21	2.34	0.42
1:B:325:LYS:O	1:B:328:GLN:HG3	2.19	0.42
1:D:373:THR:HB	2:D:665:HOH:O	2.18	0.42
1:B:297:LEU:O	1:B:301:LYS:HG2	2.19	0.42
1:C:314:MET:HA	1:D:314:MET:SD	2.59	0.42
1:A:250:LEU:HG	1:C:242:VAL:HG11	2.00	0.42
1:D:190:GLY:O	1:D:194:LYS:HD3	2.19	0.42
1:B:409:LYS:CA	1:B:409:LYS:HE2	2.34	0.42
1:D:173:GLN:HE22	1:D:453:GLN:NE2	2.06	0.42
1:D:74:GLU:OE1	1:D:78:LYS:HE3	2.19	0.42
1:C:47:TYR:CE2	1:C:113:ARG:HB2	2.54	0.42
1:C:65:ILE:O	1:C:69:LEU:HG	2.20	0.42
1:D:141:LEU:HD13	1:D:182:LEU:HD13	2.01	0.42
1:B:119:VAL:HG13	1:B:120:VAL:N	2.34	0.42
1:C:385:PRO:HD2	1:C:388:GLN:HG3	2.02	0.41
1:B:38:GLU:O	1:B:42:GLN:HG3	2.20	0.41
1:D:74:GLU:CG	1:D:78:LYS:HE3	2.51	0.41
1:C:201:LEU:HD13	1:C:218:LEU:HD13	2.02	0.41
1:D:189:LEU:HD23	1:D:189:LEU:O	2.21	0.41
1:D:163:LEU:HA	1:D:163:LEU:HD12	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:311:SER:O	1:D:315:VAL:HG23	2.20	0.41
1:A:189:LEU:C	1:A:189:LEU:HD23	2.41	0.41
1:C:463:LYS:HD2	2:C:633:HOH:O	2.20	0.41
1:B:163:LEU:HA	1:B:163:LEU:HD12	1.83	0.41
1:B:297:LEU:HD21	1:C:26:LEU:HB3	2.03	0.41
1:B:59:LYS:HB2	1:B:59:LYS:HE3	1.93	0.41
1:A:289:LYS:NZ	1:A:291:ASN:HD21	2.19	0.41
1:B:311:SER:O	1:B:315:VAL:HG23	2.21	0.40
1:A:257:LYS:HG3	1:C:319:LEU:O	2.20	0.40
1:B:206:LEU:C	1:B:206:LEU:HD23	2.41	0.40
1:A:163:LEU:HD12	1:A:163:LEU:HA	1.86	0.40
1:D:281:THR:HB	1:D:293:ASP:OD1	2.21	0.40
1:C:88:GLU:CD	1:C:88:GLU:H	2.25	0.40
1:A:294:SER:HA	1:D:23:MET:HE1	2.03	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	447/474 (94%)	439 (98%)	6 (1%)	2 (0%)	39	33
1	B	447/474 (94%)	437 (98%)	8 (2%)	2 (0%)	39	33
1	C	448/474 (94%)	435 (97%)	11 (2%)	2 (0%)	39	33
1	D	448/474 (94%)	437 (98%)	10 (2%)	1 (0%)	52	48
All	All	1790/1896 (94%)	1748 (98%)	35 (2%)	7 (0%)	39	33

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	206	LEU

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Mol	Chain	Res	Type
1	B	206	LEU
1	C	206	LEU
1	D	206	LEU
1	A	386	PHE
1	C	104	ASP
1	B	200	PRO

### 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	388/411 (94%)	379 (98%)	9 (2%)	58	60
1	B	387/411 (94%)	376 (97%)	11 (3%)	51	50
1	C	386/411 (94%)	378 (98%)	8 (2%)	61	63
1	D	389/411 (95%)	378 (97%)	11 (3%)	51	50
All	All	1550/1644 (94%)	1511 (98%)	39 (2%)	55	55

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	LEU
1	A	141	LEU
1	A	173	GLN
1	A	247	PHE
1	A	250	LEU
1	A	262	LEU
1	A	326	ASP
1	A	363	GLU
1	A	457	LEU
1	B	59	LYS
1	B	109	LEU
1	B	138	LEU
1	B	141	LEU
1	B	173	GLN
1	B	201	LEU

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Mol	Chain	Res	Type
1	B	247	PHE
1	B	250	LEU
1	B	262	LEU
1	B	326	ASP
1	B	387	ARG
1	C	138	LEU
1	C	141	LEU
1	C	201	LEU
1	C	247	PHE
1	C	250	LEU
1	C	262	LEU
1	C	387	ARG
1	C	388	GLN
1	D	25	LYS
1	D	109	LEU
1	D	138	LEU
1	D	141	LEU
1	D	161	ASP
1	D	173	GLN
1	D	237	GLU
1	D	247	PHE
1	D	250	LEU
1	D	293	ASP
1	D	326	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	HIS
1	A	110	HIS
1	A	118	GLN
1	A	140	GLN
1	A	167	GLN
1	A	173	GLN
1	A	230	ASN
1	A	291	ASN
1	A	328	GLN
1	A	388	GLN
1	A	467	GLN
1	B	118	GLN
1	B	173	GLN
1	B	288	GLN

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Mol	Chain	Res	Type
1	B	356	GLN
1	B	388	GLN
1	B	390	HIS
1	B	421	GLN
1	B	467	GLN
1	C	118	GLN
1	C	230	ASN
1	D	110	HIS
1	D	118	GLN
1	D	167	GLN
1	D	173	GLN
1	D	230	ASN
1	D	388	GLN
1	D	428	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 carbohydrates 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 ⓘ

### 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, 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	449/474 (94%)	-0.19	7 (1%) 74 75	6, 16, 41, 61	0
1	B	449/474 (94%)	-0.28	2 (0%) 93 93	7, 19, 35, 48	0
1	C	450/474 (94%)	-0.13	9 (2%) 68 69	7, 17, 43, 54	0
1	D	450/474 (94%)	-0.38	2 (0%) 93 93	5, 18, 34, 49	0
All	All	1798/1896 (94%)	-0.24	20 (1%) 82 83	5, 17, 39, 61	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	467	GLN	3.3
1	A	283	SER	3.1
1	C	82	VAL	3.0
1	D	80	VAL	2.9
1	C	65	ILE	2.7
1	C	468	ALA	2.6
1	A	465	LYS	2.5
1	C	80	VAL	2.4
1	A	412	LEU	2.4
1	A	404	GLY	2.3
1	C	70	GLU	2.3
1	C	282	GLY	2.3
1	C	83	VAL	2.3
1	B	284	SER	2.2
1	C	283	SER	2.2
1	D	19	THR	2.1
1	B	404	GLY	2.1
1	A	414	ASP	2.1
1	A	166	ALA	2.1
1	C	81	PHE	2.0

## 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 carbohydrates 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.