



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:58 AM GMT

PDB ID : 2F1M  
Title : Conformational flexibility in the multidrug efflux system protein AcrA  
Authors : Mikolosko, J.; Bobyk, K.; Zgurskaya, H.I.; Ghosh, P.  
Deposited on : 2005-11-14  
Resolution : 2.71 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

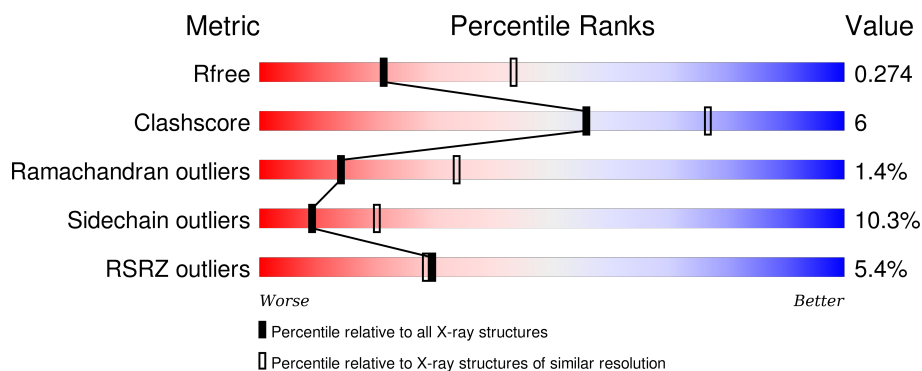
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.71 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	2439 (2.74-2.70)
Clashscore	102246	2771 (2.74-2.70)
Ramachandran outliers	100387	2726 (2.74-2.70)
Sidechain outliers	100360	2727 (2.74-2.70)
RSRZ outliers	91569	2443 (2.74-2.70)

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	277	
1	B	277	
1	C	277	
1	D	277	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7000 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 Acriflavine resistance protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	Se	20	0	0
			1728	1076	300	348	4			
1	B	223	Total	C	N	O	Se	25	0	0
			1680	1049	292	336	3			
1	C	247	Total	C	N	O	Se	0	0	0
			1870	1161	327	377	5			
1	D	224	Total	C	N	O	Se	21	0	0
			1686	1053	294	336	3			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	44	MSE	-	INITIATING METHIONINE	UNP P0AE06
A	223	MSE	PHE	ENGINEERED	UNP P0AE06
A	224	MSE	LEU	ENGINEERED	UNP P0AE06
A	287	MSE	LEU	ENGINEERED	UNP P0AE06
A	288	MSE	LEU	ENGINEERED	UNP P0AE06
A	291	MSE	MET	MODIFIED RESIDUE	UNP P0AE06
A	313	LEU	-	CLONING ARTIFACT	UNP P0AE06
A	314	GLU	-	CLONING ARTIFACT	UNP P0AE06
A	315	HIS	-	CLONING ARTIFACT	UNP P0AE06
A	316	HIS	-	CLONING ARTIFACT	UNP P0AE06
A	317	HIS	-	CLONING ARTIFACT	UNP P0AE06
A	318	HIS	-	CLONING ARTIFACT	UNP P0AE06
A	319	HIS	-	CLONING ARTIFACT	UNP P0AE06
A	320	HIS	-	CLONING ARTIFACT	UNP P0AE06
B	44	MSE	-	INITIATING METHIONINE	UNP P0AE06
B	223	MSE	PHE	ENGINEERED	UNP P0AE06
B	224	MSE	LEU	ENGINEERED	UNP P0AE06
B	287	MSE	LEU	ENGINEERED	UNP P0AE06
B	288	MSE	LEU	ENGINEERED	UNP P0AE06
B	291	MSE	MET	MODIFIED RESIDUE	UNP P0AE06
B	313	LEU	-	CLONING ARTIFACT	UNP P0AE06

