



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 6, 2017 – 02:43 PM EST

PDB ID : 4I83
Title : Crystal Structure of (3R)-Hydroxymyristoyl-ACP Dehydratase from *Neisseria meningitidis* FAM18
Authors : Khandokar, Y.B.; Forwood, J.K.
Deposited on : unknown
Resolution : 2.60 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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20030345

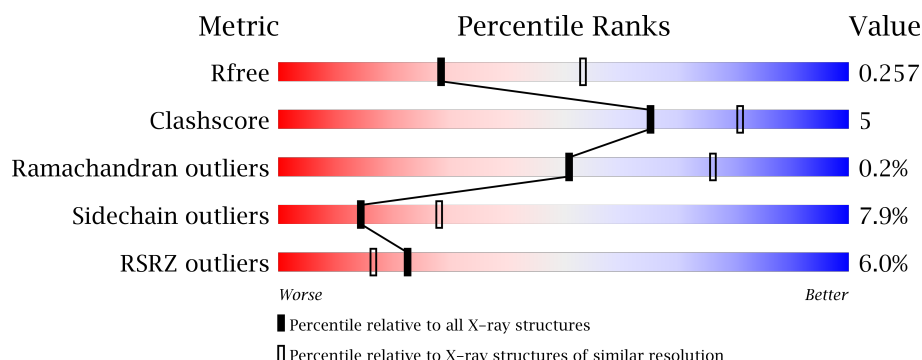
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	100719	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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. 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	152	<div> <div>3%</div> <div>82%</div> <div>13%</div> <div>5%</div> </div>
1	B	152	<div> <div>7%</div> <div>80%</div> <div>14%</div> <div>• 5%</div> </div>
1	C	152	<div> <div>7%</div> <div>80%</div> <div>14%</div> <div>• 5%</div> </div>
1	D	152	<div> <div>5%</div> <div>80%</div> <div>14%</div> <div>• 5%</div> </div>
1	E	152	<div> <div>7%</div> <div>80%</div> <div>13%</div> <div>•• 5%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	152	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: a small red segment at the beginning labeled '7%', a large green segment labeled '82%', a yellow segment labeled '13%', and a small grey segment at the end labeled '5%'.

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6756 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 3-hydroxyacyl-[acyl-carrier-protein] dehydratase FabZ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	144	Total	C	N	O	S	6	0	0
			1126	729	191	200	6			
1	C	144	Total	C	N	O	S	6	0	0
			1126	729	191	200	6			
1	B	144	Total	C	N	O	S	12	0	0
			1126	729	191	200	6			
1	D	144	Total	C	N	O	S	6	0	0
			1126	729	191	200	6			
1	E	144	Total	C	N	O	S	12	0	0
			1126	729	191	200	6			
1	F	144	Total	C	N	O	S	6	0	0
			1126	729	191	200	6			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP A1KRL1
A	-1	ASN	-	EXPRESSION TAG	UNP A1KRL1
A	0	ALA	-	EXPRESSION TAG	UNP A1KRL1
C	-2	SER	-	EXPRESSION TAG	UNP A1KRL1
C	-1	ASN	-	EXPRESSION TAG	UNP A1KRL1
C	0	ALA	-	EXPRESSION TAG	UNP A1KRL1
B	-2	SER	-	EXPRESSION TAG	UNP A1KRL1
B	-1	ASN	-	EXPRESSION TAG	UNP A1KRL1
B	0	ALA	-	EXPRESSION TAG	UNP A1KRL1
D	-2	SER	-	EXPRESSION TAG	UNP A1KRL1
D	-1	ASN	-	EXPRESSION TAG	UNP A1KRL1
D	0	ALA	-	EXPRESSION TAG	UNP A1KRL1
E	-2	SER	-	EXPRESSION TAG	UNP A1KRL1
E	-1	ASN	-	EXPRESSION TAG	UNP A1KRL1
E	0	ALA	-	EXPRESSION TAG	UNP A1KRL1
F	-2	SER	-	EXPRESSION TAG	UNP A1KRL1
F	-1	ASN	-	EXPRESSION TAG	UNP A1KRL1

