Full wwPDB X-ray Structure Validation Report (i)

Jun 22, 2024 – 10:07 AM EDT

PDB ID	:	6B03
Title	:	The crystal structure of the ferredoxin protease FusC in complex with its substrate plant ferredoxin
Authors	:	Grinter, R.
Deposited on	:	2017-09-13
Resolution	:	2.70 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 (i) 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 (1)) were used in the production of this report:

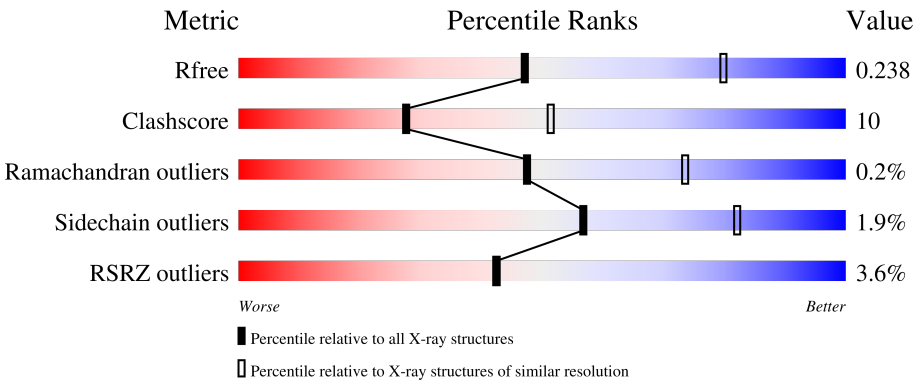
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 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 <=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	924	% 76% 20% • •
2	C	104	2% 5% • 93%
2	F	104	21% 28% 10% • 61%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7620 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 Putative zinc protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	896	7199	4519	1290	1368	22	0	11	0

- Molecule 2 is a protein called Ferredoxin-2, chloroplastic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	F	41	315	200	43	71	1	0	0	0
			Total	C	N	O				
2	C	7	56	34	9	13		0	0	0

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	95	ALA	-	expression tag	UNP P16972
F	96	ILE	-	expression tag	UNP P16972
F	97	MET	-	expression tag	UNP P16972
F	98	LEU	-	expression tag	UNP P16972
F	99	GLU	-	expression tag	UNP P16972
F	100	HIS	-	expression tag	UNP P16972
F	101	HIS	-	expression tag	UNP P16972
F	102	HIS	-	expression tag	UNP P16972
F	103	HIS	-	expression tag	UNP P16972
F	104	HIS	-	expression tag	UNP P16972
F	105	HIS	-	expression tag	UNP P16972
C	144	ALA	-	expression tag	UNP P16972
C	145	ILE	-	expression tag	UNP P16972
C	146	MET	-	expression tag	UNP P16972
C	147	LEU	-	expression tag	UNP P16972
C	148	GLU	-	expression tag	UNP P16972
C	149	HIS	-	expression tag	UNP P16972
C	150	HIS	-	expression tag	UNP P16972

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Chain	Residue	Modelled	Actual	Comment	Reference
C	151	HIS	-	expression tag	UNP P16972
C	152	HIS	-	expression tag	UNP P16972
C	153	HIS	-	expression tag	UNP P16972
C	154	HIS	-	expression tag	UNP P16972

- Molecule 3 is water.

