



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 17, 2023 – 03:44 PM EDT

PDB ID : 2FYZ
Title : Structural of Mumps virus fusion protein core
Authors : Lou, Z.; Xu, Y.; Liu, Y.; Rao, Z.
Deposited on : 2006-02-08
Resolution : 2.20 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Xtrriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

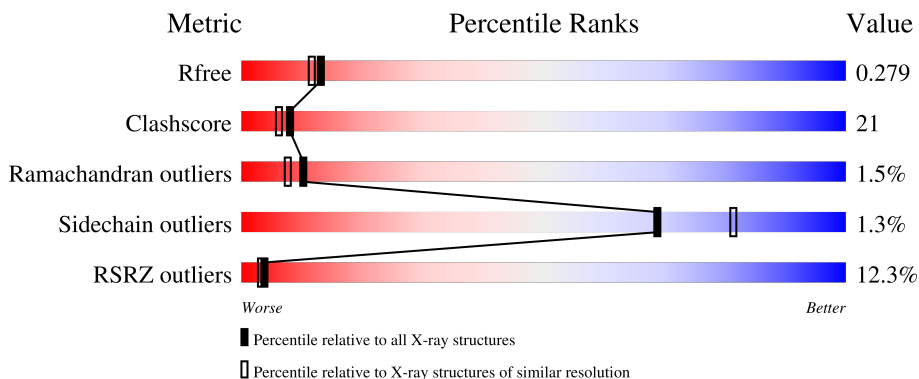
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


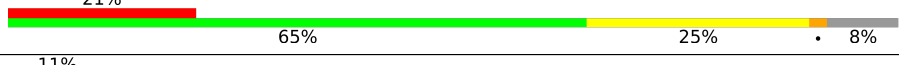



The reported resolution of this entry is 2.20 Å.

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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	63	
1	C	63	
1	E	63	
2	B	48	
2	D	48	

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Mol	Chain	Length	Quality of chain
2	F	48	 <p>A horizontal bar chart representing the quality of the chain. The bar is divided into three segments: a green segment on the left labeled '56%', a yellow segment in the middle labeled '17%', and a grey segment on the right labeled '25%'. A small black dot is positioned at the end of the grey segment.</p>

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 2330 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 Fusion glycoprotein F0.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	57	Total 428	C 261	N 82	O 83	S 2	0	0	0
1	C	58	Total 426	C 262	N 79	O 83	S 2	0	0	0
1	E	62	Total 452	C 278	N 86	O 87	S 1	0	0	0

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	119	GLY	-	cloning artifact	UNP P11236
A	120	PRO	-	cloning artifact	UNP P11236
A	121	LEU	-	cloning artifact	UNP P11236
A	122	GLY	-	cloning artifact	UNP P11236
A	123	SER	-	cloning artifact	UNP P11236
C	119	GLY	-	cloning artifact	UNP P11236
C	120	PRO	-	cloning artifact	UNP P11236
C	121	LEU	-	cloning artifact	UNP P11236
C	122	GLY	-	cloning artifact	UNP P11236
C	123	SER	-	cloning artifact	UNP P11236
E	119	GLY	-	cloning artifact	UNP P11236
E	120	PRO	-	cloning artifact	UNP P11236
E	121	LEU	-	cloning artifact	UNP P11236
E	122	GLY	-	cloning artifact	UNP P11236
E	123	SER	-	cloning artifact	UNP P11236

- Molecule 2 is a protein called Fusion glycoprotein F0.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	32	Total 246	C 151	N 43	O 52	0	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	31	Total	C	N	O	0	0	0
			242	149	42	51			
2	F	36	Total	C	N	O	0	0	0
			273	167	50	56			