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	ALA	-	EXPRESSION TAG	UNP A1KRL1

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 ($RSRZ > 2$). 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: 3-hydroxyacyl-[acyl-carrier-protein] dehydratase FabZ

Chain A: 




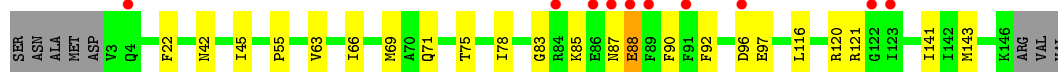
- Molecule 1: 3-hydroxyacyl-[acyl-carrier-protein] dehydratase FabZ

Chain C: 




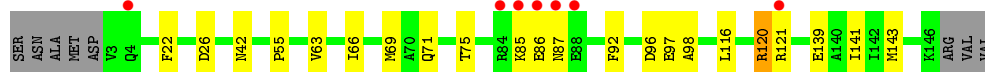
- Molecule 1: 3-hydroxyacyl-[acyl-carrier-protein] dehydratase FabZ

Chain B: 




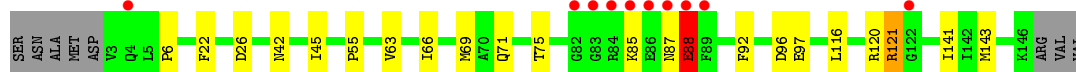
- Molecule 1: 3-hydroxyacyl-[acyl-carrier-protein] dehydratase FabZ

Chain D: 

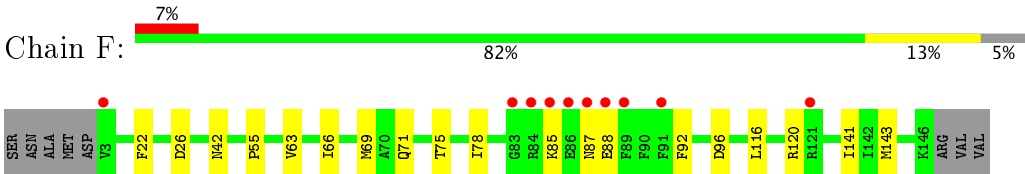


- Molecule 1: 3-hydroxyacyl-[acyl-carrier-protein] dehydratase FabZ

Chain E: 



- Molecule 1: 3-hydroxyacyl-[acyl-carrier-protein] dehydratase FabZ



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.44Å 103.98Å 135.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 2.60 19.96 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (19.96-2.60) 100.0 (19.96-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.59Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.224 , 0.256 0.225 , 0.257	Depositor DCC
R_{free} test set	1617 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	46.5	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 35.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6756	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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	0.64	1/1148 (0.1%)	0.74	1/1550 (0.1%)
1	B	0.73	2/1148 (0.2%)	0.70	0/1550
1	C	0.61	1/1148 (0.1%)	0.69	1/1550 (0.1%)
1	D	0.79	1/1148 (0.1%)	0.73	2/1550 (0.1%)
1	E	0.61	1/1148 (0.1%)	0.72	1/1550 (0.1%)
1	F	0.64	1/1148 (0.1%)	0.68	1/1550 (0.1%)
All	All	0.67	7/6888 (0.1%)	0.71	6/9300 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	120	ARG	CB-CG	-17.08	1.06	1.52
1	B	121	ARG	CB-CG	13.74	1.89	1.52
1	F	120	ARG	CB-CG	-10.77	1.23	1.52
1	A	120	ARG	CB-CG	-8.17	1.30	1.52
1	B	120	ARG	CB-CG	-8.09	1.30	1.52
1	E	88	GLU	CD-OE2	6.75	1.33	1.25
1	C	120	ARG	CB-CG	-5.51	1.37	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	120	ARG	CA-CB-CG	8.88	132.95	113.40
1	A	120	ARG	CB-CG-CD	8.24	133.01	111.60
1	D	120	ARG	CB-CG-CD	5.55	126.04	111.60
1	E	120	ARG	CA-CB-CG	-5.47	101.36	113.40
1	F	120	ARG	CB-CG-CD	5.38	125.58	111.60
1	C	120	ARG	CB-CG-CD	5.20	125.12	111.60