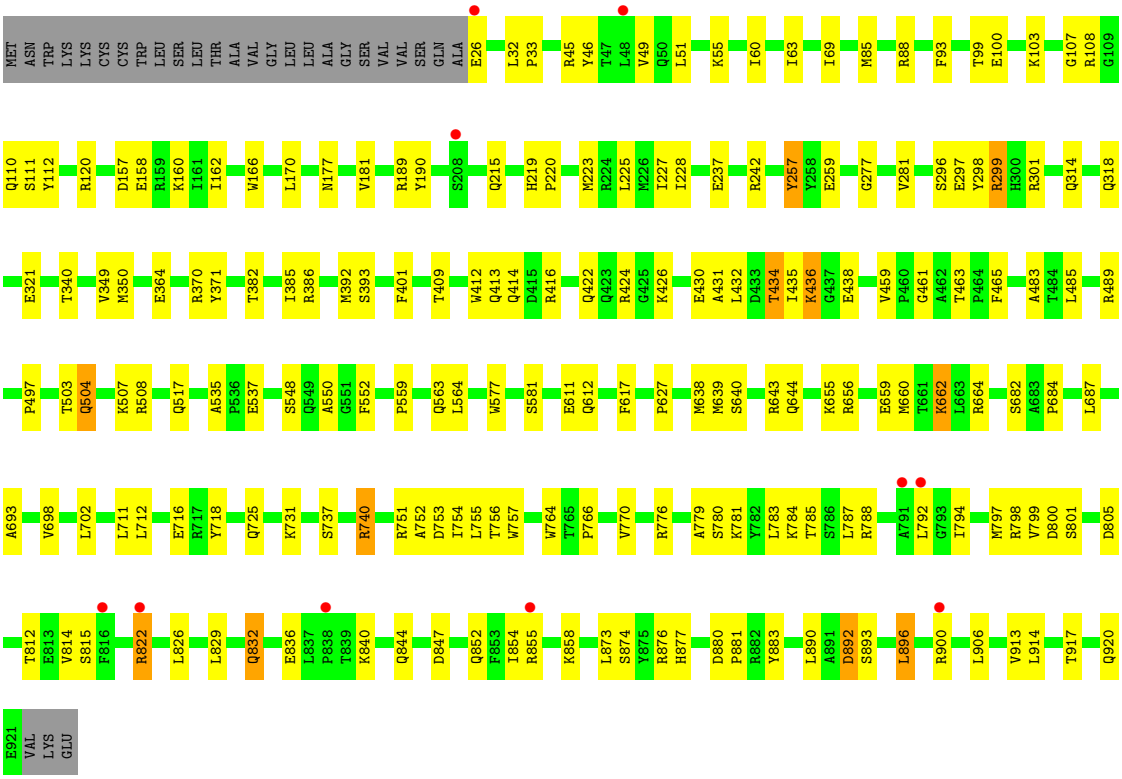
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	50	50	50	0	0

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.

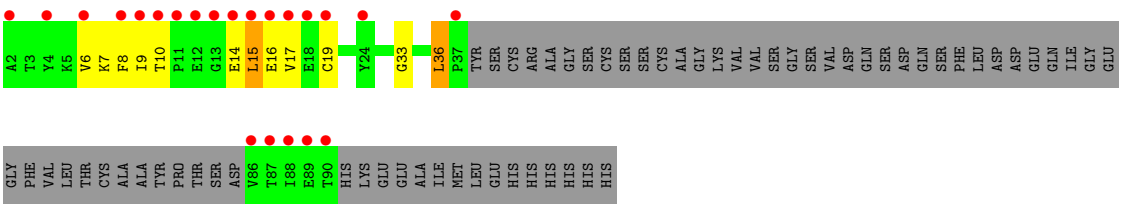
• Molecule 1: Putative zinc protease

Chain A:



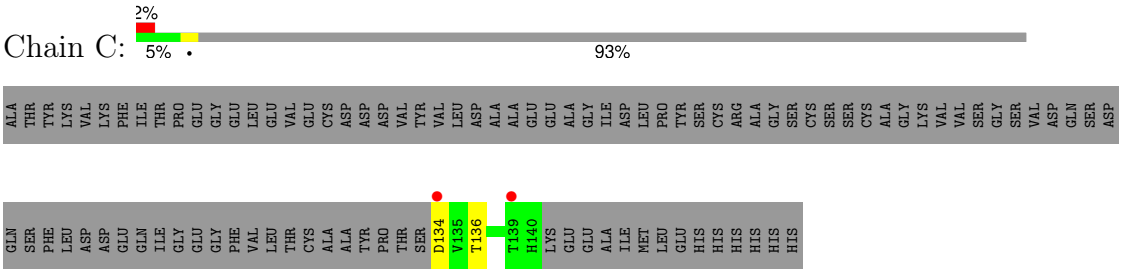
• Molecule 2: Ferredoxin-2, chloroplastic

Chain F:



• Molecule 2: Ferredoxin-2, chloroplastic

Chain C:


2%
5%
93%
ALA THR TYR LYS VAL LYS PHE ILE THR PRO GLU GLY GLU LEU GLU VAL GLU CYS ASP ASP ASP VAL TYR VAL LEU ASP ALA ALA GLU GLU ALA GLY ILE ASP LEU PRO TYR SER CYS ARG ALA GLY SER CYS SER SER CYS ALA GLY LYS VAL VAL SER GLY SER VAL ASP GLN SER ASP
GLN SER PHE LEU ASP ASP GLU GLN ILE GLY GLU GLY PHE VAL LEU THR CYS ALA ALA TYR PRO THR SER D134 V135 T136 T139 H140 LYS GLU GLU ALA ILE MET LEU GLU HIS HIS HIS HIS HIS HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α, β, γ	82.58Å 127.26Å 133.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.29 – 2.70 41.92 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (41.29-2.70) 99.9 (41.92-2.70)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.46 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.196 , 0.243 0.197 , 0.238	Depositor DCC
R_{free} test set	2001 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å^2)	63.2	Xtriage
Anisotropy	0.520	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 41.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.012 for -h,l,k	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	7620	wwPDB-VP
Average B, all atoms (Å^2)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.47	0/7347	0.67	0/9962
2	C	0.53	0/56	0.56	0/76
2	F	0.58	0/318	0.81	1/433 (0.2%)
All	All	0.47	0/7721	0.67	1/10471 (0.0%)

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	3

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	15	LEU	CA-CB-CG	7.35	132.22	115.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	299	ARG	Sidechain
1	A	503	THR	Peptide
1	A	822	ARG	Peptide

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	7199	0	7185	138	0
2	C	56	0	50	1	0
2	F	315	0	297	14	0
3	A	50	0	0	6	0
All	All	7620	0	7532	151	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 10.