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Chain	Residue	Modelled	Actual	Comment	Reference
B	314	GLU	-	CLONING ARTIFACT	UNP P0AE06
B	315	HIS	-	CLONING ARTIFACT	UNP P0AE06
B	316	HIS	-	CLONING ARTIFACT	UNP P0AE06
B	317	HIS	-	CLONING ARTIFACT	UNP P0AE06
B	318	HIS	-	CLONING ARTIFACT	UNP P0AE06
B	319	HIS	-	CLONING ARTIFACT	UNP P0AE06
B	320	HIS	-	CLONING ARTIFACT	UNP P0AE06
C	44	MSE	-	INITIATING METHIONINE	UNP P0AE06
C	223	MSE	PHE	ENGINEERED	UNP P0AE06
C	224	MSE	LEU	ENGINEERED	UNP P0AE06
C	287	MSE	LEU	ENGINEERED	UNP P0AE06
C	288	MSE	LEU	ENGINEERED	UNP P0AE06
C	291	MSE	MET	MODIFIED RESIDUE	UNP P0AE06
C	313	LEU	-	CLONING ARTIFACT	UNP P0AE06
C	314	GLU	-	CLONING ARTIFACT	UNP P0AE06
C	315	HIS	-	CLONING ARTIFACT	UNP P0AE06
C	316	HIS	-	CLONING ARTIFACT	UNP P0AE06
C	317	HIS	-	CLONING ARTIFACT	UNP P0AE06
C	318	HIS	-	CLONING ARTIFACT	UNP P0AE06
C	319	HIS	-	CLONING ARTIFACT	UNP P0AE06
C	320	HIS	-	CLONING ARTIFACT	UNP P0AE06
D	44	MSE	-	INITIATING METHIONINE	UNP P0AE06
D	223	MSE	PHE	ENGINEERED	UNP P0AE06
D	224	MSE	LEU	ENGINEERED	UNP P0AE06
D	287	MSE	LEU	ENGINEERED	UNP P0AE06
D	288	MSE	LEU	ENGINEERED	UNP P0AE06
D	291	MSE	MET	MODIFIED RESIDUE	UNP P0AE06
D	313	LEU	-	CLONING ARTIFACT	UNP P0AE06
D	314	GLU	-	CLONING ARTIFACT	UNP P0AE06
D	315	HIS	-	CLONING ARTIFACT	UNP P0AE06
D	316	HIS	-	CLONING ARTIFACT	UNP P0AE06
D	317	HIS	-	CLONING ARTIFACT	UNP P0AE06
D	318	HIS	-	CLONING ARTIFACT	UNP P0AE06
D	319	HIS	-	CLONING ARTIFACT	UNP P0AE06
D	320	HIS	-	CLONING ARTIFACT	UNP P0AE06

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	14	Total O 14 14	0	0
2	B	10	Total O 10 10	0	0

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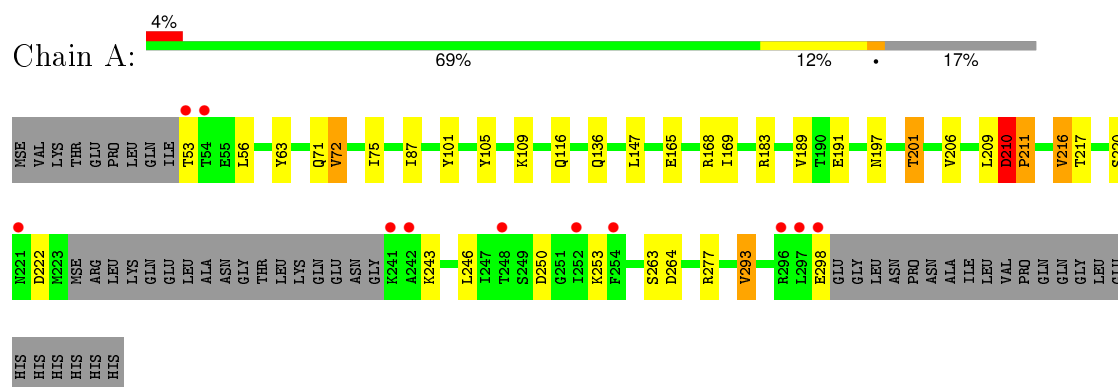
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	8	Total	O	0	0
			8	8		
2	D	4	Total	O	0	0
			4	4		

