Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 08:29 PM GMT

PDB ID	:	4TKN
Title	:	Structure of the SNX17 FERM domain bound to the second NPxF motif of KRIT1
Authors	:	Stiegler, A.L.; Zhang, R.; Liu, W.; Boggon, T.J.
Deposited on	:	2014-05-27
Resolution	:	3.00 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

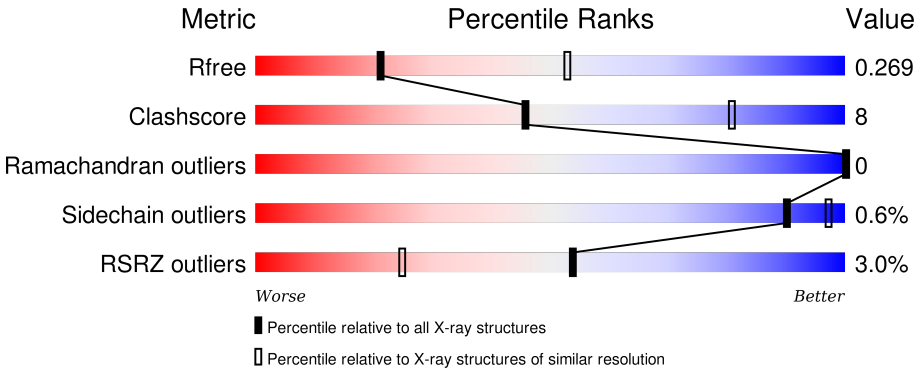
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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 <=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	286	% 73% 16% 10%
1	B	286	3% 72% 15% 14%
1	C	286	4% 76% 14% • 10%
2	D	19	26% 26% 47%
2	E	19	16% 32% 53%

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Mol	Chain	Length	Quality of chain
2	F	19	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6419 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 Sorting nexin-17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	256	Total	C	N	O	S	0	0	0
			2075	1321	351	392	11			
1	B	247	Total	C	N	O	S	0	0	0
			1996	1272	335	379	10			
1	C	257	Total	C	N	O	S	0	0	0
			2081	1326	350	394	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	106	GLY	-	expression tag	UNP Q15036
A	107	SER	-	expression tag	UNP Q15036
B	106	GLY	-	expression tag	UNP Q15036
B	107	SER	-	expression tag	UNP Q15036
C	106	GLY	-	expression tag	UNP Q15036
C	107	SER	-	expression tag	UNP Q15036

- Molecule 2 is a protein called Krev interaction trapped protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	10	Total	C	N	O	S	0	0	0
			77	51	11	14	1			
2	E	9	Total	C	N	O	S	0	0	0
			71	48	10	12	1			
2	F	9	Total	C	N	O	S	0	0	0
			71	48	10	12	1			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	219	GLY	-	expression tag	UNP O00522
D	220	PRO	-	expression tag	UNP O00522

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Chain	Residue	Modelled	Actual	Comment	Reference
D	221	LEU	-	expression tag	UNP O00522
D	222	GLY	-	expression tag	UNP O00522
D	223	SER	-	expression tag	UNP O00522
D	224	PRO	-	expression tag	UNP O00522
E	219	GLY	-	expression tag	UNP O00522
E	220	PRO	-	expression tag	UNP O00522
E	221	LEU	-	expression tag	UNP O00522
E	222	GLY	-	expression tag	UNP O00522
E	223	SER	-	expression tag	UNP O00522
E	224	PRO	-	expression tag	UNP O00522
F	219	GLY	-	expression tag	UNP O00522
F	220	PRO	-	expression tag	UNP O00522
F	221	LEU	-	expression tag	UNP O00522
F	222	GLY	-	expression tag	UNP O00522
F	223	SER	-	expression tag	UNP O00522
F	224	PRO	-	expression tag	UNP O00522

- Molecule 3 is water.

