



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:37 AM GMT

PDB ID : 2NN3  
Title : structure of pro-sf-caspase-1  
Authors : Fisher, A.J.; Ni, L.  
Deposited on : 2006-10-23  
Resolution : 3.00 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

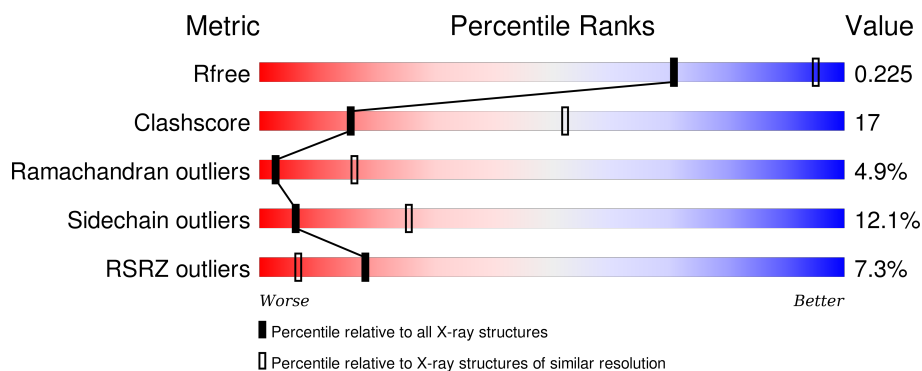
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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	C	310	 2% 44% 27% 5% • 24%
1	D	310	 9% 49% 27% 6% • 16%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3872 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	C	235	Total	C	N	O	S	0	0	0
			1852	1186	314	338	14			
1	D	260	Total	C	N	O	S	0	0	0
			2014	1281	345	373	15			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	136	ALA	HIS	ENGINEERED	UNP P89116
C	300	GLY	-	EXPRESSION TAG	UNP P89116
C	301	GLY	-	EXPRESSION TAG	UNP P89116
C	302	GLY	-	EXPRESSION TAG	UNP P89116
C	303	LEU	-	EXPRESSION TAG	UNP P89116
C	304	GLU	-	EXPRESSION TAG	UNP P89116
C	305	HIS	-	EXPRESSION TAG	UNP P89116
C	306	HIS	-	EXPRESSION TAG	UNP P89116
C	307	HIS	-	EXPRESSION TAG	UNP P89116
C	308	HIS	-	EXPRESSION TAG	UNP P89116
C	309	HIS	-	EXPRESSION TAG	UNP P89116
C	310	HIS	-	EXPRESSION TAG	UNP P89116
D	136	ALA	HIS	ENGINEERED	UNP P89116
D	300	GLY	-	EXPRESSION TAG	UNP P89116
D	301	GLY	-	EXPRESSION TAG	UNP P89116
D	302	GLY	-	EXPRESSION TAG	UNP P89116
D	303	LEU	-	EXPRESSION TAG	UNP P89116
D	304	GLU	-	EXPRESSION TAG	UNP P89116
D	305	HIS	-	EXPRESSION TAG	UNP P89116
D	306	HIS	-	EXPRESSION TAG	UNP P89116
D	307	HIS	-	EXPRESSION TAG	UNP P89116
D	308	HIS	-	EXPRESSION TAG	UNP P89116
D	309	HIS	-	EXPRESSION TAG	UNP P89116
D	310	HIS	-	EXPRESSION TAG	UNP P89116