There are 33 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	438	ASN	-	cloning artifact	UNP P11236
B	439	MET	-	cloning artifact	UNP P11236
B	440	SER	-	cloning artifact	UNP P11236
B	441	SER	-	cloning artifact	UNP P11236
B	442	GLY	-	cloning artifact	UNP P11236
B	443	GLY	-	cloning artifact	UNP P11236
B	444	ARG	-	cloning artifact	UNP P11236
B	445	GLY	-	cloning artifact	UNP P11236
B	446	GLY	-	cloning artifact	UNP P11236
B	463	THR	ALA	engineered mutation	UNP P11236
B	478	ILE	ASN	engineered mutation	UNP P11236
D	438	ASN	-	cloning artifact	UNP P11236
D	439	MET	-	cloning artifact	UNP P11236
D	440	SER	-	cloning artifact	UNP P11236
D	441	SER	-	cloning artifact	UNP P11236
D	442	GLY	-	cloning artifact	UNP P11236
D	443	GLY	-	cloning artifact	UNP P11236
D	444	ARG	-	cloning artifact	UNP P11236
D	445	GLY	-	cloning artifact	UNP P11236
D	446	GLY	-	cloning artifact	UNP P11236
D	463	THR	ALA	engineered mutation	UNP P11236
D	478	ILE	ASN	engineered mutation	UNP P11236
F	438	ASN	-	cloning artifact	UNP P11236
F	439	MET	-	cloning artifact	UNP P11236
F	440	SER	-	cloning artifact	UNP P11236
F	441	SER	-	cloning artifact	UNP P11236
F	442	GLY	-	cloning artifact	UNP P11236
F	443	GLY	-	cloning artifact	UNP P11236
F	444	ARG	-	cloning artifact	UNP P11236
F	445	GLY	-	cloning artifact	UNP P11236
F	446	GLY	-	cloning artifact	UNP P11236
F	463	THR	ALA	engineered mutation	UNP P11236
F	478	ILE	ASN	engineered mutation	UNP P11236

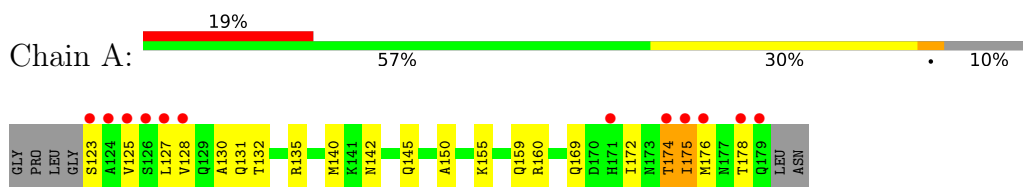
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	50	Total O 50 50	0	0
3	B	43	Total O 43 43	0	0
3	C	44	Total O 44 44	0	0
3	D	43	Total O 43 43	0	0
3	E	36	Total O 36 36	0	0
3	F	47	Total O 47 47	0	0

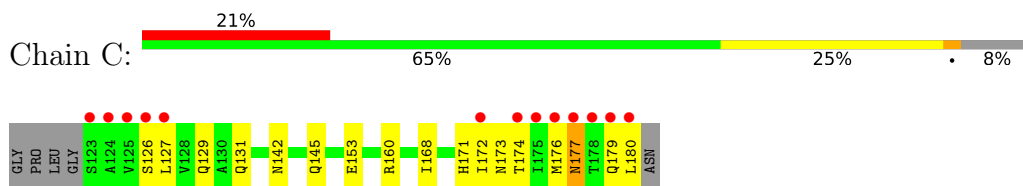
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