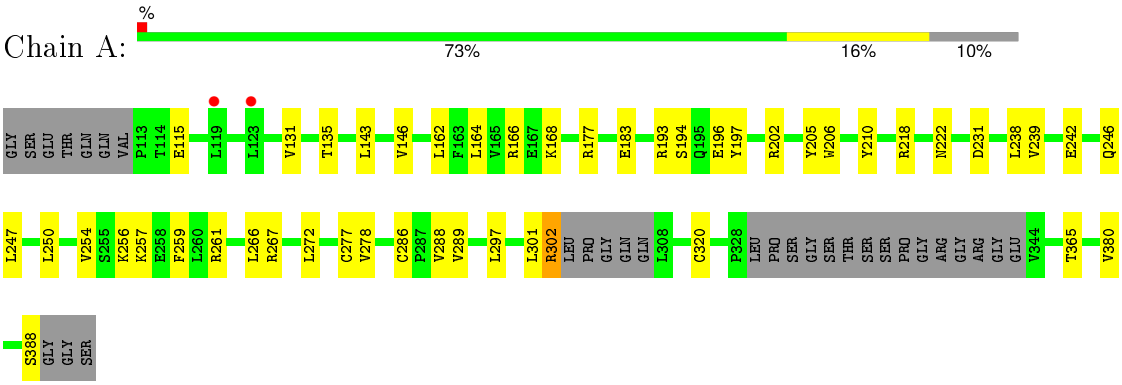
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	15	Total O 15 15	0	0
3	B	17	Total O 17 17	0	0
3	C	16	Total O 16 16	0	0

3 Residue-property plots

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.

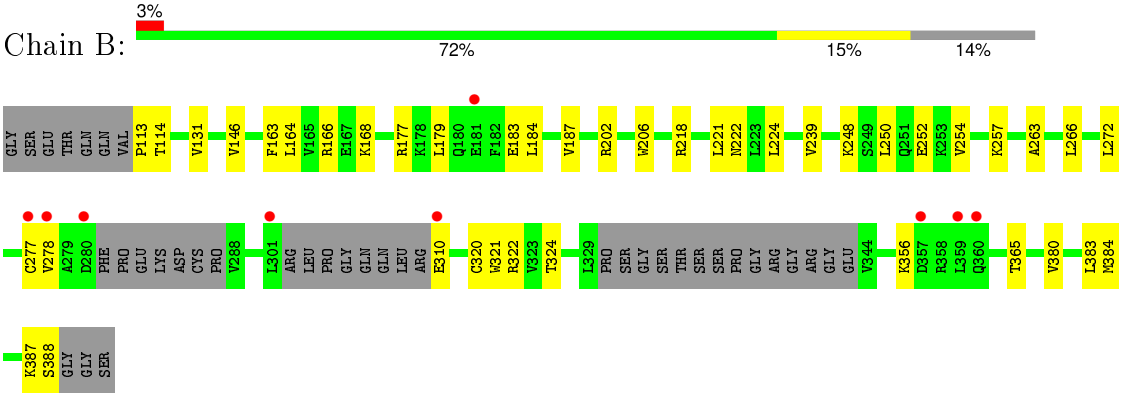
- Molecule 1: Sorting nexin-17

Chain A:



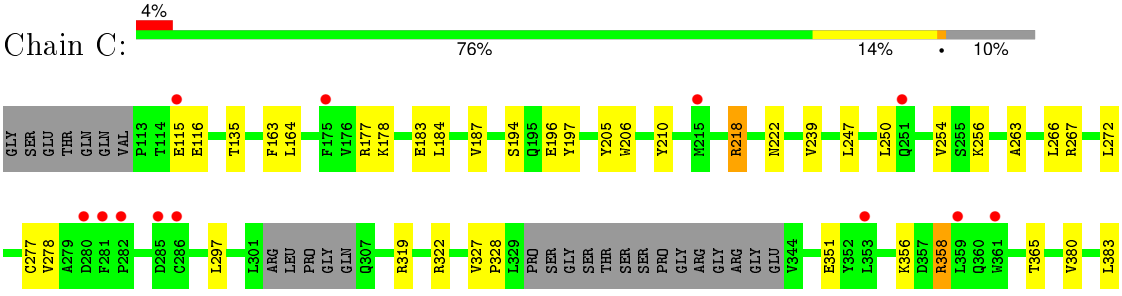
- Molecule 1: Sorting nexin-17

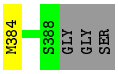
Chain B:



- Molecule 1: Sorting nexin-17

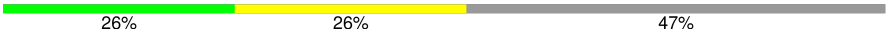
Chain C:





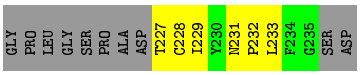
• Molecule 2: Krev interaction trapped protein 1

Chain D:



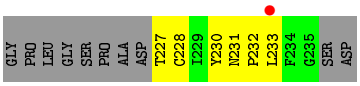
• Molecule 2: Krev interaction trapped protein 1

Chain E:



• Molecule 2: Krev interaction trapped protein 1

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	117.92Å 191.20Å 46.59Å 90.00° 110.25° 90.00°	Depositor
Resolution (Å)	47.88 – 3.00 47.88 – 3.00	Depositor EDS
% Data completeness (in resolution range)	92.5 (47.88-3.00) 92.2 (47.88-3.00)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.65 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.218 , 0.268 0.218 , 0.269	Depositor DCC
R_{free} test set	1793 reflections (10.02%)	DCC
Wilson B-factor (Å^2)	86.8	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 73.7	EDS
Estimated twinning fraction	0.052 for -h-2*l,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	1 of 19141 reflections (0.005%)	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6419	wwPDB-VP
Average B, all atoms (Å^2)	126.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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, 0.375 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	0.32	0/2108	0.61	0/2844
1	B	0.32	0/2025	0.62	0/2731
1	C	0.30	0/2114	0.62	0/2853
2	D	0.28	0/79	0.51	0/107
2	E	0.31	0/73	0.63	0/99
2	F	0.27	0/73	0.58	0/99
All	All	0.31	0/6472	0.62	0/8733

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

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	2075	0	2103	32	1
1	B	1996	0	2024	34	0
1	C	2081	0	2109	31	1
2	D	77	0	72	4	0
2	E	71	0	66	7	0
2	F	71	0	67	5	0
3	A	15	0	0	0	0
3	B	17	0	0	0	0
3	C	16	0	0	0	0
All	All	6419	0	6441	99	1

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 8.