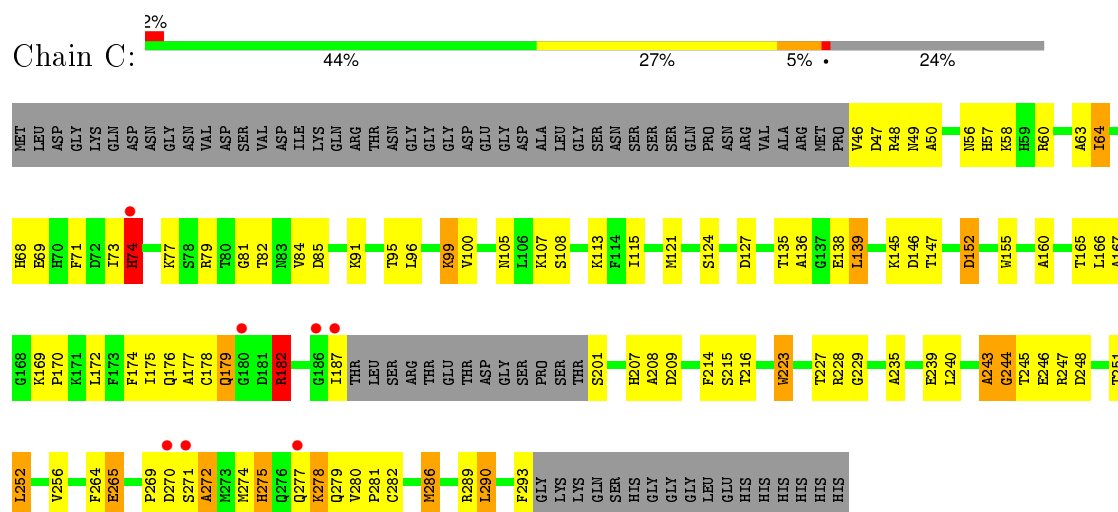
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	3	Total 3	O 3	0	0
2	D	3	Total 3	O 3	0	0

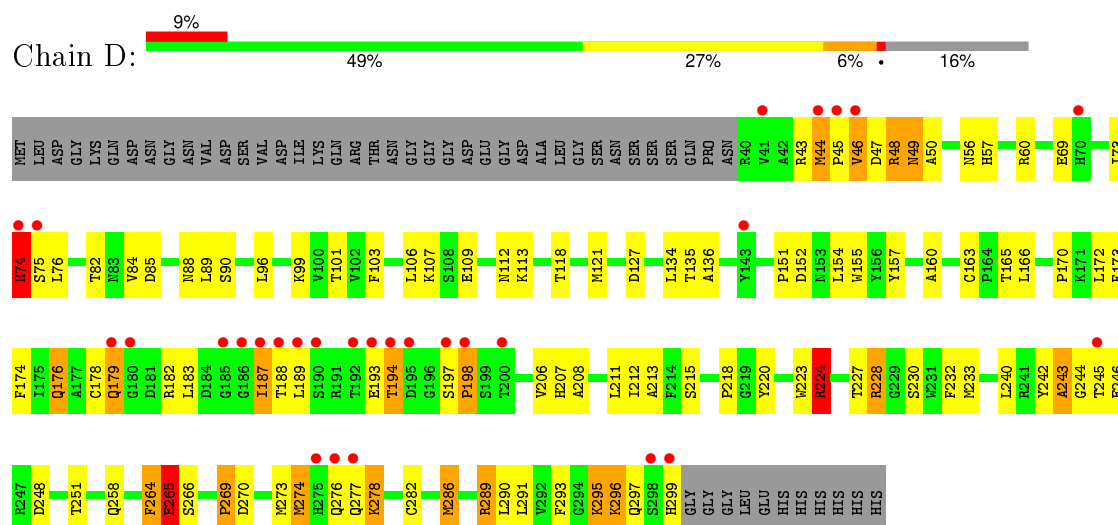
### 3 Residue-property plots

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 ( $\text{RSRZ} > 2$ ). 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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	106.22Å 106.22Å 113.59Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.34 – 3.00 48.33 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.5 (48.34-3.00) 97.6 (48.33-3.00)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.55 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.228 , 0.275 0.228 , 0.225	Depositor DCC
$R_{free}$ test set	721 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	108.1	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 135.2	EDS
Estimated twinning fraction	0.049 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 14298 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3872	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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	C	0.87	5/1897 (0.3%)	0.77	2/2572 (0.1%)
1	D	1.04	7/2061 (0.3%)	1.03	5/2798 (0.2%)
All	All	0.96	12/3958 (0.3%)	0.92	7/5370 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	43	ARG	CZ-NH1	22.03	1.61	1.33
1	D	224	ARG	CZ-NH1	16.26	1.54	1.33
1	C	74	HIS	CE1-NE2	15.16	1.67	1.32
1	C	187	ILE	C-O	14.16	1.50	1.23
1	D	224	ARG	NE-CZ	13.32	1.50	1.33
1	D	74	HIS	CE1-NE2	11.74	1.59	1.32
1	C	74	HIS	CG-ND1	9.39	1.59	1.38
1	D	74	HIS	CG-ND1	9.06	1.58	1.38
1	C	74	HIS	CG-CD2	6.60	1.47	1.35
1	D	224	ARG	CZ-NH2	5.80	1.40	1.33
1	D	74	HIS	CG-CD2	5.22	1.44	1.35
1	C	74	HIS	CD2-NE2	-5.17	1.26	1.38