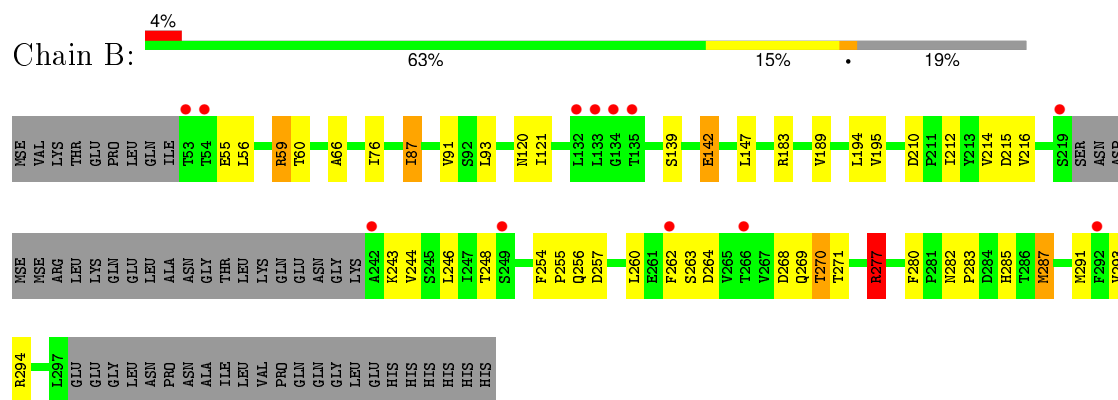
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

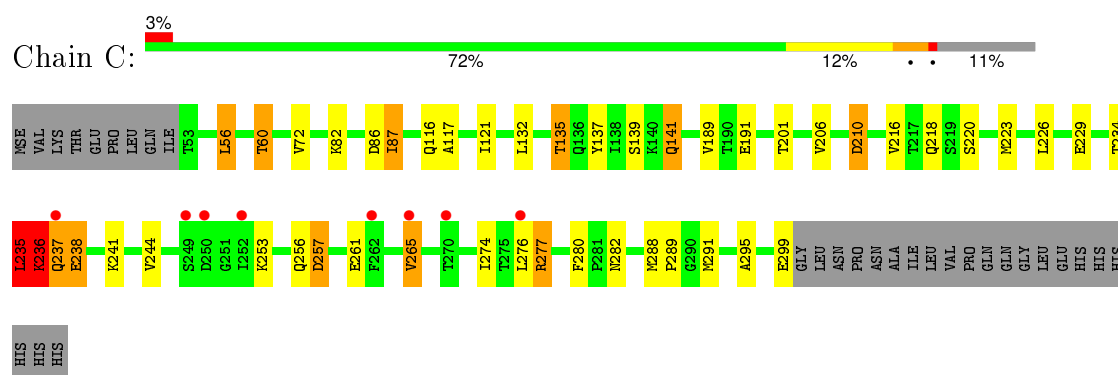
#### • Molecule 1: Acriflavine resistance protein A