All (99) 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:177:ARG:NH1	1:B:183:GLU:OE1	2.16	0.78
1:C:322:ARG:HB2	2:F:228:CYS:HB3	1.67	0.75
1:B:310:GLU:OE2	1:C:267:ARG:NH2	2.21	0.74
1:B:166:ARG:HH12	1:B:168:LYS:HD2	1.52	0.73
1:C:177:ARG:HH12	1:C:183:GLU:CD	1.97	0.67
1:C:164:LEU:HD11	1:C:197:TYR:HB3	1.77	0.67
1:A:250:LEU:HD13	1:A:259:PHE:HA	1.78	0.65
1:A:218:ARG:O	1:A:222:ASN:ND2	2.33	0.62
1:C:222:ASN:OD1	1:C:256:LYS:NZ	2.33	0.61
1:C:277:CYS:HB2	1:C:365:THR:O	2.02	0.60
1:C:218:ARG:HD2	1:C:222:ASN:HD21	1.67	0.59
2:D:227:THR:OG1	2:D:228:CYS:N	2.35	0.59
1:C:177:ARG:NH1	1:C:183:GLU:OE1	2.35	0.59
2:F:227:THR:OG1	2:F:228:CYS:N	2.35	0.59
1:B:166:ARG:NH1	1:B:168:LYS:HD2	2.17	0.58
1:A:288:VAL:HG12	1:A:301:LEU:HD13	1.86	0.57
1:A:239:VAL:HG12	1:A:266:LEU:HD22	1.87	0.57
1:C:218:ARG:O	1:C:222:ASN:ND2	2.37	0.57
1:A:250:LEU:O	1:A:254:VAL:N	2.37	0.57
1:A:177:ARG:NH1	1:A:183:GLU:OE2	2.34	0.56
1:B:356:LYS:O	1:C:358:ARG:NH2	2.38	0.56
1:C:239:VAL:HG12	1:C:266:LEU:HD22	1.87	0.55
1:B:239:VAL:HG12	1:B:266:LEU:HD22	1.88	0.55
1:B:184:LEU:HB2	1:B:187:VAL:HG22	1.88	0.54
1:B:383:LEU:HB2	2:E:233:LEU:HD11	1.89	0.54
1:C:115:GLU:HG3	1:C:135:THR:HG21	1.89	0.54
1:A:250:LEU:HB2	1:A:259:PHE:HD1	1.72	0.54
1:B:250:LEU:O	1:B:254:VAL:N	2.40	0.53
1:A:206:TRP:HH2	1:A:272:LEU:HD12	1.73	0.53
1:B:321:TRP:CE2	2:E:229:ILE:HD11	2.43	0.53
1:A:143:LEU:HD13	1:A:162:LEU:HG	1.90	0.52
1:A:289:VAL:HG23	1:A:302:ARG:HG2	1.92	0.52
1:A:238:LEU:HD12	1:A:267:ARG:HE	1.74	0.52
1:A:202:ARG:NH2	1:A:231:ASP:OD1	2.35	0.51
1:A:166:ARG:HH12	1:A:168:LYS:HG3	1.73	0.51
1:C:278:VAL:HG22	1:C:365:THR:HB	1.91	0.51
1:A:247:LEU:HD11	1:A:266:LEU:HD11	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:321:TRP:HH2	1:B:380:VAL:HG21	1.74	0.51
1:A:380:VAL:HG13	2:D:233:LEU:HD12	1.93	0.51
1:B:356:LYS:HB2	1:C:358:ARG:HH21	1.76	0.50
2:E:227:THR:OG1	2:E:228:CYS:N	2.32	0.50
1:A:278:VAL:HG22	1:A:365:THR:HB	1.94	0.50
1:C:383:LEU:HB2	2:F:233:LEU:HD11	1.93	0.50
1:A:164:LEU:HD11	1:A:197:TYR:HB3	1.93	0.49
1:C:319:ARG:NH1	1:C:351:GLU:OE1	2.45	0.48
1:A:222:ASN:OD1	1:A:256:LYS:NZ	2.46	0.48
1:C:327:VAL:HG12	1:C:328:PRO:HD2	1.95	0.48
1:A:288:VAL:HG12	1:A:301:LEU:CD1	2.44	0.47
1:C:250:LEU:O	1:C:254:VAL:N	2.45	0.47
1:B:320:CYS:HB2	2:E:229:ILE:O	2.15	0.47
1:B:163:PHE:CE1	1:B:202:ARG:HB3	2.49	0.47
1:A:277:CYS:HB2	1:A:365:THR:O	2.14	0.47
1:B:168:LYS:HE3	1:B:168:LYS:HB3	1.79	0.46
1:B:388:SER:HB3	2:D:230:TYR:HB2	1.97	0.46
1:C:115:GLU:HG3	1:C:135:THR:CG2	2.46	0.46
1:B:322:ARG:HG2	1:B:324:THR:HG23	1.98	0.46
1:B:380:VAL:O	1:B:384:MET:HG2	2.16	0.46
1:B:218:ARG:O	1:B:222:ASN:ND2	2.49	0.46
1:A:286:CYS:SG	1:A:301:LEU:HD12	2.56	0.46
1:C:356:LYS:HD3	1:C:356:LYS:HA	1.57	0.46
1:A:388:SER:HB2	2:E:228:CYS:O	2.17	0.45
1:B:248:LYS:O	1:B:252:GLU:HG3	2.17	0.45
1:C:163:PHE:CE2	1:C:178:LYS:HE2	2.51	0.45
1:A:115:GLU:HG3	1:A:135:THR:CG2	2.48	0.44
1:C:194:SER:HB3	1:C:196:GLU:HG2	1.98	0.44
1:B:221:LEU:HD11	1:B:257:LYS:HG3	2.00	0.44
2:F:231:ASN:HA	2:F:232:PRO:HD2	1.91	0.43
1:C:184:LEU:HB2	1:C:187:VAL:HG22	1.98	0.43
1:A:205:TYR:CD1	1:A:210:TYR:HB2	2.53	0.43
1:C:205:TYR:CD1	1:C:210:TYR:HB2	2.54	0.43
1:C:206:TRP:HH2	1:C:272:LEU:HD12	1.84	0.42
1:B:263:ALA:HA	1:B:266:LEU:HD12	2.00	0.42
1:B:277:CYS:HB2	1:B:365:THR:O	2.19	0.42
1:A:320:CYS:HB2	2:D:229:ILE:O	2.19	0.42
1:A:297:LEU:HA	1:A:297:LEU:HD23	1.82	0.42
1:B:224:LEU:HA	1:B:224:LEU:HD12	1.81	0.42
1:A:242:GLU:O	1:A:246:GLN:HG3	2.19	0.42
1:A:257:LYS:O	1:A:261:ARG:HG2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:VAL:HG11	1:A:146:VAL:HA	2.02	0.41
1:A:194:SER:HB3	1:A:196:GLU:HG2	2.01	0.41
1:C:263:ALA:HA	1:C:266:LEU:CD1	2.50	0.41
1:A:242:GLU:HG3	1:A:246:GLN:OE1	2.20	0.41
1:B:131:VAL:HG11	1:B:146:VAL:HA	2.02	0.41
1:B:184:LEU:O	1:B:187:VAL:HG22	2.21	0.41
1:B:206:TRP:CH2	1:B:272:LEU:HD12	2.56	0.41
1:C:247:LEU:HD11	1:C:266:LEU:HD11	2.03	0.41
1:B:206:TRP:HH2	1:B:272:LEU:HD12	1.86	0.41
1:C:184:LEU:O	1:C:187:VAL:HG22	2.20	0.41
1:B:113:PRO:HB2	1:B:114:THR:H	1.64	0.41
1:C:319:ARG:NH1	2:F:230:TYR:HE1	2.19	0.40
1:B:278:VAL:CG2	1:B:365:THR:HB	2.51	0.40
2:E:231:ASN:HA	2:E:232:PRO:HD2	1.94	0.40
1:A:301:LEU:HD13	1:A:301:LEU:HA	1.84	0.40
1:C:297:LEU:HD23	1:C:297:LEU:HA	1.89	0.40
1:C:380:VAL:O	1:C:384:MET:HG2	2.22	0.40
1:B:278:VAL:HG22	1:B:365:THR:HB	2.04	0.40
1:B:263:ALA:HA	1:B:266:LEU:CD1	2.52	0.40
1:B:384:MET:SD	2:E:233:LEU:HG	2.61	0.40
1:B:164:LEU:HB2	1:B:179:LEU:HD21	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:ARG:NH1	1:C:116:GLU:O[4_556]	2.15	0.05