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	224	ARG	NE-CZ-NH2	-23.37	108.61	120.30
1	D	43	ARG	NE-CZ-NH2	-16.18	112.21	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	43	ARG	NE-CZ-NH1	15.10	127.85	120.30
1	D	224	ARG	NE-CZ-NH1	15.02	127.81	120.30
1	D	290	LEU	CA-CB-CG	6.72	130.76	115.30
1	C	215	SER	N-CA-C	5.21	125.07	111.00
1	C	139	LEU	CA-CB-CG	5.12	127.09	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	224	ARG	Sidechain

## 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	C	1852	0	1797	75	1
1	D	2014	0	1943	71	1
2	C	3	0	0	0	0
2	D	3	0	0	0	0
All	All	3872	0	3740	131	1

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 17.

All (131) 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:C:74:HIS:CE1	1:C:74:HIS:NE2	1.67	1.43
1:D:44:MET:HB3	1:D:45:PRO:HD3	1.16	1.11
1:C:167:ALA:HB1	1:D:276:GLN:CB	1.90	1.02
1:C:282:CYS:HA	1:D:286:MET:HE2	1.38	1.01
1:D:44:MET:CB	1:D:45:PRO:HD3	1.96	0.94
1:C:286:MET:HE2	1:D:282:CYS:HA	1.50	0.93
1:D:44:MET:HB3	1:D:45:PRO:CD	1.96	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:57:HIS:HD2	1:D:127:ASP:OD1	1.54	0.89
1:D:84:VAL:HG21	1:D:228:ARG:HB2	1.53	0.87
1:D:273:MET:HG3	1:D:274:MET:H	1.38	0.87
1:C:182:ARG:HH11	1:C:182:ARG:HA	1.53	0.73
1:C:46:VAL:HG13	1:C:47:ASP:H	1.53	0.72
1:D:178:CYS:O	1:D:179:GLN:HG2	1.91	0.71
1:C:277:GLN:HB3	1:D:207:HIS:CE1	2.27	0.70
1:D:57:HIS:CD2	1:D:127:ASP:OD1	2.43	0.70
1:D:112:ASN:HD21	1:D:157:TYR:HE1	1.40	0.70
1:D:152:ASP:HA	1:D:155:TRP:CE2	2.27	0.68
1:D:45:PRO:O	1:D:46:VAL:HB	1.92	0.68
1:C:182:ARG:HH21	1:D:206:VAL:HB	1.59	0.68
1:C:74:HIS:ND1	1:C:74:HIS:N	2.45	0.65
1:C:248:ASP:OD2	1:C:251:THR:HB	1.97	0.65
1:C:48:ARG:O	1:C:289:ARG:NH2	2.29	0.64
1:C:68:HIS:HD2	1:C:135:THR:HB	1.62	0.64
1:C:269:PRO:HG3	1:C:275:HIS:CE1	2.32	0.64
1:C:223:TRP:HB2	1:C:228:ARG:HD3	1.80	0.64
1:C:286:MET:CE	1:D:282:CYS:HA	2.27	0.61
1:C:79:ARG:NH2	1:C:82:THR:OG1	2.33	0.61
1:C:50:ALA:HB3	1:C:289:ARG:NH1	2.16	0.60
1:D:48:ARG:O	1:D:289:ARG:NH2	2.34	0.60
1:D:187:ILE:HG22	1:D:187:ILE:O	2.01	0.60
1:D:242:TYR:O	1:D:243:ALA:HB2	2.02	0.60
1:C:152:ASP:HA	1:C:155:TRP:CE2	2.37	0.59
1:C:85:ASP:OD1	1:C:229:GLY:HA2	2.02	0.59