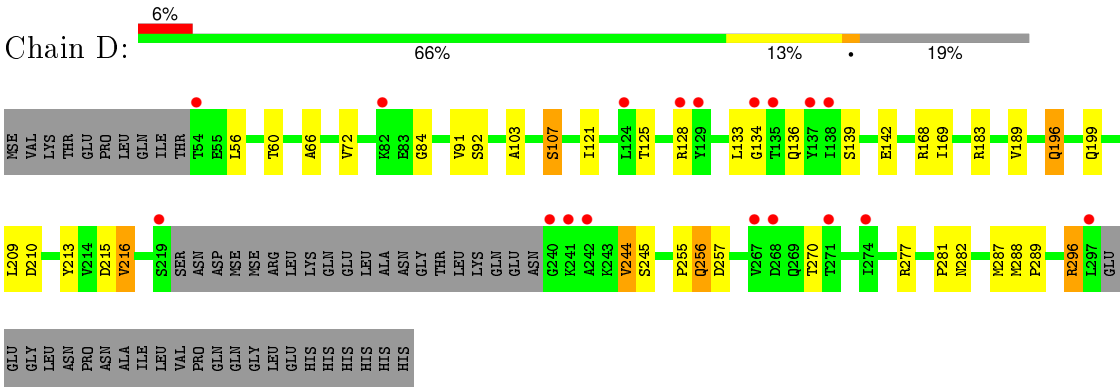
#### • Molecule 1: Acriflavine resistance protein A



#### • Molecule 1: Acriflavine resistance protein A



● Molecule 1: Acriflavine resistance protein A



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.79Å 100.03Å 332.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.71 32.44 – 2.71	Depositor EDS
% Data completeness (in resolution range)	96.7 (50.00-2.71) 96.7 (32.44-2.71)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.73 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.237 , 0.275 0.233 , 0.274	Depositor DCC
$R_{free}$ test set	1972 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	74.6	Xtriage
Anisotropy	0.259	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 66.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 39271 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	7000	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% 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	0.46	0/1746	0.77	1/2364 (0.0%)
1	B	0.45	0/1699	0.80	1/2304 (0.0%)
1	C	0.45	0/1888	0.83	3/2552 (0.1%)
1	D	0.42	0/1705	0.71	0/2310
All	All	0.44	0/7038	0.78	5/9530 (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	A	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	C	210	ASP	C-N-CD	6.69	142.46	128.40
1	C	277	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	B	277	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	A	222	ASP	CB-CG-OD2	5.25	123.02	118.30
1	C	257	ASP	CB-CG-OD2	5.05	122.84	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	210	ASP	Peptide

