Full wwPDB X-ray Structure Validation Report ⓘ

Sep 13, 2023 – 01:18 AM EDT

PDB ID	:	4OCN
Title	:	Crystal Structure of the Rpn8-Rpn11 MPN domain heterodimer, crystal form II
Authors	:	Pathare, G.R.; Bracher, A.
Deposited on	:	2014-01-09
Resolution	:	2.25 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:

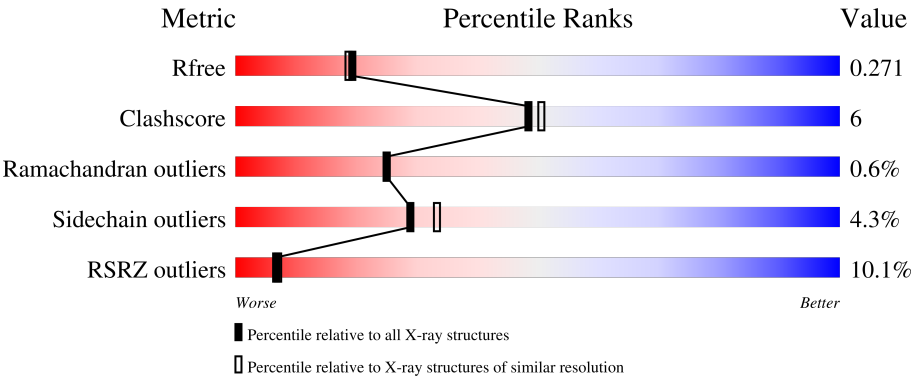
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	187	14% 75% 17% · 7%
1	D	187	12% 75% 15% · 8%
2	B	220	7% 64% 11% 25%
2	E	220	5% 59% 14% · 26%
3	C	133	5% 82% 10% 8%

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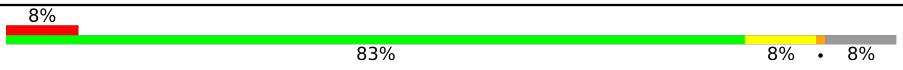
Mol	Chain	Length	Quality of chain
3	F	133	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7341 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 26S proteasome regulatory subunit RPN8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	174	Total	C	N	O	S	0	0	0
			1374	870	232	267	5			
1	D	172	Total	C	N	O	S	0	0	0
			1360	863	231	261	5			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	cloning artifact	UNP Q08723
A	0	HIS	-	cloning artifact	UNP Q08723
A	177	GLY	-	SEE REMARK 999	UNP Q08723
A	178	SER	-	SEE REMARK 999	UNP Q08723
A	179	GLY	-	SEE REMARK 999	UNP Q08723
A	180	GLY	-	SEE REMARK 999	UNP Q08723
A	181	SER	-	SEE REMARK 999	UNP Q08723
A	182	GLY	-	SEE REMARK 999	UNP Q08723
A	183	GLY	-	SEE REMARK 999	UNP Q08723
A	184	SER	-	SEE REMARK 999	UNP Q08723
A	185	GLY	-	SEE REMARK 999	UNP Q08723
D	-1	GLY	-	cloning artifact	UNP Q08723
D	0	HIS	-	cloning artifact	UNP Q08723
D	177	GLY	-	SEE REMARK 999	UNP Q08723
D	178	SER	-	SEE REMARK 999	UNP Q08723
D	179	GLY	-	SEE REMARK 999	UNP Q08723
D	180	GLY	-	SEE REMARK 999	UNP Q08723
D	181	SER	-	SEE REMARK 999	UNP Q08723
D	182	GLY	-	SEE REMARK 999	UNP Q08723
D	183	GLY	-	SEE REMARK 999	UNP Q08723
D	184	SER	-	SEE REMARK 999	UNP Q08723
D	185	GLY	-	SEE REMARK 999	UNP Q08723

- Molecule 2 is a protein called 26S proteasome regulatory subunit RPN11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	165	Total	C	N	O	S	0	0	0
			1296	827	220	237	12			
2	E	162	Total	C	N	O	S	0	0	0
			1265	809	214	230	12			

- Molecule 3 is a protein called Nb1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	122	Total	C	N	O	S	0	0	0
			955	598	173	180	4			
3	F	123	Total	C	N	O	S	0	0	0
			956	601	169	182	4			