1:C:107:LYS:HB3	1:C:146:ASP:OD2	2.01	0.59
1:C:57:HIS:CD2	1:C:127:ASP:OD1	2.55	0.59
1:D:273:MET:HG3	1:D:274:MET:HG2	1.85	0.59
1:D:103:PHE:HB3	1:D:106:LEU:CD1	2.33	0.58
1:C:282:CYS:HA	1:D:286:MET:CE	2.23	0.58
1:D:174:PHE:HD2	1:D:213:ALA:HB3	1.69	0.57
1:C:248:ASP:OD2	1:C:251:THR:CB	2.54	0.56
1:D:49:ASN:N	1:D:49:ASN:OD1	2.38	0.56
1:D:103:PHE:HB3	1:D:106:LEU:HD11	1.86	0.56
1:D:174:PHE:CD2	1:D:213:ALA:HB3	2.40	0.56
1:C:239:GLU:O	1:C:243:ALA:HB3	2.07	0.55
1:D:248:ASP:OD2	1:D:251:THR:HB	2.07	0.55
1:C:178:CYS:O	1:C:179:GLN:HG2	2.07	0.55
1:C:49:ASN:HA	1:D:258:GLN:NE2	2.22	0.54
1:D:118:THR:O	1:D:121:MET:HG3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:277:GLN:HG2	1:D:208:ALA:HB3	1.90	0.53
1:C:146:ASP:OD1	1:C:147:THR:N	2.40	0.53
1:C:170:PRO:HB2	1:C:172:LEU:HD11	1.89	0.53
1:D:176:GLN:HE22	1:D:230:SER:CB	2.21	0.53
1:C:277:GLN:CG	1:D:208:ALA:HB3	2.38	0.53
1:D:163:CYS:SG	1:D:166:LEU:HD12	2.49	0.53
1:C:290:LEU:HD21	1:D:251:THR:HG23	1.91	0.53
1:D:44:MET:CB	1:D:45:PRO:CD	2.72	0.52
1:C:178:CYS:O	1:C:179:GLN:CG	2.57	0.52
1:C:105:ASN:HB3	1:C:145:LYS:HD3	1.92	0.52
1:C:68:HIS:CD2	1:C:135:THR:HB	2.42	0.51
1:D:135:THR:HG23	1:D:136:ALA:O	2.10	0.51
1:C:269:PRO:HG3	1:C:275:HIS:HE1	1.77	0.50
1:D:73:ILE:HG22	1:D:74:HIS:O	2.11	0.50
1:C:280:VAL:HB	1:C:281:PRO:HD2	1.93	0.50
1:C:166:LEU:HA	1:C:169:LYS:HD2	1.91	0.50
1:C:57:HIS:CD2	1:C:57:HIS:N	2.78	0.50
1:D:188:THR:HG22	1:D:189:LEU:H	1.77	0.50
1:D:69:GLU:HG3	1:D:82:THR:HG22	1.93	0.50
1:C:175:ILE:HG21	1:C:214:PHE:CE1	2.47	0.49
1:C:228:ARG:HG3	1:C:228:ARG:O	2.13	0.48
1:C:179:GLN:HA	1:C:179:GLN:OE1	2.13	0.48
1:C:172:LEU:N	1:C:172:LEU:HD12	2.29	0.48
1:C:64:ILE:HD11	1:C:115:ILE:HD13	1.95	0.48
1:C:264:PHE:CE2	1:C:279:GLN:HB3	2.49	0.48
1:C:84:VAL:HG21	1:C:228:ARG:HB2	1.95	0.48
1:C:240:LEU:HD23	1:C:252:LEU:CD2	2.43	0.48
1:C:69:GLU:HG3	1:C:79:ARG:O	2.14	0.48
1:D:215:SER:HB2	1:D:232:PHE:CD2	2.48	0.48
1:C:146:ASP:OD1	1:C:147:THR:HG22	2.15	0.47
1:D:178:CYS:O	1:D:179:GLN:CG	2.60	0.47
1:C:135:THR:O	1:C:177:ALA:HA	2.14	0.47
1:C:99:LYS:HE3	1:C:100:VAL:N	2.30	0.47
1:D:155:TRP:CZ3	1:D:212:ILE:HD13	2.50	0.47
1:C:121:MET:O	1:C:165:THR:HG21	2.14	0.47