## 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	1728	0	1735	17	0
1	B	1680	0	1692	23	0
1	C	1870	0	1881	30	0
1	D	1686	0	1701	20	0
2	A	14	0	0	1	0
2	B	10	0	0	0	0
2	C	8	0	0	0	0
2	D	4	0	0	0	0
All	All	7000	0	7009	86	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 (86) 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:91:VAL:HG22	1:D:91:VAL:HG22	1.49	0.95
1:B:277:ARG:HH11	1:B:277:ARG:HG3	1.32	0.92
1:C:236:LYS:HD2	1:C:237:GLN:H	1.37	0.88
1:D:60:THR:HG22	1:D:289:PRO:HA	1.60	0.83
1:C:191:GLU:O	1:D:66:ALA:HA	1.79	0.83
1:C:56:LEU:HB3	1:C:218:GLN:HE21	1.47	0.80
1:D:215:ASP:OD1	1:D:277:ARG:HD3	1.81	0.79
1:D:196:GLN:H	1:D:199:GLN:HE21	1.31	0.78
1:B:87:ILE:HG13	1:B:93:LEU:HD21	1.73	0.71
1:D:196:GLN:H	1:D:199:GLN:NE2	1.90	0.70
1:C:139:SER:HB2	1:C:141:GLN:HE21	1.57	0.70
1:B:139:SER:OG	1:B:142:GLU:HB2	1.92	0.69
1:B:248:THR:HG22	1:B:287:MSE:HE1	1.76	0.68
1:A:191:GLU:O	1:B:66:ALA:HA	1.94	0.68
1:C:236:LYS:CD	1:C:237:GLN:H	2.10	0.63
1:C:60:THR:HG21	1:C:288:MSE:O	2.00	0.62
1:A:63:TYR:HB2	1:A:211:PRO:O	2.02	0.60
1:C:135:THR:HG23	1:C:137:TYR:H	1.68	0.58
1:A:210:ASP:HB3	1:A:211:PRO:CD	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:LEU:HG	1:B:280:PHE:CE2	2.40	0.56
1:D:213:TYR:CG	1:D:277:ARG:HD2	2.41	0.56
1:A:210:ASP:HB3	1:A:211:PRO:HD2	1.87	0.55
1:D:209:LEU:HD13	1:D:288:MSE:HE1	1.89	0.55
1:D:256:GLN:HB2	1:D:281:PRO:HG2	1.88	0.55
1:B:55:GLU:HG3	1:B:294:ARG:HE	1.72	0.54
1:A:209:LEU:C	1:A:210:ASP:O	2.46	0.53
1:B:246:LEU:HD22	1:B:293:VAL:CG1	2.37	0.53
1:B:87:ILE:HG13	1:B:93:LEU:CD2	2.37	0.53
1:C:244:VAL:HG12	1:C:280:PHE:HZ	1.74	0.53
1:A:246:LEU:HD22	1:A:293:VAL:CG1	2.39	0.53
1:D:245:SER:OG	1:D:296:ARG:HG3	2.08	0.53
1:C:60:THR:HG23	1:C:289:PRO:HA	1.92	0.52
1:A:210:ASP:O	1:A:211:PRO:C	2.48	0.52
1:B:210:ASP:O	1:B:282:ASN:N	2.43	0.51
1:C:132:LEU:O	1:C:135:THR:HG22	2.11	0.51
1:C:237:GLN:O	1:C:238:GLU:HB2	2.11	0.51
1:D:103:ALA:O	1:D:107:SER:HB2	2.11	0.50
1:C:265:VAL:HG12	1:C:276:LEU:CD2	2.41	0.50
1:C:265:VAL:HG12	1:C:276:LEU:HD23	1.93	0.50
1:B:283:PRO:HB2	1:D:168:ARG:CZ	2.42	0.50
1:D:56:LEU:HB3	1:D:216:VAL:HG11	1.93	0.50
1:D:244:VAL:O	1:D:257:ASP:HB2	2.11	0.50
1:A:201:THR:HG23	2:A:29:HOH:O	2.11	0.50