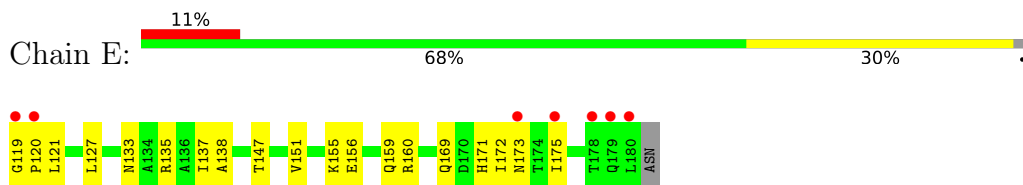
- Molecule 1: Fusion glycoprotein F0



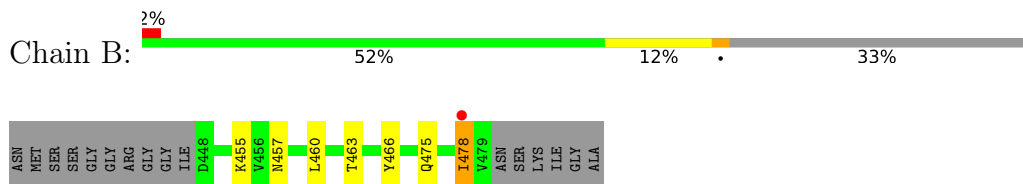
- Molecule 1: Fusion glycoprotein F0



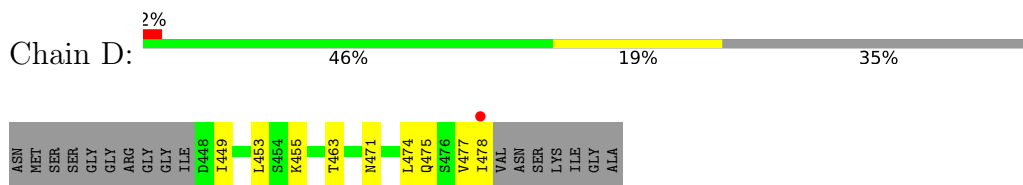
- Molecule 1: Fusion glycoprotein F0



- Molecule 2: Fusion glycoprotein F0



- Molecule 2: Fusion glycoprotein F0



- Molecule 2: Fusion glycoprotein F0

Chain F:  56% 17% 25%

ASN	MET	SER	SER	GLY	E444	G445	G446	I447	V456	M462	K468	E469	V477	I478	V479	ASN	SER	LYS	ILE	GLY	ALA
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	161.17Å 60.81Å 40.15Å 90.00° 98.36° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 40.02 – 2.21	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.20) 88.9 (40.02-2.21)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.04 (at 2.22Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.231 , 0.278 0.232 , 0.279	Depositor DCC
R_{free} test set	905 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	18.3	Xtrriage
Anisotropy	0.133	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 67.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2330	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.25% 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 [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	0.30	0/429	0.47	0/580
1	C	0.32	0/427	0.47	0/579
1	E	0.35	0/454	0.54	0/616
2	B	0.35	0/247	0.54	0/332
2	D	0.34	0/243	0.50	0/327
2	F	0.33	0/274	0.52	0/367
All	All	0.33	0/2074	0.50	0/2801

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	428	0	441	31	0
1	C	426	0	435	22	0
1	E	452	0	463	21	0
2	B	246	0	243	14	0
2	D	242	0	242	9	0
2	F	273	0	273	9	0
3	A	50	0	0	3	0
3	B	43	0	0	4	0
3	C	44	0	0	3	0
3	D	43	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	36	0	0	0	0
3	F	47	0	0	2	0
All	All	2330	0	2097	89	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 21.

