



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 10:06 am BST

PDB ID : 6KN0  
Title : caspase-1 P20/P10 C285A in complex with human GSDMD-C domain  
Authors : Ding, J.; Sun, Q.  
Deposited on : 2019-08-02  
Resolution : 2.79 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.

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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  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

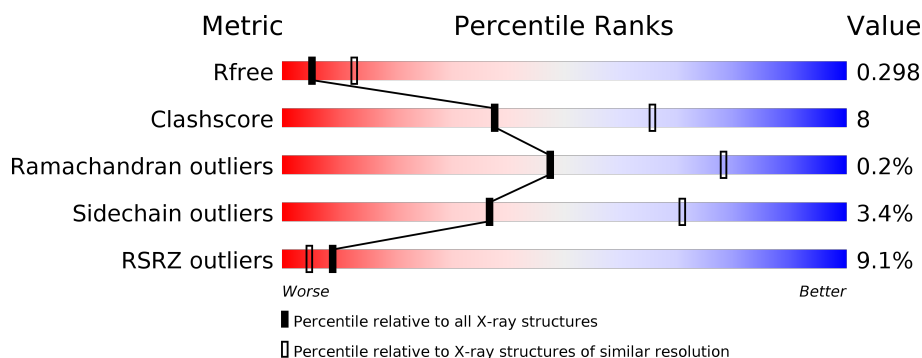
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.79 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	167	<div> <div>4%</div> <div>84%</div> <div>15%</div> <div>.</div> </div>
1	C	167	<div> <div>8%</div> <div>81%</div> <div>19%</div> </div>
2	B	88	<div> <div>8%</div> <div>81%</div> <div>18%</div> <div>.</div> </div>
2	D	88	<div> <div>9%</div> <div>68%</div> <div>32%</div> </div>
3	E	198	<div> <div>11%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>
3	F	198	<div> <div>13%</div> <div>79%</div> <div>17%</div> <div>..</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6974 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 Caspase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	167	Total	C	N	O	S	0	0	0
			1312	824	229	249	10			
1	C	167	Total	C	N	O	S	0	0	0
			1312	824	229	249	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	285	ALA	CYS	engineered mutation	UNP P29466
C	285	ALA	CYS	engineered mutation	UNP P29466

- Molecule 2 is a protein called Caspase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	88	Total	C	N	O	S	0	0	0
			716	456	126	127	7			
2	D	88	Total	C	N	O	S	0	0	0
			719	457	126	129	7			

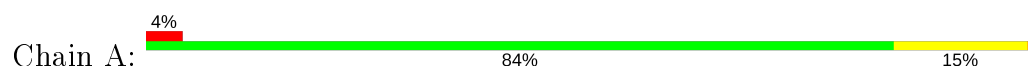
- Molecule 3 is a protein called Gasdermin-D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	195	Total	C	N	O	S	0	0	0
			1468	935	234	290	9			
3	F	192	Total	C	N	O	S	0	0	0
			1447	922	230	286	9			

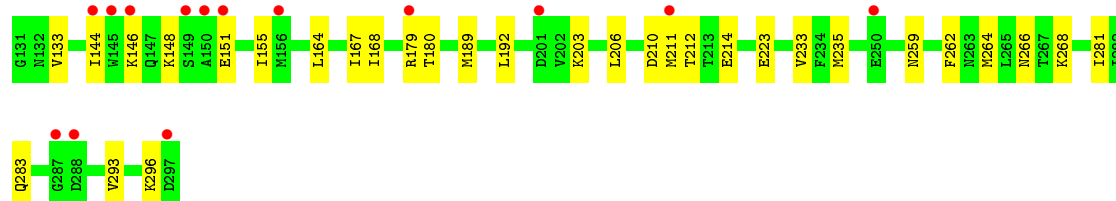
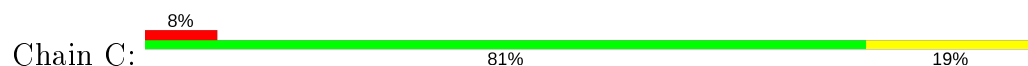
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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 ( $\text{RSRZ} > 2$ ). 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.