All (151) 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)
2:F:7:LYS:HD3	2:F:16:GLU:HA	1.55	0.86
2:F:7:LYS:HE2	2:F:8:PHE:H	1.42	0.85
1:A:832:GLN:OE1	3:A:1001:HOH:O	2.01	0.79
1:A:436:LYS:HE2	1:A:438:GLU:H	1.48	0.78
1:A:756:THR:HG22	1:A:914:LEU:HD12	1.65	0.78
1:A:296:SER:HB3	1:A:299:ARG:HB2	1.66	0.78
1:A:318:GLN:HE22	1:A:489:ARG:H	1.33	0.76
1:A:85:MET:HE2	1:A:88:ARG:HG3	1.68	0.75
1:A:799:VAL:HG22	1:A:814:VAL:HG12	1.69	0.73
1:A:93:PHE:CD1	1:A:100:GLU:HG3	2.24	0.73
1:A:314:GLN:HE21	1:A:318:GLN:HE21	1.38	0.71
1:A:640:SER:O	1:A:644:GLN:HG3	1.90	0.71
1:A:755:LEU:HD22	1:A:815:SER:HB2	1.71	0.71
1:A:784:LYS:HG2	1:A:797:MET:CE	2.21	0.69
2:F:7:LYS:CD	2:F:16:GLU:HA	2.24	0.65
1:A:737[A]:SER:O	1:A:740:ARG:NH1	2.25	0.63
1:A:314:GLN:HE21	1:A:318:GLN:NE2	1.96	0.63
1:A:725:GLN:NE2	3:A:1011:HOH:O	2.31	0.63
1:A:497:PRO:HB2	1:A:577:TRP:CE2	2.33	0.63
1:A:297:GLU:OE1	1:A:297:GLU:N	2.23	0.63
1:A:414:GLN:HG3	1:A:416:ARG:HH11	1.64	0.62
2:F:7:LYS:HE2	2:F:8:PHE:N	2.11	0.62
1:A:393:SER:HB2	1:A:426:LYS:HG2	1.82	0.62
1:A:781:LYS:HD3	1:A:784:LYS:HE3	1.82	0.61
2:F:7:LYS:HE3	2:F:15:LEU:O	2.02	0.60
1:A:801:SER:HB2	1:A:812:THR:HG22	1.83	0.59
1:A:435:ILE:O	1:A:435:ILE:HG22	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:852:GLN:OE1	1:A:855:ARG:NH2	2.34	0.59
1:A:299:ARG:HG3	1:A:431:ALA:HB1	1.84	0.59
1:A:259:GLU:HG3	1:A:340:THR:HG23	1.86	0.58
1:A:461:GLY:O	3:A:1002:HOH:O	2.18	0.57
1:A:784:LYS:C	1:A:788:ARG:HE	2.07	0.57
1:A:892:ASP:N	1:A:892:ASP:OD1	2.37	0.57
1:A:754:ILE:C	1:A:755:LEU:HD23	2.25	0.56
1:A:794:ILE:HG12	1:A:826:LEU:HD21	1.88	0.56
1:A:896:LEU:HD21	1:A:900:ARG:HH11	1.70	0.56
1:A:382:THR:HG23	1:A:432:LEU:HD22	1.87	0.56
1:A:801:SER:CB	1:A:812:THR:HG22	2.35	0.56
1:A:99:THR:O	1:A:103:LYS:HE2	2.06	0.55
1:A:920:GLN:NE2	3:A:1016:HOH:O	2.40	0.55
1:A:784:LYS:HG2	1:A:797:MET:HE2	1.88	0.55
1:A:611:GLU:OE2	1:A:612:GLN:NE2	2.41	0.54
1:A:85:MET:HE2	1:A:85:MET:HA	1.89	0.54
1:A:844:GLN:HA	1:A:847:ASP:OD1	2.07	0.54
1:A:656:ARG:O	1:A:660:MET:HG3	2.08	0.53
1:A:409:THR:O	1:A:413:GLN:HG3	2.08	0.53
1:A:177:ASN:O	1:A:181:VAL:HG23	2.09	0.53
1:A:409:THR:HG22	1:A:413:GLN:HE21	1.73	0.53
1:A:242:ARG:HG3	1:A:242:ARG:HH11	1.74	0.53
1:A:805:ASP:HB2	1:A:877:HIS:NE2	2.23	0.53
1:A:107:GLY:H	1:A:111:SER:HB2	1.74	0.52
1:A:85:MET:CE	1:A:88:ARG:HG3	2.39	0.52
1:A:854:ILE:O	1:A:858:LYS:HG2	2.08	0.52
1:A:281:VAL:HG13	1:A:349:VAL:HG21	1.92	0.52
1:A:780:SER:O	1:A:784:LYS:HG3	2.10	0.51
1:A:364:GLU:HG2	3:A:1014:HOH:O	2.09	0.51
1:A:798:ARG:NH1	1:A:800:ASP:OD2	2.44	0.51
1:A:459:VAL:HG12	1:A:463:THR:HB	1.92	0.51
1:A:627:PRO:HG2	1:A:687:LEU:HD22	1.93	0.51
1:A:659:GLU:HA	1:A:662:LYS:HE3	1.92	0.51
1:A:764:TRP:CZ2	1:A:873:LEU:HD21	2.46	0.51
1:A:617:PHE:HB3	1:A:718:TYR:HB2	1.94	0.50
1:A:836:GLU:HB3	1:A:840:LYS:HE3	1.92	0.50
1:A:504:GLN:OE1	1:A:504:GLN:N	2.44	0.50
1:A:550:ALA:HB1	1:A:693:ALA:HA	1.94	0.50
1:A:783:LEU:HD22	1:A:799:VAL:HG21	1.94	0.50
1:A:655:LYS:O	1:A:659:GLU:HG3	2.12	0.49
1:A:158:GLU:HG3	1:A:162:ILE:HD11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:564:LEU:HD21	1:A:638[B]:MET:HE1	1.94	0.49
1:A:787:LEU:HD11	1:A:829:LEU:HB3	1.94	0.49
1:A:639:MET:O	1:A:643:ARG:HG2	2.12	0.49
2:F:6:VAL:O	2:F:7:LYS:HD2	2.13	0.49
1:A:370:ARG:NH2	1:A:483:ALA:O	2.43	0.49
1:A:756:THR:HA	1:A:913:VAL:O	2.11	0.49
1:A:60:ILE:HG12	1:A:227:ILE:HG12	1.95	0.49
1:A:754:ILE:O	1:A:755:LEU:HD23	2.13	0.49
1:A:507:LYS:HD2	1:A:508:ARG:O	2.13	0.48