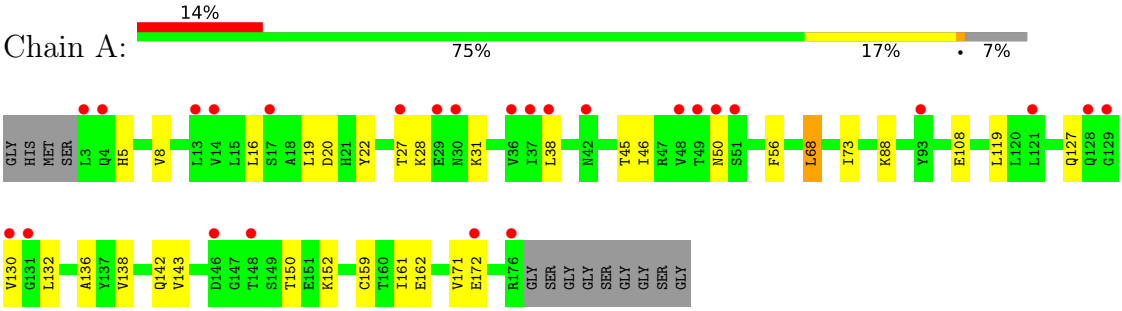
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	47	Total	O	0	0
			47	47		
4	B	23	Total	O	0	0
			23	23		
4	C	20	Total	O	0	0
			20	20		
4	D	16	Total	O	0	0
			16	16		
4	E	29	Total	O	0	0
			29	29		

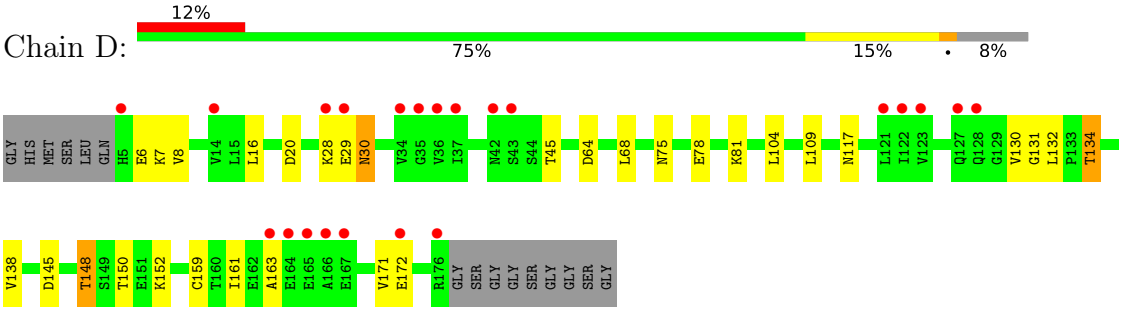
3 Residue-property plots (i)

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.

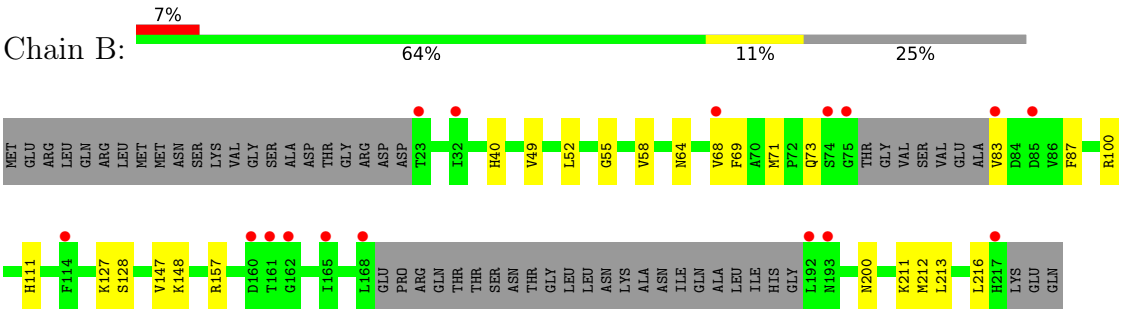
• Molecule 1: 26S proteasome regulatory subunit RPN8