1:A:165:GLU:OE1	1:A:168:ARG:NH1	2.44	0.50
1:A:56:LEU:HD23	1:A:217:THR:O	2.12	0.49
1:C:60:THR:CG2	1:C:288:MSE:O	2.60	0.49
1:C:135:THR:CG2	1:C:137:TYR:H	2.26	0.49
1:D:84:GLY:O	1:D:183:ARG:HD2	2.13	0.48
1:C:244:VAL:CG1	1:C:295:ALA:HB1	2.44	0.48
1:D:139:SER:HB3	1:D:142:GLU:HB2	1.94	0.48
1:C:135:THR:HG23	1:C:137:TYR:N	2.29	0.48
1:D:121:ILE:O	1:D:125:THR:HG23	2.14	0.47
1:A:56:LEU:HD22	1:A:216:VAL:HG13	1.96	0.47
1:A:243:LYS:HB2	1:A:298:GLU:HB2	1.97	0.47
1:C:265:VAL:CG1	1:C:276:LEU:CD2	2.93	0.46
1:D:210:ASP:O	1:D:282:ASN:N	2.45	0.46
1:C:135:THR:HG23	1:C:137:TYR:HD2	1.82	0.45
1:C:244:VAL:HG12	1:C:280:PHE:CZ	2.52	0.45
1:C:117:ALA:O	1:C:121:ILE:HG12	2.16	0.45
1:A:71:GLN:O	1:A:197:ASN:ND2	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:ILE:HD11	1:B:195:VAL:HG11	1.97	0.45
1:B:264:ASP:N	1:B:264:ASP:OD1	2.50	0.44
1:D:287:MSE:O	1:D:288:MSE:HE2	2.17	0.44
1:D:56:LEU:CB	1:D:216:VAL:HG11	2.48	0.44
1:B:287:MSE:HE2	1:B:291:MSE:SE	2.68	0.43
1:B:212:ILE:HG21	1:B:287:MSE:HG2	1.99	0.43
1:C:210:ASP:O	1:C:282:ASN:N	2.51	0.43
1:C:234:THR:H	1:C:235:LEU:HD23	1.83	0.43
1:B:246:LEU:HD22	1:B:293:VAL:HG11	2.01	0.43
1:C:82:LYS:HA	1:C:82:LYS:HD3	1.87	0.43
1:A:105:TYR:CZ	1:A:168:ARG:HG3	2.54	0.43
1:A:210:ASP:O	1:A:211:PRO:O	2.37	0.42
1:B:215:ASP:OD1	1:B:277:ARG:NH1	2.53	0.42
1:C:265:VAL:CG1	1:C:276:LEU:HD21	2.49	0.42
1:A:72:VAL:HG13	1:A:101:TYR:CZ	2.54	0.42
1:C:236:LYS:CG	1:C:237:GLN:H	2.33	0.41
1:B:60:THR:HG22	1:B:214:VAL:HG22	2.01	0.41
1:B:254:PHE:HA	1:B:255:PRO:HD3	1.82	0.41
1:C:241:LYS:HD3	1:C:261:GLU:HB3	2.03	0.41
1:C:244:VAL:HG13	1:C:295:ALA:HB1	2.03	0.41
1:C:87:ILE:HA	1:C:87:ILE:HD12	1.78	0.41
1:C:60:THR:HG22	1:C:291:MSE:H	1.85	0.41
1:A:105:TYR:CZ	1:A:109:LYS:HD2	2.55	0.41
1:B:269:GLN:C	1:B:271:THR:H	2.24	0.41
1:B:244:VAL:O	1:B:257:ASP:HB2	2.22	0.40
1:B:59:ARG:HH11	1:B:59:ARG:HB2	1.86	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	225/277 (81%)	213 (95%)	8 (4%)	4 (2%)	11	25
1	B	219/277 (79%)	209 (95%)	7 (3%)	3 (1%)	14	34
1	C	245/277 (88%)	236 (96%)	6 (2%)	3 (1%)	16	38
1	D	220/277 (79%)	208 (94%)	9 (4%)	3 (1%)	14	34
All	All	909/1108 (82%)	866 (95%)	30 (3%)	13 (1%)	14	34

