



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 07:30 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 validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

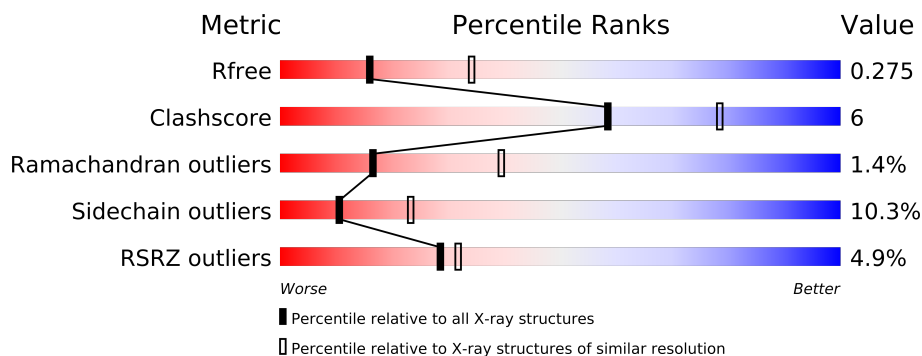
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	1770 (2.74-2.70)
Clashscore	79885	2183 (2.74-2.70)
Ramachandran outliers	78287	2147 (2.74-2.70)
Sidechain outliers	78261	2148 (2.74-2.70)
RSRZ outliers	66119	1772 (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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

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 hydrogen and 0 are deuterium.

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.

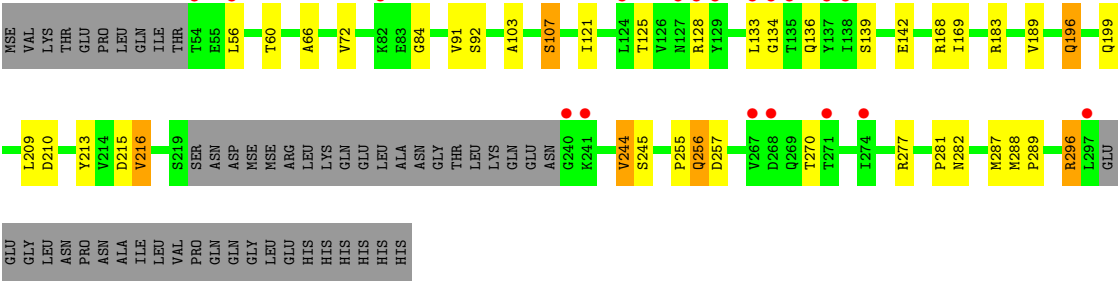
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	14	Total O 14 14	0	0
2	B	10	Total O 10 10	0	0

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

Chain D:



GLU
GLY
LEU
ASN
PRO
ASN
ALA
ILE
LEU
VAL
PRO
GLN
GLN
GLY
LEU
GLU
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	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
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.237 , 0.275 0.233 , 0.275	Depositor DCC
R_{free} test set	1972 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	74.6	Xtriage
Anisotropy	0.259	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 76.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 39271 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7000	wwPDB-VP
Average B, all atoms (Å ²)	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.*

¹Intensities estimated from amplitudes.

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(°)	Ideal(°)
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 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

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

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (86) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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	Distance(Å)	Clash(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
1:B:76:ILE:HD11	1:B:195:VAL:HG11	1.97	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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

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	225/277 (81%)	213 (95%)	8 (4%)	4 (2%)	13	30
1	B	219/277 (79%)	209 (95%)	7 (3%)	3 (1%)	16	39
1	C	245/277 (88%)	236 (96%)	6 (2%)	3 (1%)	19	44

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	220/277 (79%)	208 (94%)	9 (4%)	3 (1%)	16	39
All	All	909/1108 (82%)	866 (95%)	30 (3%)	13 (1%)	16	39

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%)	12	26
1	B	178/220 (81%)	159 (89%)	19 (11%)	10	21
1	C	199/220 (90%)	173 (87%)	26 (13%)	6	14
1	D	178/220 (81%)	165 (93%)	13 (7%)	20	44
All	All	739/880 (84%)	663 (90%)	76 (10%)	10	23

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

Mol	Chain	Res	Type
1	A	53	THR
1	A	72	VAL

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Mol	Chain	Res	Type
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
1	C	135	THR

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Mol	Chain	Res	Type
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
1	C	218	GLN

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Mol	Chain	Res	Type
1	C	237	GLN
1	D	197	ASN
1	D	199	GLN
1	D	208	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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 ⓘ

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, 95th 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 ²)	Q<0.9
1	A	229/277 (82%)	0.10	11 (4%) 29 33	50, 62, 67, 72	5 (2%)
1	B	223/277 (80%)	0.20	11 (4%) 28 32	48, 61, 68, 74	6 (2%)
1	C	247/277 (89%)	0.11	4 (1%) 68 74	52, 61, 67, 70	0
1	D	224/277 (80%)	0.38	19 (8%) 11 12	56, 63, 69, 71	5 (2%)
All	All	923/1108 (83%)	0.19	45 (4%) 28 32	48, 62, 68, 74	16 (1%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	134	GLY	7.8
1	D	137	TYR	4.8
1	B	135	THR	4.5
1	A	298	GLU	4.4
1	A	241	LYS	4.4
1	B	292	PHE	3.9
1	A	53	THR	3.8
1	B	54	THR	3.7
1	C	250	ASP	3.5
1	C	249	SER	3.4
1	D	135	THR	3.4
1	D	54	THR	3.2
1	A	242	ALA	3.2
1	D	274	ILE	3.1
1	B	219	SER	3.0
1	D	268	ASP	3.0
1	A	252	ILE	3.0
1	A	248	THR	3.0
1	B	262	PHE	3.0
1	D	134	GLY	3.0
1	D	124	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	297	LEU	3.0
1	D	138	ILE	2.7
1	B	249	SER	2.7
1	A	297	LEU	2.7
1	B	132	LEU	2.7
1	B	242	ALA	2.7
1	D	271	THR	2.6
1	B	53	THR	2.5
1	A	270	THR	2.4
1	D	240	GLY	2.3
1	A	54	THR	2.3
1	C	270	THR	2.3
1	B	133	LEU	2.3
1	D	129	TYR	2.3
1	D	128	ARG	2.3
1	D	267	VAL	2.2
1	A	254	PHE	2.2
1	C	237	GLN	2.2
1	A	296	ARG	2.1
1	D	241	LYS	2.1
1	D	133	LEU	2.1
1	D	127	ASN	2.1
1	D	56	LEU	2.1
1	D	82	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.