1:A:46:TYR:HA	1:A:225:LEU:O	2.13	0.48
1:A:301[B]:ARG:NH2	2:C:136:THR:HG23	2.28	0.48
1:A:660:MET:HE2	1:A:755:LEU:HD12	1.95	0.48
1:A:33:PRO:HB3	1:A:51:LEU:HD23	1.94	0.48
1:A:46:TYR:OH	1:A:237:GLU:OE2	2.25	0.48
1:A:660:MET:CE	1:A:755:LEU:HD12	2.44	0.48
1:A:638[B]:MET:SD	1:A:638[B]:MET:N	2.86	0.47
1:A:548:SER:HB3	1:A:698:VAL:HG22	1.95	0.47
1:A:890:LEU:O	1:A:893:SER:OG	2.31	0.47
1:A:26:GLU:CB	1:A:401:PHE:H	2.28	0.47
1:A:756:THR:HG22	1:A:914:LEU:CD1	2.40	0.46
2:F:9:ILE:HD13	2:F:14:GLU:HG3	1.98	0.46
1:A:215:GLN:O	1:A:219:HIS:NE2	2.46	0.46
1:A:49:VAL:HG22	1:A:228:ILE:HG12	1.98	0.46
1:A:63:ILE:HD13	1:A:120[B]:ARG:HD2	1.98	0.46
1:A:99:THR:O	1:A:103:LYS:HG2	2.16	0.46
1:A:298:TYR:OH	1:A:424[B]:ARG:HG3	2.15	0.46
1:A:779:ALA:O	1:A:783:LEU:HG	2.15	0.46
1:A:108:ARG:HG2	1:A:112:TYR:CZ	2.51	0.45
1:A:785:THR:N	1:A:788:ARG:HE	2.14	0.45
2:F:7:LYS:HD3	2:F:16:GLU:CA	2.38	0.45
1:A:497:PRO:HD3	1:A:581:SER:HB2	1.99	0.45
1:A:780:SER:O	1:A:784:LYS:HE2	2.16	0.45
1:A:321:GLU:OE2	1:A:489:ARG:HD3	2.17	0.45
1:A:858:LYS:HD3	1:A:858:LYS:N	2.32	0.45
2:F:7:LYS:HE2	2:F:7:LYS:HA	1.99	0.45
1:A:682:SER:HB2	1:A:684:PRO:HD2	1.99	0.44
1:A:414:GLN:CG	1:A:416:ARG:HH11	2.30	0.44
1:A:783:LEU:CD2	1:A:799:VAL:HG21	2.48	0.44
1:A:787:LEU:HD11	1:A:829:LEU:CB	2.47	0.44
1:A:32:LEU:HD12	1:A:33:PRO:HD2	2.00	0.44
1:A:430:GLU:O	1:A:430:GLU:HG2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:766:PRO:O	1:A:770:VAL:HG13	2.18	0.44
1:A:277:GLY:O	1:A:643:ARG:HD2	2.17	0.44
1:A:805:ASP:OD2	1:A:876:ARG:NH1	2.46	0.44
1:A:436:LYS:HE3	1:A:436:LYS:HB2	1.80	0.44
1:A:751:ARG:HG2	1:A:752:ALA:N	2.33	0.44
1:A:69:ILE:HD12	1:A:190:TYR:CE1	2.53	0.44
1:A:314:GLN:NE2	1:A:318:GLN:HE21	2.11	0.44
1:A:434:THR:O	1:A:434:THR:OG1	2.30	0.44
1:A:535:ALA:HB1	1:A:537:GLU:OE1	2.17	0.43
1:A:559:PRO:O	1:A:563:GLN:HG3	2.17	0.43
1:A:392:MET:O	1:A:422:GLN:HA	2.18	0.43
1:A:63:ILE:HG13	1:A:412:TRP:CH2	2.54	0.43
1:A:108:ARG:HG2	1:A:112:TYR:OH	2.19	0.43
1:A:49:VAL:CG2	1:A:228:ILE:HG12	2.48	0.43
1:A:371:TYR:OH	1:A:485:LEU:O	2.29	0.43
1:A:220:PRO:HA	1:A:223:MET:HE2	2.01	0.43
1:A:712:LEU:O	1:A:716:GLU:HG3	2.18	0.43
1:A:55:LYS:HA	1:A:55:LYS:HD3	1.90	0.43
1:A:643:ARG:HD2	2:F:33:GLY:HA3	2.02	0.42
1:A:874:SER:HB2	1:A:883:TYR:HB3	2.00	0.42
1:A:259:GLU:HG3	1:A:340:THR:CG2	2.48	0.42
1:A:157:ASP:O	1:A:160:LYS:HE3	2.19	0.42
2:F:7:LYS:HD2	2:F:7:LYS:HA	1.89	0.42
2:F:10:THR:HG21	2:F:36:LEU:HD12	2.01	0.42
1:A:45:ARG:NH1	3:A:1025:HOH:O	2.51	0.42
1:A:664:ARG:HD2	1:A:757:TRP:CD2	2.55	0.42
1:A:385:ILE:HG22	1:A:432:LEU:HD11	2.02	0.41
1:A:801:SER:OG	1:A:906:LEU:HD22	2.19	0.41
2:F:7:LYS:HA	2:F:7:LYS:CE	2.49	0.41
1:A:794:ILE:HG12	1:A:826:LEU:CD2	2.50	0.41
1:A:349:VAL:HG12	1:A:350:MET:O	2.20	0.41
1:A:702:LEU:HD13	1:A:711:LEU:HD21	2.03	0.41
1:A:792:LEU:HA	1:A:822:ARG:HG3	2.03	0.41
2:F:8:PHE:HE2	2:F:17:VAL:HG11	1.85	0.41
1:A:459:VAL:HG21	1:A:465:PHE:CD1	2.56	0.41
1:A:166:TRP:CH2	1:A:170:LEU:HD11	2.57	0.40
1:A:414:GLN:HG3	1:A:416:ARG:NH1	2.34	0.40
1:A:517:GLN:HG2	1:A:881:PRO:O	2.21	0.40
1:A:189:ARG:HD2	1:A:257:TYR:CD2	2.56	0.40
1:A:880:ASP:HB2	1:A:881:PRO:HD2	2.02	0.40
1:A:854:ILE:HG23	1:A:858:LYS:NZ	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:753:ASP:HB2	1:A:917:THR:HB	2.04	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	905/924 (98%)	878 (97%)	26 (3%)	1 (0%)	51 78
2	C	5/104 (5%)	5 (100%)	0	0	100 100
2	F	37/104 (36%)	33 (89%)	3 (8%)	1 (3%)	5 12
All	All	947/1132 (84%)	916 (97%)	29 (3%)	2 (0%)	47 73