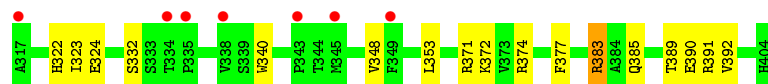
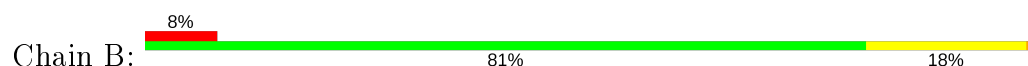
- Molecule 1: Caspase-1



- Molecule 1: Caspase-1



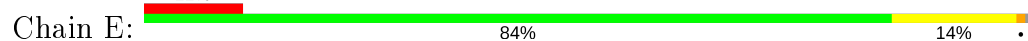
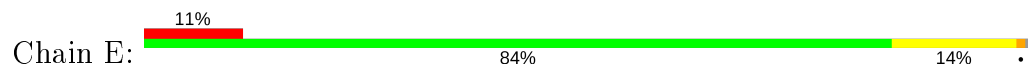
- Molecule 2: Caspase-1

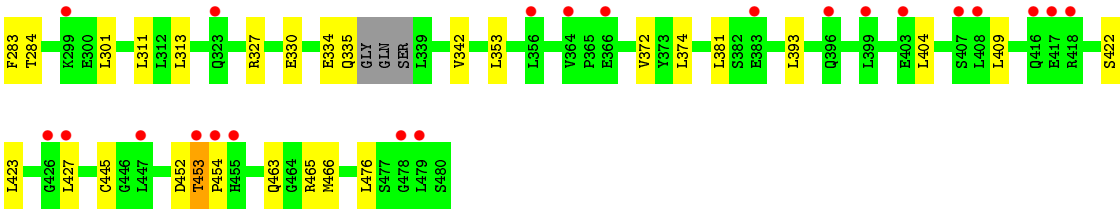


- Molecule 2: Caspase-1

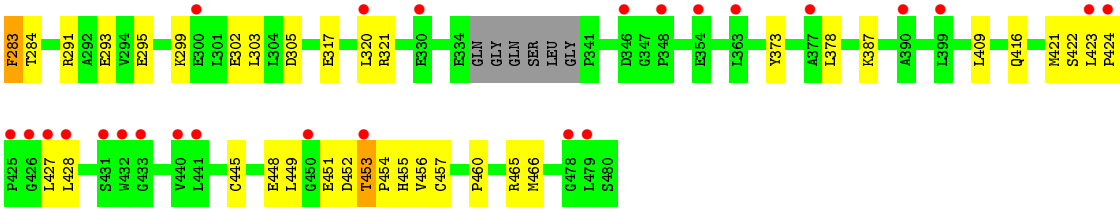
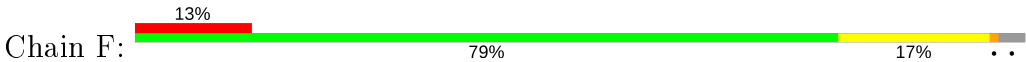


- Molecule 3: Gasdermin-D





● Molecule 3: Gasdermin-D



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	141.04Å 94.89Å 76.08Å 90.00° 116.43° 90.00°	Depositor
Resolution (Å)	47.45 – 2.79 47.45 – 2.79	Depositor EDS
% Data completeness (in resolution range)	97.3 (47.45-2.79) 97.3 (47.45-2.79)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.12 (at 2.81Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, $R_{free}$	0.240 , 0.300 0.240 , 0.298	Depositor DCC
$R_{free}$ test set	1998 reflections (9.16%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	67.3	Xtriage
Anisotropy	0.574	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 46.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	6974	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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	0.23	0/1334	0.42	0/1796
1	C	0.23	0/1334	0.41	0/1796
2	B	0.26	0/736	0.44	0/991
2	D	0.24	0/739	0.40	0/995
3	E	0.26	0/1492	0.44	0/2032
3	F	0.26	0/1471	0.44	0/2003
All	All	0.25	0/7106	0.43	0/9613

There are no bond length outliers.