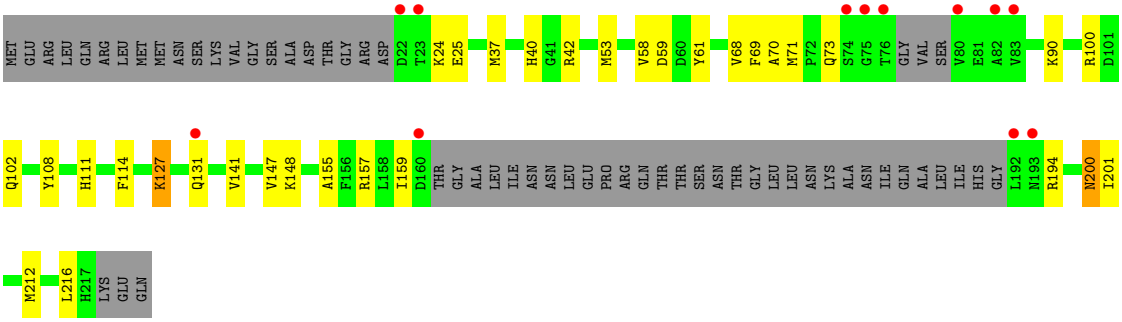
• Molecule 1: 26S proteasome regulatory subunit RPN8



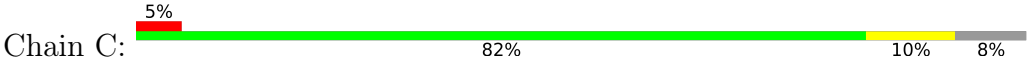
• Molecule 2: 26S proteasome regulatory subunit RPN11

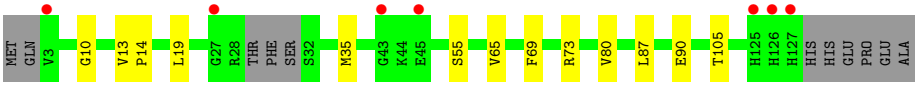


• Molecule 2: 26S proteasome regulatory subunit RPN11

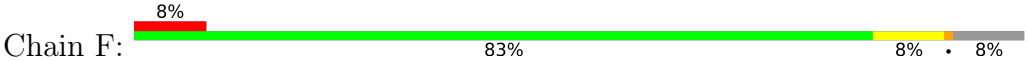


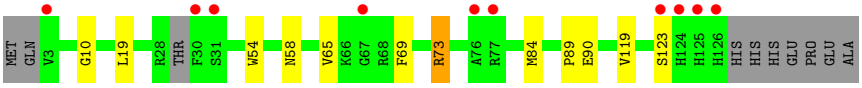
• Molecule 3: Nb1





• Molecule 3: Nb1





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	80.04Å 80.04Å 386.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.25 29.86 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-2.25) 99.7 (29.86-2.25)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	1.86 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.232 , 0.281 0.226 , 0.271	Depositor DCC
R_{free} test set	3108 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å^2)	54.4	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 36.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7341	wwPDB-VP
Average B, all atoms (Å^2)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% 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

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	0.49	0/1399	0.58	0/1896
1	D	0.44	0/1385	0.55	0/1876
2	B	0.51	0/1319	0.66	0/1779
2	E	0.47	0/1288	0.62	0/1738
3	C	0.46	0/979	0.56	0/1324
3	F	0.47	0/980	0.56	0/1326
All	All	0.47	0/7350	0.59	0/9939

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	1374	0	1354	19	0
1	D	1360	0	1353	18	0
2	B	1296	0	1299	19	0
2	E	1265	0	1263	30	0
3	C	955	0	901	7	0
3	F	956	0	897	7	0
4	A	47	0	0	0	0
4	B	23	0	0	0	0
4	C	20	0	0	0	0
4	D	16	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	29	0	0	0	0
All	All	7341	0	7067	86	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 6.