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	1126	0	1159	11	0
1	B	1126	0	1159	17	0
1	C	1126	0	1159	12	0
1	D	1126	0	1159	12	0
1	E	1126	0	1159	13	0
1	F	1126	0	1159	12	0
All	All	6756	0	6954	64	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 5.

All (64) 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:83:GLY:HA3	1:B:88:GLU:OE1	1.77	0.84
1:B:78:ILE:HA	1:B:88:GLU:OE2	1.87	0.74
1:C:66:ILE:HD13	1:C:69:MET:CE	2.19	0.73
1:B:66:ILE:HD13	1:B:69:MET:CE	2.20	0.71
1:A:66:ILE:HD13	1:A:69:MET:CE	2.20	0.71
1:E:66:ILE:HD13	1:E:69:MET:CE	2.20	0.71
1:F:66:ILE:HD13	1:F:69:MET:CE	2.21	0.71
1:D:66:ILE:HD13	1:D:69:MET:CE	2.20	0.71
1:F:78:ILE:HG12	1:F:88:GLU:OE1	1.91	0.69
1:D:66:ILE:HA	1:D:69:MET:HE2	1.76	0.66
1:F:96:ASP:HB3	1:F:141:ILE:HB	1.79	0.65
1:A:96:ASP:HB3	1:A:141:ILE:HB	1.77	0.64
1:A:66:ILE:HA	1:A:69:MET:HE2	1.79	0.64
1:C:66:ILE:HD13	1:C:69:MET:HE2	1.79	0.63
1:E:87:ASN:O	1:E:88:GLU:CD	2.37	0.63
1:B:66:ILE:HA	1:B:69:MET:HE2	1.81	0.62
1:E:66:ILE:HD13	1:E:69:MET:HE2	1.82	0.61
1:C:96:ASP:HB3	1:C:141:ILE:HB	1.82	0.61
1:A:42:ASN:HB3	1:D:42:ASN:HB3	1.83	0.61
1:B:66:ILE:HD13	1:B:69:MET:HE2	1.81	0.60
1:F:66:ILE:HD13	1:F:69:MET:HE2	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:ASP:HB3	1:D:141:ILE:HB	1.84	0.60
1:E:66:ILE:HA	1:E:69:MET:HE2	1.83	0.60
1:F:66:ILE:HA	1:F:69:MET:HE2	1.84	0.59
1:A:66:ILE:HD13	1:A:69:MET:HE2	1.85	0.58
1:C:42:ASN:HB3	1:E:42:ASN:HB3	1.85	0.58
1:B:96:ASP:HB3	1:B:141:ILE:HB	1.85	0.58
1:D:66:ILE:HD13	1:D:69:MET:HE2	1.86	0.57
1:C:66:ILE:HA	1:C:69:MET:HE2	1.87	0.56
1:B:42:ASN:HB3	1:F:42:ASN:HB3	1.87	0.55
1:B:78:ILE:HG12	1:B:88:GLU:OE2	2.06	0.55
1:C:96:ASP:O	1:D:96:ASP:O	2.25	0.53
1:E:87:ASN:O	1:E:88:GLU:OE2	2.29	0.50
1:E:96:ASP:O	1:F:96:ASP:O	2.30	0.50
1:D:71:GLN:O	1:D:75:THR:HG23	2.12	0.49
1:F:96:ASP:CB	1:F:141:ILE:HB	2.41	0.49
1:A:71:GLN:O	1:A:75:THR:HG23	2.13	0.49
1:A:96:ASP:O	1:B:96:ASP:O	2.31	0.49
1:D:96:ASP:CB	1:D:141:ILE:HB	2.43	0.48
1:B:96:ASP:CB	1:B:141:ILE:HB	2.43	0.48
1:C:71:GLN:O	1:C:75:THR:HG23	2.14	0.47
1:C:63:VAL:HB	1:D:63:VAL:HB	1.96	0.47
1:B:78:ILE:CG1	1:B:88:GLU:OE2	2.63	0.47
1:E:71:GLN:O	1:E:75:THR:HG23	2.15	0.47
1:D:66:ILE:HD13	1:D:69:MET:HE1	1.95	0.46
1:B:71:GLN:O	1:B:75:THR:HG23	2.15	0.46
1:F:71:GLN:O	1:F:75:THR:HG23	2.16	0.45
1:E:96:ASP:HB3	1:E:141:ILE:HB	1.99	0.45
1:A:63:VAL:HB	1:B:63:VAL:HB	1.99	0.44
1:A:66:ILE:HD13	1:A:69:MET:HE1	1.97	0.44
1:A:96:ASP:CB	1:A:141:ILE:HB	2.44	0.44
1:B:66:ILE:HD13	1:B:69:MET:HE1	1.99	0.43
1:C:96:ASP:CB	1:C:141:ILE:HB	2.46	0.43
1:B:78:ILE:CA	1:B:88:GLU:OE2	2.63	0.42
1:B:88:GLU:HG3	1:B:90:PHE:CE1	2.55	0.42
1:F:78:ILE:CG1	1:F:88:GLU:OE1	2.65	0.42
1:E:63:VAL:HB	1:F:63:VAL:HB	2.01	0.42
1:A:45:ILE:HG22	1:D:26:ASP:OD1	2.20	0.41
1:C:26:ASP:OD1	1:E:45:ILE:HG22	2.21	0.41
1:B:45:ILE:HG22	1:F:26:ASP:OD1	2.21	0.41
1:C:121:ARG:HB2	1:C:121:ARG:HE	1.65	0.41
1:C:45:ILE:HG22	1:E:26:ASP:OD1	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:ALA:HA	1:D:139:GLU:O	2.21	0.40
1:E:66:ILE:HD13	1:E:69:MET:HE1	2.00	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	142/152 (93%)	133 (94%)	8 (6%)	1 (1%)	25	49
1	B	142/152 (93%)	134 (94%)	8 (6%)	0	100	100
1	C	142/152 (93%)	132 (93%)	10 (7%)	0	100	100
1	D	142/152 (93%)	135 (95%)	7 (5%)	0	100	100
1	E	142/152 (93%)	133 (94%)	8 (6%)	1 (1%)	25	49
1	F	142/152 (93%)	133 (94%)	9 (6%)	0	100	100
All	All	852/912 (93%)	800 (94%)	50 (6%)	2 (0%)	51	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	87	ASN
1	E	121	ARG