There are no bond angle outliers.

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	1312	0	1331	21	0
1	C	1312	0	1331	22	0
2	B	716	0	691	27	0
2	D	719	0	693	24	0
3	E	1468	0	1477	17	0
3	F	1447	0	1456	26	0
All	All	6974	0	6979	107	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 8.

All (107) 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)
3:E:453:THR:HB	3:E:454:PRO:CD	1.71	1.19
3:F:453:THR:HG22	3:F:454:PRO:HD3	1.19	1.15
3:E:453:THR:HB	3:E:454:PRO:HD3	1.30	1.13
3:E:452:ASP:O	3:E:453:THR:OG1	1.69	1.07
2:B:340:TRP:CE2	2:B:383:ARG:HB3	1.91	1.05
2:B:383:ARG:HD2	2:B:383:ARG:H	1.26	1.00
2:B:340:TRP:CD2	2:B:383:ARG:HB3	2.00	0.96
3:E:453:THR:CB	3:E:454:PRO:CD	2.44	0.95
3:E:452:ASP:C	3:E:453:THR:HG1	1.72	0.92
3:F:453:THR:HG22	3:F:454:PRO:CD	1.98	0.92
3:F:422:SER:OG	3:F:454:PRO:HA	1.69	0.91
2:B:340:TRP:CD2	2:B:383:ARG:CB	2.59	0.84
2:B:340:TRP:CG	2:B:383:ARG:HB2	2.13	0.83
2:B:340:TRP:CD1	2:B:383:ARG:HB2	2.15	0.82
3:F:422:SER:OG	3:F:455:HIS:N	2.19	0.76
3:E:453:THR:HB	3:E:454:PRO:HD2	1.66	0.73
2:B:340:TRP:CG	2:B:383:ARG:CB	2.71	0.73
3:F:422:SER:OG	3:F:454:PRO:CA	2.38	0.72
2:B:340:TRP:CE2	2:B:383:ARG:CB	2.72	0.70
1:A:273:LEU:HA	1:A:276:LYS:HD2	1.78	0.65
1:C:212:THR:HG23	1:C:264:MET:HE1	1.77	0.65
1:C:148:LYS:HB3	1:C:151:GLU:HB2	1.80	0.63
3:E:335:GLN:HA	3:E:372:VAL:HG21	1.81	0.62
3:F:423:LEU:HB2	3:F:424:PRO:HD2	1.82	0.62
2:D:374:ARG:NH1	2:D:389:THR:OG1	2.32	0.61
2:B:383:ARG:H	2:B:383:ARG:CD	1.94	0.61
1:A:235:MET:HG2	1:A:283:GLN:HB3	1.82	0.59
1:A:172:GLU:HB3	1:A:178:ARG:HD3	1.84	0.59
3:F:299:LYS:HD2	3:F:302:GLU:HG3	1.85	0.58
2:B:374:ARG:NH2	2:D:394:LEU:O	2.37	0.57
2:B:391:ARG:H	2:D:390:GLU:HG2	1.70	0.57
3:F:422:SER:OG	3:F:454:PRO:C	2.43	0.57
3:E:327:ARG:NH1	3:E:330:GLU:OE1	2.38	0.56
1:A:286:ARG:NH2	2:B:390:GLU:OE2	2.28	0.56
2:D:334:THR:OG1	2:D:337:ASN:OD1	2.24	0.55