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	250/286 (87%)	247 (99%)	3 (1%)	0	100	100
1	B	239/286 (84%)	236 (99%)	3 (1%)	0	100	100
1	C	251/286 (88%)	248 (99%)	3 (1%)	0	100	100
2	D	8/19 (42%)	8 (100%)	0	0	100	100
2	E	7/19 (37%)	7 (100%)	0	0	100	100
2	F	7/19 (37%)	7 (100%)	0	0	100	100
All	All	762/915 (83%)	753 (99%)	9 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	236/258 (92%)	235 (100%)	1 (0%)	93	98
1	B	227/258 (88%)	226 (100%)	1 (0%)	93	98
1	C	237/258 (92%)	235 (99%)	2 (1%)	86	96
2	D	9/15 (60%)	9 (100%)	0	100	100
2	E	8/15 (53%)	8 (100%)	0	100	100
2	F	8/15 (53%)	8 (100%)	0	100	100
All	All	725/819 (88%)	721 (99%)	4 (1%)	90	97

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	302	ARG
1	B	387	LYS
1	C	218	ARG
1	C	358	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA

There are no RNA molecules 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, 95^{th} 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	256/286 (89%)	0.03	2 (0%) 87 67	80, 117, 167, 193	0
1	B	247/286 (86%)	0.18	9 (3%) 46 20	83, 117, 169, 192	0
1	C	257/286 (89%)	0.22	12 (4%) 35 14	87, 129, 188, 241	0
2	D	10/19 (52%)	-0.01	0 100 100	107, 122, 140, 152	0
2	E	9/19 (47%)	0.45	0 100 100	104, 122, 133, 135	0
2	F	9/19 (47%)	0.54	1 (11%) 7 3	141, 159, 171, 173	0
All	All	788/915 (86%)	0.15	24 (3%) 54 25	80, 123, 174, 241	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	278	VAL	6.1
1	C	285	ASP	5.5
1	C	280	ASP	4.8
1	B	277	CYS	4.0
1	B	310	GLU	3.9
1	C	286	CYS	3.9
1	C	282	PRO	3.8
1	C	353	LEU	3.0
1	C	115	GLU	3.0
1	C	359	LEU	2.9
1	B	357	ASP	2.8
1	C	281	PHE	2.7
1	A	123	LEU	2.6
1	C	361	TRP	2.5
1	C	175	PHE	2.4
1	B	359	LEU	2.3
1	C	215	MET	2.3
1	B	181	GLU	2.3
1	B	280	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	119	LEU	2.2
1	C	251	GLN	2.1
1	B	301	LEU	2.1
1	B	360	GLN	2.0
2	F	233	LEU	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.