5.3.2 Protein sidechains [i](#)

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	120/127 (94%)	111 (92%)	9 (8%)	16	31
1	B	120/127 (94%)	111 (92%)	9 (8%)	16	31
1	C	120/127 (94%)	109 (91%)	11 (9%)	11	20
1	D	120/127 (94%)	109 (91%)	11 (9%)	11	20
1	E	120/127 (94%)	110 (92%)	10 (8%)	13	25
1	F	120/127 (94%)	113 (94%)	7 (6%)	23	46
All	All	720/762 (94%)	663 (92%)	57 (8%)	14	28

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	22	PHE
1	A	55	PRO
1	A	85	LYS
1	A	86	GLU
1	A	92	PHE
1	A	97	GLU
1	A	116	LEU
1	A	143	MET
1	C	4	GLN
1	C	22	PHE
1	C	55	PRO
1	C	85	LYS
1	C	87	ASN
1	C	88	GLU
1	C	92	PHE
1	C	97	GLU
1	C	116	LEU
1	C	121	ARG
1	C	143	MET
1	B	22	PHE
1	B	55	PRO
1	B	85	LYS
1	B	87	ASN
1	B	88	GLU
1	B	92	PHE
1	B	97	GLU
1	B	116	LEU
1	B	143	MET
1	D	22	PHE