2:B:340:TRP:CD1	2:B:383:ARG:CB	2.89	0.55
3:E:404:LEU:HA	3:E:427:LEU:HD13	1.89	0.54
1:C:180:THR:O	2:D:341:ARG:NH2	2.41	0.54
1:A:294:TRP:HD1	3:F:303:LEU:HB3	1.71	0.54
3:F:320:LEU:HD21	3:F:378:LEU:HD13	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:133:VAL:HG13	2:D:358:GLN:HA	1.90	0.53
3:F:291:ARG:NH2	3:F:295:GLU:OE2	2.41	0.53
1:A:294:TRP:HE1	3:F:303:LEU:HD23	1.73	0.53
2:D:361:ALA:HA	2:D:369:ILE:HD11	1.91	0.53
3:F:453:THR:CG2	3:F:454:PRO:HD3	2.13	0.53
2:B:371:ARG:NH2	1:C:151:GLU:OE1	2.42	0.53
2:D:340:TRP:CG	2:D:383:ARG:HB3	2.44	0.52
2:B:372:LYS:NZ	1:C:151:GLU:OE2	2.41	0.52
3:F:409:LEU:HD21	3:F:466:MET:HB3	1.91	0.52
3:E:393:LEU:HD21	3:E:476:LEU:HD13	1.92	0.51
3:E:453:THR:CB	3:E:454:PRO:HD2	2.33	0.51
1:C:268:LYS:HG3	3:F:305:ASP:HA	1.93	0.51
3:F:453:THR:CG2	3:F:454:PRO:CD	2.83	0.51
3:E:422:SER:HB3	3:E:454:PRO:HA	1.93	0.51
1:C:281:ILE:HD13	2:D:353:LEU:HD21	1.93	0.50
2:B:377:PHE:O	2:D:322:HIS:NE2	2.43	0.49
1:C:179:ARG:HH21	2:D:341:ARG:HD3	1.78	0.49
3:E:409:LEU:HD21	3:E:466:MET:HB3	1.95	0.48
1:C:164:LEU:HD11	1:C:223:GLU:HB2	1.94	0.48
1:C:168:ILE:HG21	1:C:211:MET:HG2	1.96	0.48
3:E:445:CYS:HA	3:E:465:ARG:HD2	1.96	0.47
1:C:235:MET:HG2	1:C:283:GLN:HB3	1.96	0.47
1:A:293:VAL:HG21	1:C:268:LYS:HB2	1.97	0.47
1:A:165:ALA:HB1	1:A:193:LEU:HD13	1.96	0.47
1:C:192:LEU:HD22	2:D:354:ILE:HG12	1.97	0.47
1:A:250:GLU:OE1	1:A:250:GLU:N	2.41	0.46
1:A:151:GLU:OE2	2:D:372:LYS:NZ	2.48	0.46
1:A:281:ILE:HD13	2:B:353:LEU:HD21	1.97	0.46
1:A:283:GLN:HA	2:B:332:SER:HB3	1.98	0.46
1:A:266:ASN:HB2	2:B:323:ILE:O	2.16	0.46
3:F:283:PHE:HB3	3:F:284:THR:H	1.53	0.46
3:F:449:LEU:HD22	3:F:456:VAL:HG22	1.96	0.46
1:C:266:ASN:HB2	2:D:323:ILE:O	2.15	0.46
2:B:371:ARG:HE	2:D:396:ARG:HA	1.79	0.46
2:D:348:VAL:HG21	2:D:383:ARG:HH22	1.80	0.46
1:C:144:ILE:HG12	2:D:399:TYR:CZ	2.51	0.45