All (86) 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:E:37:MET:HE1	2:E:108:TYR:HB3	1.59	0.83
2:E:37:MET:CE	2:E:108:TYR:HB3	2.09	0.83
1:D:145:ASP:O	1:D:148:THR:HB	1.82	0.79
2:E:69:PHE:HE1	2:E:71:MET:CE	1.95	0.79
1:A:127:GLN:HE22	2:B:212:MET:HA	1.48	0.77
2:E:59:ASP:HB3	2:E:61:TYR:H	1.51	0.74
2:E:155:ALA:HB2	2:E:201:ILE:HD11	1.68	0.74
3:C:65:VAL:HG13	3:C:69:PHE:HB2	1.72	0.71
1:D:68:LEU:HD23	1:D:109:LEU:HD23	1.71	0.71
1:A:68:LEU:HD23	1:A:73:ILE:HD11	1.72	0.71
1:D:8:VAL:HB	1:D:159:CYS:HB3	1.73	0.70
2:B:69:PHE:HE1	2:B:71:MET:CE	2.06	0.68
1:A:143:VAL:HG13	1:A:150:THR:HG22	1.78	0.66
1:A:68:LEU:CD2	1:A:73:ILE:HD11	2.26	0.66
2:E:69:PHE:CE1	2:E:71:MET:CE	2.79	0.64
2:E:37:MET:HE1	2:E:108:TYR:CB	2.30	0.62
1:D:16:LEU:HD23	2:E:212:MET:HE1	1.84	0.60
2:E:100:ARG:HD3	2:E:102:GLN:OE1	2.01	0.60
3:C:35:MET:HG3	3:C:80:VAL:HG21	1.82	0.59
2:B:69:PHE:CE1	2:B:71:MET:HE3	2.37	0.58
1:A:127:GLN:HG2	2:B:211:LYS:HD3	1.86	0.57
2:E:37:MET:HE3	2:E:53:MET:SD	2.45	0.57
2:E:37:MET:CE	2:E:108:TYR:CB	2.81	0.57
2:E:24:LYS:HD2	2:E:200:ASN:HD22	1.69	0.57
2:E:37:MET:HE2	2:E:108:TYR:HB3	1.86	0.56
1:D:78:GLU:HA	1:D:81:LYS:HE2	1.86	0.56
2:B:69:PHE:HE1	2:B:71:MET:HE3	1.72	0.55
2:B:69:PHE:HE1	2:B:71:MET:HE1	1.70	0.55
3:F:89:PRO:HG3	3:F:123:SER:HA	1.89	0.55
1:D:16:LEU:HD23	2:E:212:MET:CE	2.37	0.55
2:B:69:PHE:CE1	2:B:71:MET:CE	2.90	0.54
2:E:111:HIS:HB3	2:E:114:PHE:HB2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:52:LEU:HD12	2:B:69:PHE:CZ	2.41	0.54
3:F:54:TRP:CE3	3:F:73:ARG:HG3	2.44	0.53
1:D:130:VAL:HG12	1:D:131:GLY:H	1.74	0.52
1:A:16:LEU:HD11	2:B:213:LEU:HD21	1.91	0.52
1:D:20:ASP:OD1	2:E:100:ARG:NH2	2.43	0.52
4:D:503:HOH:O	2:E:42:ARG:NH2	2.40	0.52
2:E:37:MET:HE2	2:E:108:TYR:CD1	2.46	0.51
3:C:90:GLU:OE1	3:C:90:GLU:N	2.44	0.51
1:D:16:LEU:HA	2:E:212:MET:HE1	1.93	0.50
2:E:127:LYS:O	2:E:131:GLN:HG2	2.12	0.50
1:D:117:ASN:ND2	1:D:138:VAL:HG13	2.27	0.50
3:F:10:GLY:HA2	3:F:19:LEU:HD13	1.93	0.50
1:D:75:ASN:OD1	2:E:90:LYS:NZ	2.44	0.49
1:A:22:TYR:CE1	1:A:27:THR:HB	2.48	0.49
1:D:7:LYS:HB2	1:D:45:THR:HG22	1.95	0.49
3:F:65:VAL:HG13	3:F:69:PHE:HB2	1.95	0.49
1:A:8:VAL:HB	1:A:159:CYS:HB3	1.95	0.48
1:A:119:LEU:HD11	1:A:136:ALA:HB1	1.95	0.48
1:A:172:GLU:OE1	2:B:148:LYS:HG2	2.14	0.48
3:C:13:VAL:HG11	3:C:87:LEU:HD13	1.96	0.48
2:B:49:VAL:HG12	2:B:73:GLN:NE2	2.29	0.48
2:E:69:PHE:HE1	2:E:71:MET:HE1	1.78	0.47
3:C:13:VAL:HG22	3:C:14:PRO:HD2	1.96	0.47
3:F:58:ASN:HD22	3:F:58:ASN:H	1.63	0.47
2:B:40:HIS:CE1	2:B:49:VAL:HB	2.50	0.46
1:D:29:GLU:O	1:D:30:ASN:CB	2.64	0.46
2:B:71:MET:HE2	2:B:87:PHE:CD1	2.50	0.46
1:A:5:HIS:CG	1:A:46:ILE:HD11	2.51	0.46
1:A:171:VAL:HG11	2:B:216:LEU:HD23	1.98	0.44
1:D:171:VAL:HG21	2:E:216:LEU:O	2.16	0.44
3:C:13:VAL:CG2	3:C:14:PRO:HD2	2.48	0.44
2:E:37:MET:HE2	2:E:108:TYR:CG	2.53	0.44
3:F:84:MET:HE1	3:F:119:VAL:HG21	1.99	0.44
1:A:132:LEU:HD22	1:A:161:ILE:HD12	2.00	0.43
2:E:40:HIS:CE1	2:E:70:ALA:HB1	2.54	0.43
1:A:19:LEU:HD22	2:B:212:MET:HG3	2.00	0.43
3:F:58:ASN:H	3:F:58:ASN:ND2	2.17	0.43
2:E:25:GLU:OE1	2:E:157:ARG:NH2	2.52	0.42
2:B:49:VAL:HG12	2:B:73:GLN:HE22	1.84	0.42
1:A:20:ASP:OD1	2:B:100:ARG:NH2	2.53	0.42
2:B:83:VAL:HG13	2:B:83:VAL:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:PHE:CD1	1:A:68:LEU:HD12	2.55	0.41
2:E:69:PHE:HE1	2:E:71:MET:HE2	1.80	0.41
1:A:108:GLU:OE2	1:A:152:LYS:NZ	2.49	0.41
1:D:104:LEU:HD22	1:D:152:LYS:HG2	2.01	0.41
1:A:38:LEU:HA	1:A:88:LYS:O	2.21	0.41
2:E:69:PHE:CE1	2:E:71:MET:HE2	2.54	0.41
1:D:130:VAL:HG12	1:D:131:GLY:N	2.36	0.41
3:C:10:GLY:HA2	3:C:19:LEU:HD13	2.03	0.40
1:D:132:LEU:HB3	1:D:134:THR:HG22	2.03	0.40
2:E:37:MET:HE2	2:E:108:TYR:CB	2.49	0.40
1:A:38:LEU:HD12	1:A:50:ASN:HB3	2.04	0.40
2:B:55:GLY:HA3	2:B:64:ASN:O	2.20	0.40
1:D:172:GLU:HB2	2:E:148:LYS:HE3	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	172/187 (92%)	165 (96%)	6 (4%)	1 (1%)	25 25
1	D	170/187 (91%)	162 (95%)	6 (4%)	2 (1%)	13 9
2	B	159/220 (72%)	154 (97%)	4 (2%)	1 (1%)	25 25
2	E	156/220 (71%)	150 (96%)	5 (3%)	1 (1%)	25 25
3	C	118/133 (89%)	114 (97%)	4 (3%)	0	100 100
3	F	119/133 (90%)	114 (96%)	5 (4%)	0	100 100
All	All	894/1080 (83%)	859 (96%)	30 (3%)	5 (1%)	25 25