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Mol	Chain	Res	Type
1	D	55	PRO
1	D	85	LYS
1	D	86	GLU
1	D	87	ASN
1	D	92	PHE
1	D	97	GLU
1	D	116	LEU
1	D	120	ARG
1	D	121	ARG
1	D	143	MET
1	E	6	PRO
1	E	22	PHE
1	E	55	PRO
1	E	85	LYS
1	E	88	GLU
1	E	92	PHE
1	E	97	GLU
1	E	116	LEU
1	E	121	ARG
1	E	143	MET
1	F	22	PHE
1	F	55	PRO
1	F	85	LYS
1	F	87	ASN
1	F	92	PHE
1	F	116	LEU
1	F	143	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	109	GLN
1	C	24	GLN
1	C	51	GLN
1	C	109	GLN
1	B	24	GLN
1	B	51	GLN
1	B	87	ASN
1	B	109	GLN
1	D	24	GLN
1	D	109	GLN

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Mol	Chain	Res	Type
1	E	24	GLN
1	E	51	GLN
1	E	87	ASN
1	E	109	GLN
1	F	24	GLN
1	F	51	GLN
1	F	87	ASN
1	F	109	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, 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	144/152 (94%)	-0.19	4 (2%)	53 46	31, 42, 69, 80	1 (0%)
1	B	144/152 (94%)	-0.04	10 (6%)	18 12	31, 47, 80, 116	2 (1%)
1	C	144/152 (94%)	-0.01	11 (7%)	15 10	33, 49, 76, 128	1 (0%)
1	D	144/152 (94%)	0.01	7 (4%)	30 24	30, 44, 75, 129	1 (0%)
1	E	144/152 (94%)	0.06	10 (6%)	18 12	33, 51, 84, 124	2 (1%)
1	F	144/152 (94%)	0.17	10 (6%)	18 12	32, 49, 79, 130	1 (0%)
All	All	864/912 (94%)	0.00	52 (6%)	23 17	30, 47, 80, 130	8 (0%)

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	88	GLU	8.1
1	F	85	LYS	7.5
1	C	86	GLU	7.3
1	D	88	GLU	6.9
1	B	88	GLU	6.6
1	F	86	GLU	5.6
1	E	84	ARG	5.6
1	D	86	GLU	5.4
1	D	87	ASN	5.0
1	E	86	GLU	4.7
1	B	84	ARG	4.6
1	F	121	ARG	4.5
1	E	85	LYS	4.5
1	B	86	GLU	4.4
1	C	84	ARG	4.2
1	B	87	ASN	4.0
1	E	88	GLU	3.9
1	C	121	ARG	3.9
1	D	85	LYS	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	121	ARG	3.8
1	D	84	ARG	3.7
1	C	87	ASN	3.7
1	D	4	GLN	3.6
1	E	87	ASN	3.5
1	A	121	ARG	3.3
1	B	122	GLY	3.2
1	C	85	LYS	3.1
1	E	89	PHE	3.1
1	C	88	GLU	3.0
1	F	87	ASN	3.0
1	E	122	GLY	3.0
1	F	84	ARG	2.8
1	C	4	GLN	2.7
1	C	3	VAL	2.7
1	E	4	GLN	2.7
1	F	91	PHE	2.7
1	F	89	PHE	2.6
1	B	91	PHE	2.6
1	F	3	VAL	2.5
1	B	89	PHE	2.5
1	E	82	GLY	2.5
1	A	88	GLU	2.5
1	C	96	ASP	2.5
1	B	4	GLN	2.4
1	E	83	GLY	2.4
1	B	123	ILE	2.3
1	C	135	GLN	2.2
1	A	72	ALA	2.2
1	B	96	ASP	2.2
1	C	5	LEU	2.1
1	A	122	GLY	2.1
1	F	83	GLY	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.