1:A:259:ASN:HB3	2:B:391:ARG:HH22	1.81	0.45
2:B:322:HIS:NE2	2:D:377:PHE:O	2.49	0.45
1:C:262:PHE:HZ	2:D:329:ALA:HB2	1.80	0.45
3:E:381:LEU:HD23	3:E:465:ARG:HG3	1.99	0.45
3:F:445:CYS:HA	3:F:465:ARG:HG2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:371:ARG:HB3	2:B:371:ARG:HH11	1.83	0.44
1:C:203:LYS:NZ	1:C:214:GLU:OE1	2.33	0.44
1:A:164:LEU:HD23	1:A:164:LEU:HA	1.87	0.44
2:B:392:VAL:HB	2:D:389:THR:HB	2.00	0.44
1:C:206:LEU:HB3	1:C:210:ASP:HB2	2.01	0.43
3:F:416:GLN:HA	3:F:460:PRO:HB3	1.99	0.43
2:D:340:TRP:CD1	2:D:383:ARG:HB3	2.53	0.43
1:A:135:LEU:H	1:A:135:LEU:HD23	1.84	0.43
2:B:323:ILE:HG23	2:B:324:GLU:HG2	2.00	0.42
1:C:296:LYS:HE2	3:E:311:LEU:HD11	2.01	0.42
2:B:348:VAL:HB	2:B:385:GLN:NE2	2.35	0.42
3:E:342:VAL:HG11	3:E:353:LEU:HD11	2.00	0.42
3:F:387:LYS:HE3	3:F:387:LYS:HB2	1.89	0.42
1:A:164:LEU:HD12	1:A:224:HIS:CE1	2.55	0.42
1:C:155:ILE:HG13	2:D:401:PHE:HD2	1.85	0.42
1:A:204:LYS:HB3	1:A:204:LYS:HE2	1.92	0.42
1:A:211:MET:HB2	1:A:211:MET:HE3	1.96	0.42
3:F:293:GLU:OE2	3:F:373:TYR:OH	2.33	0.41
1:A:134:LYS:HD2	1:A:134:LYS:HA	1.91	0.41
1:C:167:ILE:HG12	1:C:233:VAL:HB	2.01	0.41
3:F:448:GLU:HB2	3:F:457:CYS:SG	2.61	0.41
2:B:389:THR:HB	2:D:392:VAL:HB	2.03	0.41
2:D:365:ASP:HB3	2:D:399:TYR:CE1	2.56	0.41
1:A:294:TRP:NE1	3:F:303:LEU:HD23	2.35	0.41
3:F:427:LEU:HA	3:F:427:LEU:HD23	1.94	0.40
3:F:317:GLU:O	3:F:321:ARG: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	165/167 (99%)	152 (92%)	13 (8%)	0	100	100
1	C	165/167 (99%)	156 (94%)	9 (6%)	0	100	100
2	B	86/88 (98%)	83 (96%)	3 (4%)	0	100	100
2	D	86/88 (98%)	84 (98%)	2 (2%)	0	100	100
3	E	191/198 (96%)	187 (98%)	3 (2%)	1 (0%)	29	61
3	F	188/198 (95%)	182 (97%)	5 (3%)	1 (0%)	29	61
All	All	881/906 (97%)	844 (96%)	35 (4%)	2 (0%)	47	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	453	THR
3	F	453	THR