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	210	ASP
1	C	236	LYS
1	C	238	GLU
1	D	136	GLN
1	A	136	GLN
1	B	263	SER
1	B	268	ASP
1	A	211	PRO
1	A	263	SER
1	C	235	LEU
1	D	255	PRO
1	B	270	THR
1	D	134	GLY

### 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	184/220 (84%)	166 (90%)	18 (10%)	10	22
1	B	178/220 (81%)	159 (89%)	19 (11%)	8	18
1	C	199/220 (90%)	173 (87%)	26 (13%)	5	11
1	D	178/220 (81%)	165 (93%)	13 (7%)	17	39
All	All	739/880 (84%)	663 (90%)	76 (10%)	9	20

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	THR
1	A	72	VAL
1	A	75	ILE
1	A	87	ILE
1	A	116	GLN
1	A	147	LEU
1	A	169	ILE
1	A	183	ARG
1	A	189	VAL
1	A	201	THR
1	A	206	VAL
1	A	216	VAL
1	A	220	SER
1	A	250	ASP
1	A	253	LYS
1	A	264	ASP
1	A	277	ARG
1	A	293	VAL
1	B	56	LEU
1	B	59	ARG
1	B	87	ILE
1	B	120	ASN
1	B	121	ILE
1	B	142	GLU
1	B	147	LEU
1	B	183	ARG
1	B	189	VAL
1	B	194	LEU
1	B	216	VAL
1	B	243	LYS
1	B	256	GLN
1	B	260	LEU
1	B	262	PHE
1	B	270	THR
1	B	277	ARG
1	B	285	HIS
1	B	287	MSE
1	C	56	LEU
1	C	60	THR
1	C	72	VAL
1	C	86	ASP
1	C	87	ILE
1	C	116	GLN

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Mol	Chain	Res	Type
1	C	135	THR
1	C	141	GLN
1	C	189	VAL
1	C	201	THR
1	C	206	VAL
1	C	216	VAL
1	C	220	SER
1	C	223	MSE
1	C	226	LEU
1	C	229	GLU
1	C	235	LEU
1	C	236	LYS
1	C	237	GLN
1	C	253	LYS
1	C	256	GLN
1	C	257	ASP
1	C	265	VAL
1	C	274	ILE
1	C	277	ARG
1	C	299	GLU
1	D	72	VAL
1	D	92	SER
1	D	107	SER
1	D	128	ARG
1	D	133	LEU
1	D	169	ILE
1	D	189	VAL
1	D	196	GLN
1	D	216	VAL
1	D	244	VAL
1	D	256	GLN
1	D	270	THR
1	D	296	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	154	ASN
1	A	269	GLN
1	B	207	GLN
1	C	80	ASN
1	C	141	GLN

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Mol	Chain	Res	Type
1	C	218	GLN
1	C	237	GLN
1	D	197	ASN
1	D	199	GLN
1	D	208	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	225/277 (81%)	0.18	11 (4%)	33 32	50, 62, 67, 72	5 (2%)
1	B	220/277 (79%)	0.36	12 (5%)	29 27	48, 61, 68, 74	6 (2%)
1	C	242/277 (87%)	0.31	8 (3%)	50 50	52, 61, 66, 70	0
1	D	221/277 (79%)	0.54	18 (8%)	15 13	56, 63, 69, 71	5 (2%)
All	All	908/1108 (81%)	0.34	49 (5%)	29 28	48, 61, 68, 74	16 (1%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	134	GLY	7.8
1	D	137	TYR	5.3
1	B	292	PHE	4.2
1	B	135	THR	4.1
1	A	298	GLU	4.1
1	B	54	THR	3.9
1	A	241	LYS	3.7
1	C	250	ASP	3.7
1	A	53	THR	3.7
1	A	252	ILE	3.5
1	D	135	THR	3.4
1	B	219	SER	3.2
1	C	249	SER	3.2
1	A	242	ALA	3.1
1	D	268	ASP	3.1
1	A	248	THR	3.0
1	C	270	THR	3.0
1	D	124	LEU	3.0
1	D	128	ARG	2.9
1	B	249	SER	2.7
1	D	297	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	262	PHE	2.7
1	D	274	ILE	2.7
1	B	53	THR	2.6
1	A	297	LEU	2.6
1	B	132	LEU	2.6
1	D	240	GLY	2.6
1	C	237	GLN	2.5
1	D	219	SER	2.5
1	D	241	LYS	2.5
1	A	254	PHE	2.5
1	B	266	THR	2.4
1	D	134	GLY	2.4
1	B	133	LEU	2.4
1	D	129	TYR	2.4
1	D	54	THR	2.4
1	C	252	ILE	2.3
1	C	276	LEU	2.3
1	A	54	THR	2.3
1	A	296	ARG	2.3
1	D	138	ILE	2.2
1	C	265	VAL	2.1
1	D	271	THR	2.1
1	C	262	PHE	2.1
1	D	267	VAL	2.1
1	D	82	LYS	2.1
1	B	242	ALA	2.1
1	A	221	ASN	2.0
1	D	242	ALA	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.