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	36	LEU
1	A	504	GLN

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	776/789 (98%)	763 (98%)	13 (2%)	60 84
2	C	7/88 (8%)	6 (86%)	1 (14%)	3 8

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	F	35/88 (40%)	34 (97%)	1 (3%)	42	71
All	All	818/965 (85%)	803 (98%)	15 (2%)	57	83

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	GLN
1	A	257	TYR
1	A	386	ARG
1	A	434	THR
1	A	436	LYS
1	A	552	PHE
1	A	662	LYS
1	A	731	LYS
1	A	740	ARG
1	A	776	ARG
1	A	832	GLN
1	A	892	ASP
1	A	896	LEU
2	F	19	CYS
2	C	134	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	318	GLN
1	A	567	GLN
1	A	832	GLN
2	C	140	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 (i)

6.1 Protein, DNA and RNA chains (i)

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	896/924 (96%)	-0.17	10 (1%)	80	82	37, 63, 104, 139	1 (0%)
2	C	7/104 (6%)	1.53	2 (28%)	0	0	74, 89, 115, 126	0
2	F	41/104 (39%)	2.59	22 (53%)	0	0	59, 131, 148, 160	0
All	All	944/1132 (83%)	-0.03	34 (3%)	42	42	37, 65, 112, 160	1 (0%)

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	8	PHE	9.4
2	F	9	ILE	8.4
2	F	2	ALA	7.3
2	F	88	ILE	7.2
2	F	90	THR	6.5
2	F	10	THR	5.6
2	F	86	VAL	5.5
2	F	87	THR	5.4
2	F	89	GLU	5.3
1	A	791	ALA	5.0
2	F	15	LEU	4.4
2	F	11	PRO	4.0
2	F	14	GLU	3.8
1	A	838	PRO	3.7
2	F	12	GLU	3.7
2	F	17	VAL	3.7
2	F	13	GLY	3.7
1	A	816	PHE	3.6
1	A	792	LEU	3.5
2	F	6	VAL	3.5
2	C	139	THR	3.3
1	A	822	ARG	3.2
2	F	16	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
2	F	37	PRO	3.1
2	C	134	ASP	2.9
1	A	900	ARG	2.8
2	F	18	GLU	2.5
2	F	4	TYR	2.4
1	A	26	GLU	2.2
1	A	48	LEU	2.2
2	F	24	TYR	2.1
1	A	208	SER	2.1
1	A	855	ARG	2.1
2	F	19	CYS	2.1

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.