### 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	148/148 (100%)	143 (97%)	5 (3%)	37	71
1	C	148/148 (100%)	144 (97%)	4 (3%)	44	78
2	B	78/79 (99%)	77 (99%)	1 (1%)	69	91
2	D	79/79 (100%)	76 (96%)	3 (4%)	33	67
3	E	160/162 (99%)	152 (95%)	8 (5%)	24	56
3	F	158/162 (98%)	153 (97%)	5 (3%)	39	73
All	All	771/778 (99%)	745 (97%)	26 (3%)	37	71

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	135	LEU
1	A	176	ILE
1	A	191	MET

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Mol	Chain	Res	Type
1	A	258	LEU
1	A	296	LYS
2	B	383	ARG
1	C	146	LYS
1	C	189	MET
1	C	259	ASN
1	C	293	VAL
2	D	318	ILE
2	D	378	GLU
2	D	381	ASP
3	E	283	PHE
3	E	284	THR
3	E	301	LEU
3	E	313	LEU
3	E	334	GLU
3	E	374	LEU
3	E	423	LEU
3	E	463	GLN
3	F	283	PHE
3	F	421	MET
3	F	428	LEU
3	F	451	GLU
3	F	452	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
3	F	455	HIS

### 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 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	167/167 (100%)	0.59	6 (3%) 42 32	43, 72, 98, 115	0
1	C	167/167 (100%)	0.68	14 (8%) 11 5	41, 72, 105, 119	0
2	B	88/88 (100%)	0.61	7 (7%) 12 6	42, 61, 106, 131	0
2	D	88/88 (100%)	0.70	8 (9%) 9 5	42, 63, 118, 128	0
3	E	195/198 (98%)	0.82	22 (11%) 5 3	51, 74, 111, 134	0
3	F	192/198 (96%)	0.93	25 (13%) 3 2	50, 74, 116, 137	0
All	All	897/906 (99%)	0.74	82 (9%) 9 5	41, 71, 109, 137	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	150	ALA	7.2
3	E	403	GLU	5.5
3	F	423	LEU	5.2
3	E	426	GLY	4.8
2	D	340	TRP	4.3
2	B	335	PRO	4.3
3	F	453	THR	4.1
3	F	440	VAL	3.7
3	E	356	LEU	3.7
3	F	399	LEU	3.6
2	B	345	MET	3.6
1	C	288	ASP	3.6
3	F	433	GLY	3.5
2	B	317	ALA	3.4
2	B	334	THR	3.3
3	F	390	ALA	3.2
3	E	416	GLN	3.2
1	C	287	GLY	3.1
3	F	479	LEU	3.1

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	201	ASP	3.1
3	E	323	GLN	3.1
1	C	156	MET	3.1
1	A	241	GLU	3.0
3	F	427	LEU	3.0
3	E	417	GLU	3.0
2	B	343	PRO	3.0
3	F	441	LEU	3.0
3	F	424	PRO	3.0
1	C	297	ASP	2.9
3	F	428	LEU	2.9
3	E	478	GLY	2.9
1	C	201	ASP	2.9
3	F	431	SER	2.8
2	D	404	HIS	2.8
1	A	297	ASP	2.8
3	F	346	ASP	2.8
1	C	145	TRP	2.8
3	E	299	LYS	2.8
1	C	211	MET	2.8
3	F	348	PRO	2.7
3	E	383	GLU	2.6
3	F	377	ALA	2.6
3	F	426	GLY	2.6
3	F	450	GLY	2.6
3	E	418	ARG	2.6
3	F	363	LEU	2.5
3	E	455	HIS	2.5
1	C	146	LYS	2.5
3	E	399	LEU	2.5
3	E	366	GLU	2.4
2	D	381	ASP	2.4
2	D	334	THR	2.4
3	F	354	GLU	2.4
1	C	179	ARG	2.4
3	E	427	LEU	2.4
3	F	320	LEU	2.4
1	A	250	GLU	2.4
3	E	407	SER	2.3
3	F	425	PRO	2.3
3	F	300	GLU	2.3
1	C	151	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
2	D	336	ASP	2.3
1	C	149	SER	2.2
1	C	144	ILE	2.2
1	C	250	GLU	2.2
3	E	479	LEU	2.2
3	E	453	THR	2.2
3	F	432	TRP	2.2
2	D	388	THR	2.2
3	E	454	PRO	2.2
1	A	231	PHE	2.2
2	D	341	ARG	2.1
3	E	364	VAL	2.1
2	B	349	PHE	2.1
1	A	216	GLU	2.1
2	D	335	PRO	2.0
2	B	338	VAL	2.0
3	F	330	GLU	2.0
3	F	478	GLY	2.0
3	E	396	GLN	2.0
3	E	408	LEU	2.0
3	E	447	LEU	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.