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	130	VAL
1	D	30	ASN
1	D	163	ALA
2	B	58	VAL
2	E	58	VAL

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	151/162 (93%)	144 (95%)	7 (5%)	27	30
1	D	150/162 (93%)	143 (95%)	7 (5%)	26	29
2	B	144/191 (75%)	137 (95%)	7 (5%)	25	27
2	E	139/191 (73%)	131 (94%)	8 (6%)	20	20
3	C	99/109 (91%)	96 (97%)	3 (3%)	41	50
3	F	99/109 (91%)	97 (98%)	2 (2%)	55	64
All	All	782/924 (85%)	748 (96%)	34 (4%)	29	33

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LYS
1	A	31	LYS
1	A	45	THR
1	A	68	LEU
1	A	138	VAL
1	A	142	GLN
1	A	162	GLU
2	B	68	VAL
2	B	111	HIS
2	B	127	LYS
2	B	128	SER
2	B	147	VAL
2	B	157	ARG
2	B	200	ASN

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Mol	Chain	Res	Type
3	C	55	SER
3	C	73	ARG
3	C	105	THR
1	D	6	GLU
1	D	28	LYS
1	D	64	ASP
1	D	134	THR
1	D	148	THR
1	D	150	THR
1	D	161	ILE
2	E	68	VAL
2	E	73	GLN
2	E	127	LYS
2	E	141	VAL
2	E	147	VAL
2	E	159	ILE
2	E	194	ARG
2	E	200	ASN
3	F	73	ARG
3	F	90	GLU