1:C:108:SER:H	1:C:146:ASP:CG	2.19	0.47
1:D:223:TRP:HB2	1:D:228:ARG:HD3	1.97	0.47
1:D:295:LYS:HE3	1:D:296:LYS:H	1.79	0.47
1:C:282:CYS:CA	1:D:286:MET:HE2	2.26	0.46
1:D:264:PHE:CD1	1:D:278:LYS:CB	2.99	0.46
1:D:248:ASP:OD2	1:D:251:THR:CB	2.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:178:CYS:O	1:C:179:GLN:CD	2.55	0.46
1:D:96:LEU:HD11	1:D:240:LEU:HD22	1.97	0.46
1:C:274:MET:O	1:C:275:HIS:HB3	2.17	0.45
1:C:240:LEU:HD23	1:C:252:LEU:HD21	1.99	0.45
1:D:193:GLU:HG2	1:D:194:THR:H	1.81	0.45
1:C:235:ALA:HB3	1:C:256:VAL:HG13	1.99	0.45
1:D:74:HIS:ND1	1:D:74:HIS:N	2.64	0.45
1:C:264:PHE:O	1:C:265:GLU:HB3	2.17	0.45
1:C:277:GLN:H	1:D:160:ALA:HB3	1.82	0.44
1:C:271:SER:O	1:C:272:ALA:HB3	2.16	0.44
1:D:197:SER:N	1:D:198:PRO:HD3	2.32	0.44
1:C:139:LEU:HD23	1:C:201:SER:OG	2.17	0.44
1:D:295:LYS:HG3	1:D:296:LYS:H	1.82	0.44
1:C:209:ASP:OD2	1:D:277:GLN:HG3	2.17	0.44
1:C:160:ALA:H	1:C:207:HIS:CE1	2.35	0.44
1:D:273:MET:HG3	1:D:274:MET:N	2.20	0.44
1:C:277:GLN:HB2	1:D:160:ALA:HB2	2.00	0.44
1:C:247:ARG:HB2	1:C:252:LEU:CD1	2.47	0.44
1:C:96:LEU:HA	1:C:96:LEU:HD23	1.84	0.43
1:D:47:ASP:HB3	1:D:49:ASN:H	1.83	0.43
1:D:244:GLY:O	1:D:293:PHE:HB2	2.18	0.43
1:D:218:PRO:O	1:D:220:TYR:CE1	2.71	0.43
1:C:91:LYS:HE3	1:C:91:LYS:HB2	1.86	0.43
1:C:63:ALA:O	1:C:100:VAL:HA	2.18	0.43
1:D:85:ASP:OD1	1:D:233:MET:HB2	2.18	0.43
1:D:274:MET:HG2	1:D:274:MET:H	1.66	0.43
1:C:174:PHE:C	1:C:175:ILE:HD12	2.39	0.43
1:D:269:PRO:HB2	1:D:270:ASP:H	1.69	0.43
1:C:135:THR:OG1	1:C:136:ALA:N	2.50	0.42
1:D:50:ALA:HB3	1:D:289:ARG:NH1	2.34	0.42
1:D:297:GLN:C	1:D:299:HIS:H	2.22	0.42
1:C:81:GLY:HA3	1:C:228:ARG:HA	2.01	0.42
1:D:151:PRO:O	1:D:154:LEU:HB2	2.20	0.42
1:D:170:PRO:HG3	1:D:291:LEU:HD13	2.01	0.42
1:C:71:PHE:C	1:C:73:ILE:H	2.24	0.41
1:C:244:GLY:O	1:C:293:PHE:HB2	2.20	0.41
1:D:173:PHE:CD1	1:D:173:PHE:N	2.89	0.41
1:D:85:ASP:HB3	1:D:134:LEU:HD21	2.02	0.41
1:C:177:ALA:O	1:C:216:THR:HG22	2.21	0.40
1:D:89:LEU:HD12	1:D:89:LEU:HA	1.86	0.40
1:D:264:PHE:O	1:D:265:GLU:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:208:ALA:O	1:C:209:ASP:HB2	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:223:TRP:O	1:D:218:PRO:O[5_555]	2.16	0.04