All (89) 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:E:119:GLY:O	1:E:121:LEU:N	2.01	0.92
1:A:150:ALA:HA	3:B:162:HOH:O	1.74	0.87
2:B:463:THR:HA	3:B:162:HOH:O	1.77	0.84
1:E:156:GLU:O	1:E:160:ARG:HG2	1.81	0.80
2:F:444:ARG:HG2	2:F:447:ILE:H	1.47	0.78
1:C:179:GLN:O	1:C:180:LEU:HG	1.83	0.77
1:A:142:ASN:HA	1:A:145:GLN:HE21	1.50	0.76
1:E:160:ARG:HB2	2:F:456:VAL:HG22	1.71	0.71
1:E:119:GLY:C	1:E:121:LEU:H	1.96	0.69
1:A:142:ASN:HA	1:A:145:GLN:NE2	2.07	0.69
1:A:172:ILE:O	1:A:176:MET:HB2	1.92	0.69
2:B:475:GLN:HE21	2:B:475:GLN:HA	1.58	0.69
1:A:127:LEU:O	1:A:131:GLN:HG3	1.94	0.67
2:B:478:ILE:HD11	1:E:138:ALA:HA	1.76	0.66
2:B:457:ASN:HB3	3:B:267:HOH:O	1.98	0.64
1:A:174:THR:HG22	1:A:175:ILE:HG13	1.80	0.64
2:B:466:TYR:HD1	3:B:162:HOH:O	1.80	0.63
1:C:160:ARG:HH11	2:D:455:LYS:HB3	1.63	0.63
2:F:444:ARG:HE	2:F:446:GLY:HA3	1.63	0.63
1:A:128:VAL:HA	1:A:131:GLN:OE1	1.99	0.62
1:E:169:GLN:NE2	1:E:173:ASN:HD21	1.99	0.61
1:C:172:ILE:HG22	1:C:172:ILE:O	2.01	0.60
1:A:123:SER:C	1:A:125:VAL:H	2.05	0.58
1:C:127:LEU:O	1:C:131:GLN:HG3	2.04	0.58
1:C:177:ASN:C	1:C:177:ASN:HD22	2.07	0.58
1:A:174:THR:HG22	1:A:175:ILE:N	2.19	0.57
1:E:169:GLN:NE2	1:E:173:ASN:ND2	2.53	0.56
2:F:444:ARG:HG3	2:F:446:GLY:H	1.71	0.56
1:A:155:LYS:NZ	1:A:159:GLN:NE2	2.54	0.56
1:C:142:ASN:HA	1:C:145:GLN:HE21	1.71	0.55
1:A:174:THR:HG22	1:A:175:ILE:H	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:475:GLN:HA	2:B:475:GLN:NE2	2.22	0.54
2:D:471:ASN:O	2:D:475:GLN:HG2	2.09	0.53
1:A:130:ALA:HB3	3:A:219:HOH:O	2.08	0.53
2:D:477:VAL:O	2:D:478:ILE:HB	2.08	0.53
2:F:477:VAL:O	2:F:478:ILE:HD13	2.09	0.53
1:C:173:ASN:O	1:C:177:ASN:HB3	2.08	0.52
1:C:177:ASN:C	1:C:177:ASN:ND2	2.62	0.52
2:B:478:ILE:O	2:B:478:ILE:HG22	2.08	0.52
1:C:160:ARG:NH1	2:D:455:LYS:HB3	2.25	0.51
2:F:444:ARG:HG3	2:F:446:GLY:N	2.27	0.50
1:C:145:GLN:HG2	3:C:191:HOH:O	2.11	0.50
1:A:169:GLN:HA	1:A:169:GLN:OE1	2.12	0.49
1:C:172:ILE:HA	1:C:176:MET:CG	2.42	0.49
1:C:171:HIS:O	1:C:176:MET:HG2	2.12	0.49
1:A:155:LYS:NZ	1:A:159:GLN:HE22	2.11	0.49
1:C:179:GLN:O	1:C:180:LEU:CG	2.60	0.48
1:A:127:LEU:HD12	3:A:219:HOH:O	2.12	0.48
2:B:460:LEU:O	2:B:463:THR:HB	2.13	0.48
1:A:145:GLN:HG3	2:D:474:LEU:HD21	1.95	0.47
2:B:478:ILE:HD11	1:E:138:ALA:CA	2.44	0.47
1:A:132:THR:HA	1:A:135:ARG:NH1	2.29	0.47
1:A:169:GLN:NE2	2:D:449:ILE:HD12	2.30	0.47
1:A:123:SER:C	1:A:125:VAL:N	2.68	0.47
1:E:171:HIS:CD2	1:E:175:ILE:HB	2.49	0.47
1:C:168:ILE:O	1:C:172:ILE:HG13	2.15	0.47
2:F:468:LYS:HB2	3:F:283:HOH:O	2.15	0.47
1:E:135:ARG:HG2	1:E:135:ARG:HH11	1.80	0.46
1:A:178:THR:HG22	1:A:178:THR:O	2.14	0.46
1:A:140:MET:CE	1:E:137:ILE:HG23	2.45	0.46
1:A:140:MET:HE2	1:E:137:ILE:HG23	1.97	0.46
1:A:176:MET:CE	1:E:172:ILE:HG23	2.45	0.46
1:C:153:GLU:HB2	2:D:463:THR:OG1	2.16	0.46
1:E:127:LEU:C	1:E:127:LEU:HD23	2.36	0.46
1:C:172:ILE:HA	1:C:176:MET:HG3	1.98	0.46
2:F:469:GLU:HG3	3:F:108:HOH:O	2.15	0.46
1:A:175:ILE:O	1:A:176:MET:HG2	2.16	0.45
2:B:478:ILE:HD11	1:E:138:ALA:CB	2.45	0.45
1:C:173:ASN:O	1:C:174:THR:C	2.55	0.45
1:A:155:LYS:HD3	1:A:155:LYS:C	2.37	0.45
2:B:457:ASN:OD1	1:E:159:GLN:NE2	2.46	0.44
1:C:172:ILE:O	1:C:176:MET:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:475:GLN:HE21	2:B:475:GLN:CA	2.22	0.44
1:E:133:ASN:O	1:E:137:ILE:HG13	2.18	0.43
1:C:177:ASN:HA	3:C:222:HOH:O	2.17	0.43
2:D:477:VAL:O	2:D:478:ILE:CB	2.65	0.43
1:A:155:LYS:HZ1	1:A:159:GLN:NE2	2.16	0.43
1:E:160:ARG:HB2	2:F:456:VAL:CG2	2.43	0.42
2:B:460:LEU:HB3	1:E:155:LYS:HE2	2.02	0.42
1:A:131:GLN:O	1:A:135:ARG:HG3	2.21	0.41
1:C:127:LEU:HD12	3:C:196:HOH:O	2.19	0.41
1:E:135:ARG:HG2	1:E:135:ARG:NH1	2.35	0.41
1:E:147:THR:O	1:E:151:VAL:HG23	2.21	0.41
1:A:155:LYS:HZ3	1:A:159:GLN:NE2	2.18	0.41
1:A:174:THR:CG2	1:A:175:ILE:H	2.28	0.40
3:A:219:HOH:O	1:C:126:SER:HA	2.21	0.40
2:D:453:LEU:HD23	2:D:453:LEU:HA	1.89	0.40
1:A:160:ARG:HD3	2:B:455:LYS:HB3	2.03	0.40
1:A:130:ALA:HB1	1:C:129:GLN:HG3	2.02	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	55/63 (87%)	51 (93%)	2 (4%)	2 (4%)	3	1
1	C	56/63 (89%)	50 (89%)	6 (11%)	0	100	100
1	E	60/63 (95%)	56 (93%)	3 (5%)	1 (2%)	9	6
2	B	30/48 (62%)	28 (93%)	1 (3%)	1 (3%)	4	2
2	D	29/48 (60%)	29 (100%)	0	0	100	100
2	F	34/48 (71%)	33 (97%)	1 (3%)	0	100	100
All	All	264/333 (79%)	247 (94%)	13 (5%)	4 (2%)	10	8