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	77	ASN
2	B	124	ASN
2	B	133	ASN
2	B	215	ASN
2	B	217	HIS
3	C	40	GLN
3	C	78	ASN
1	D	71	ASN
2	E	73	GLN
2	E	109	HIS
2	E	200	ASN
3	F	58	ASN
3	F	75	ASN

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 monosaccharides 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	174/187 (93%)	0.63	26 (14%) 2 2	34, 52, 82, 101	0
1	D	172/187 (91%)	0.61	22 (12%) 3 3	40, 58, 88, 107	0
2	B	165/220 (75%)	0.28	16 (9%) 7 8	31, 44, 80, 116	0
2	E	162/220 (73%)	0.24	12 (7%) 14 15	37, 51, 78, 89	0
3	C	122/133 (91%)	0.02	7 (5%) 23 25	36, 54, 81, 97	0
3	F	123/133 (92%)	0.34	10 (8%) 12 13	39, 57, 80, 100	0
All	All	918/1080 (85%)	0.37	93 (10%) 7 7	31, 53, 83, 116	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	3	LEU	7.4
1	D	166	ALA	6.6
2	B	75	GLY	6.0
2	E	192	LEU	5.1
1	A	30	ASN	5.0
3	F	30	PHE	5.0
2	B	192	LEU	4.8
3	F	123	SER	4.8
1	D	5	HIS	4.8
2	E	23	THR	4.7
1	A	29	GLU	4.7
1	A	129	GLY	4.6
2	E	76	THR	4.5
3	F	126	HIS	4.3
3	F	124	HIS	4.1
1	A	130	VAL	4.0
3	F	3	VAL	3.9
1	A	27	THR	3.8
2	B	217	HIS	3.8

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Mol	Chain	Res	Type	RSRZ
3	F	31	SER	3.7
1	D	42	ASN	3.6
3	C	127	HIS	3.6
2	E	193	ASN	3.6
1	D	123	VAL	3.5
2	B	160	ASP	3.5
1	A	13	LEU	3.4
2	E	22	ASP	3.4
1	A	14	VAL	3.4
3	C	126	HIS	3.4
2	B	161	THR	3.3
2	B	23	THR	3.2
2	E	160	ASP	3.2
1	A	37	ILE	3.2
2	E	75	GLY	3.2
3	F	125	HIS	3.2
1	A	176	ARG	3.2
1	A	146	ASP	3.0
1	D	128	GLN	3.0
1	D	163	ALA	3.0
2	B	165	ILE	3.0
3	C	3	VAL	3.0
1	D	176	ARG	3.0
1	D	165	GLU	2.9
2	E	82	ALA	2.9
1	A	128	GLN	2.9
2	B	162	GLY	2.9
1	D	127	GLN	2.9
1	A	51	SER	2.9
2	E	83	VAL	2.9
2	B	193	ASN	2.9
1	A	4	GLN	2.8
1	D	164	GLU	2.8
2	B	74	SER	2.8
3	C	27	GLY	2.8
3	F	76	ALA	2.7
1	A	42	ASN	2.7
2	B	168	LEU	2.6
1	D	28	LYS	2.6
1	D	172	GLU	2.6
2	E	80	VAL	2.6
1	A	50	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	43	SER	2.5
3	F	77	ARG	2.5
1	A	38	LEU	2.5
1	A	93	TYR	2.5
2	B	114	PHE	2.5
1	D	37	ILE	2.5
2	B	32	ILE	2.5
1	A	131	GLY	2.5
1	A	36	VAL	2.5
3	C	45	GLU	2.4
1	D	121	LEU	2.4
1	D	167	GLU	2.4
3	F	67	GLY	2.4
1	D	34	VAL	2.4
2	E	74	SER	2.4
2	B	68	VAL	2.3
3	C	43	GLY	2.3
1	A	49	THR	2.3
1	A	121	LEU	2.2
1	D	35	GLY	2.2
3	C	125	HIS	2.2
1	D	122	ILE	2.2
1	D	14	VAL	2.1
1	D	29	GLU	2.1
2	B	83	VAL	2.1
1	A	17	SER	2.1
2	B	85	ASP	2.1
1	A	172	GLU	2.1
1	D	36	VAL	2.1
2	E	131	GLN	2.1
1	A	48	VAL	2.1
1	A	148	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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.