## 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	C	231/310 (74%)	194 (84%)	26 (11%)	11 (5%)	3	17
1	D	258/310 (83%)	203 (79%)	42 (16%)	13 (5%)	3	15
All	All	489/620 (79%)	397 (81%)	68 (14%)	24 (5%)	3	16

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	265	GLU
1	C	278	LYS
1	D	44	MET
1	D	46	VAL
1	D	243	ALA
1	D	265	GLU
1	D	269	PRO
1	C	58	LYS
1	C	272	ALA
1	C	275	HIS
1	D	228	ARG
1	D	296	LYS

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Mol	Chain	Res	Type
1	C	182	ARG
1	C	243	ALA
1	D	183	LEU
1	C	56	ASN
1	D	179	GLN
1	D	264	PHE
1	C	179	GLN
1	C	244	GLY
1	D	278	LYS
1	C	223	TRP
1	D	187	ILE
1	D	198	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	C	199/267 (74%)	179 (90%)	20 (10%)	9	34
1	D	214/267 (80%)	184 (86%)	30 (14%)	4	19
All	All	413/534 (77%)	363 (88%)	50 (12%)	6	25

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	60	ARG
1	C	64	ILE
1	C	74	HIS
1	C	77	LYS
1	C	95	THR
1	C	99	LYS
1	C	113	LYS
1	C	124	SER
1	C	138	GLU
1	C	152	ASP
1	C	176	GLN

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Mol	Chain	Res	Type
1	C	182	ARG
1	C	227	THR
1	C	245	THR
1	C	246	GLU
1	C	252	LEU
1	C	270	ASP
1	C	278	LYS
1	C	286	MET
1	C	290	LEU
1	D	48	ARG
1	D	49	ASN
1	D	56	ASN
1	D	60	ARG
1	D	74	HIS
1	D	75	SER
1	D	76	LEU
1	D	88	ASN
1	D	90	SER
1	D	99	LYS
1	D	101	THR
1	D	107	LYS
1	D	109	GLU
1	D	113	LYS
1	D	165	THR
1	D	172	LEU
1	D	176	GLN
1	D	182	ARG
1	D	194	THR
1	D	211	LEU
1	D	224	ARG
1	D	227	THR
1	D	245	THR
1	D	246	GLU
1	D	265	GLU
1	D	266	SER
1	D	274	MET
1	D	286	MET
1	D	289	ARG
1	D	295	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	57	HIS
1	C	67	ASN
1	C	68	HIS
1	C	153	ASN
1	C	176	GLN
1	C	207	HIS
1	C	279	GLN
1	D	57	HIS
1	D	59	HIS
1	D	68	HIS
1	D	88	ASN
1	D	112	ASN
1	D	153	ASN
1	D	176	GLN
1	D	258	GLN
1	D	279	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	C	235/310 (75%)	-0.06	7 (2%) 54 25	84, 101, 123, 138	0
1	D	260/310 (83%)	0.39	29 (11%) 7 3	82, 101, 137, 141	0
All	All	495/620 (79%)	0.17	36 (7%) 18 6	82, 101, 134, 141	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	276	GLN	5.9
1	D	193	GLU	5.8
1	D	192	THR	4.6
1	D	188	THR	4.5
1	D	180	GLY	4.0
1	C	270	ASP	3.9
1	D	185	GLY	3.7
1	D	277	GLN	3.7
1	D	46	VAL	3.7
1	D	299	HIS	3.6
1	D	198	PRO	3.6
1	D	275	HIS	3.3
1	D	189	LEU	3.2
1	D	186	GLY	3.0
1	D	200	THR	3.0
1	D	179	GLN	2.9
1	D	44	MET	2.8
1	D	194	THR	2.8
1	D	45	PRO	2.7
1	D	190	SER	2.7
1	C	74	HIS	2.7
1	D	197	SER	2.6
1	D	195	ASP	2.6
1	D	41	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	277	GLN	2.6
1	C	186	GLY	2.5
1	D	187	ILE	2.4
1	D	70	HIS	2.4
1	C	187	ILE	2.3
1	D	74	HIS	2.3
1	D	75	SER	2.3
1	C	180	GLY	2.3
1	D	245	THR	2.2
1	D	143	TYR	2.2
1	C	271	SER	2.1
1	D	298	SER	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.