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	174	THR
2	B	478	ILE
1	E	120	PRO
1	A	175	ILE

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	45/49 (92%)	45 (100%)	0	100	100
1	C	44/49 (90%)	43 (98%)	1 (2%)	50	63
1	E	46/49 (94%)	46 (100%)	0	100	100
2	B	29/41 (71%)	29 (100%)	0	100	100
2	D	29/41 (71%)	29 (100%)	0	100	100
2	F	31/41 (76%)	29 (94%)	2 (6%)	17	19
All	All	224/270 (83%)	221 (99%)	3 (1%)	69	81

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	177	ASN
2	F	444	ARG
2	F	462	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	GLN
1	A	159	GLN
1	A	177	ASN
2	B	475	GLN
1	C	145	GLN

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Mol	Chain	Res	Type
1	C	177	ASN
2	D	457	ASN
2	D	461	GLN
2	D	472	HIS
1	E	131	GLN
1	E	145	GLN
1	E	169	GLN
1	E	171	HIS

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 monosaccharides 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	57/63 (90%)	0.60	12 (21%) 1 0	12, 26, 90, 94	0
1	C	58/63 (92%)	0.55	13 (22%) 0 0	12, 24, 91, 93	0
1	E	62/63 (98%)	0.28	7 (11%) 5 4	11, 23, 84, 99	0
2	B	32/48 (66%)	-0.18	1 (3%) 49 47	12, 23, 43, 62	0
2	D	31/48 (64%)	-0.10	1 (3%) 47 45	12, 25, 51, 65	0
2	F	36/48 (75%)	-0.38	0 100 100	12, 20, 29, 39	0
All	All	276/333 (82%)	0.22	34 (12%) 4 3	11, 23, 90, 99	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	119	GLY	9.7
1	C	175	ILE	8.7
1	A	175	ILE	7.8
1	A	125	VAL	7.3
1	E	178	THR	7.0
1	E	175	ILE	5.5
1	A	176	MET	5.2
1	C	174	THR	5.1
1	C	123	SER	5.0
1	C	176	MET	4.9
1	C	179	GLN	4.8
1	A	179	GLN	4.5
1	E	180	LEU	4.4
1	C	178	THR	4.0
1	E	179	GLN	3.9
1	C	126	SER	3.9
1	C	177	ASN	3.6
1	E	173	ASN	3.4
1	A	124	ALA	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	126	SER	3.4
1	A	123	SER	3.2
1	C	124	ALA	3.2
1	C	180	LEU	2.8
1	C	125	VAL	2.8
2	B	478	ILE	2.7
1	A	127	LEU	2.7
1	E	120	PRO	2.6
1	A	174	THR	2.3
1	C	172	ILE	2.3
1	A	128	VAL	2.3
1	C	127	LEU	2.2
1	A	171	HIS	2.2
2	D	478	ILE	2.1
1	A	178	THR	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 monosaccharides 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.