



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:17 PM GMT

PDB ID : 1SYX
Title : The crystal structure of a binary U5 snRNP complex
Authors : Nielsen, T.K.; Liu, S.; Luhrmann, R.; Ficner, R.
Deposited on : 2004-04-02
Resolution : 2.35 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
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

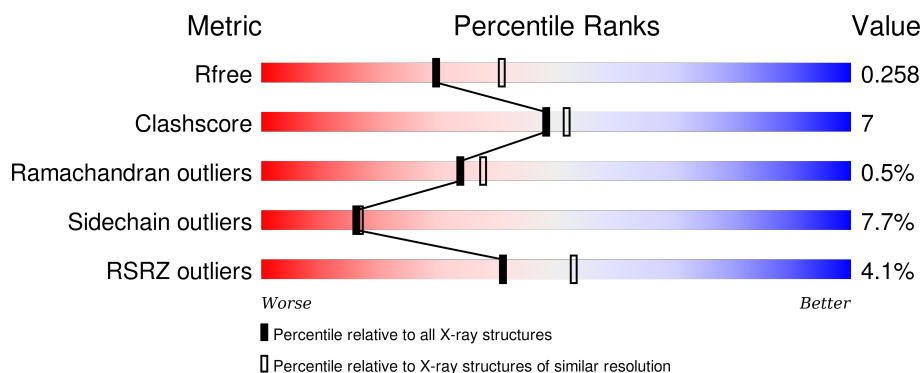
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1406 (2.36-2.32)
Clashscore	102246	1509 (2.36-2.32)
Ramachandran outliers	100387	1490 (2.36-2.32)
Sidechain outliers	100360	1491 (2.36-2.32)
RSRZ outliers	91569	1412 (2.36-2.32)

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	142	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 74%, yellow 16%, orange 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 74% 16% • • 5% </div> </div>
1	C	142	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, green 72%, yellow 20%, orange 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 72% 20% • • 5% </div> </div>
1	E	142	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 73%, yellow 18%, orange 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 73% 18% • 5% </div> </div>
2	B	86	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 5%, green 55%, yellow 15%, orange 28%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 5% 55% 15% • 28% </div> </div>
2	D	86	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 9%, green 47%, yellow 24%, orange 28%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 9% 47% 24% • 28% </div> </div>

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Mol	Chain	Length	Quality of chain	
				
2	F	86		

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5082 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 Spliceosomal U5 snRNP-specific 15 kDa protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	135	Total	C	N	O	S	16	0	0
			1112	714	184	204	10			
1	C	135	Total	C	N	O	S	18	0	0
			1112	714	184	204	10			
1	E	135	Total	C	N	O	S	11	0	0
			1112	714	184	204	10			

- Molecule 2 is a protein called CD2 antigen cytoplasmic tail-binding protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	62	Total	C	N	O	S	0	0	0
			518	335	79	101	3			
2	D	62	Total	C	N	O	S	7	0	0
			518	335	79	101	3			
2	F	62	Total	C	N	O	S	5	0	0
			518	335	79	101	3			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	ALA	-	CLONING ARTIFACT	UNP O95400
B	2	GLU	-	CLONING ARTIFACT	UNP O95400
B	3	GLU	-	CLONING ARTIFACT	UNP O95400
B	4	GLU	-	CLONING ARTIFACT	UNP O95400
B	5	LEU	-	CLONING ARTIFACT	UNP O95400
B	6	GLU	-	CLONING ARTIFACT	UNP O95400
D	1	ALA	-	CLONING ARTIFACT	UNP O95400
D	2	GLU	-	CLONING ARTIFACT	UNP O95400
D	3	GLU	-	CLONING ARTIFACT	UNP O95400
D	4	GLU	-	CLONING ARTIFACT	UNP O95400
D	5	LEU	-	CLONING ARTIFACT	UNP O95400
D	6	GLU	-	CLONING ARTIFACT	UNP O95400

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Chain	Residue	Modelled	Actual	Comment	Reference
F	1	ALA	-	CLONING ARTIFACT	UNP O95400
F	2	GLU	-	CLONING ARTIFACT	UNP O95400
F	3	GLU	-	CLONING ARTIFACT	UNP O95400
F	4	GLU	-	CLONING ARTIFACT	UNP O95400
F	5	LEU	-	CLONING ARTIFACT	UNP O95400
F	6	GLU	-	CLONING ARTIFACT	UNP O95400

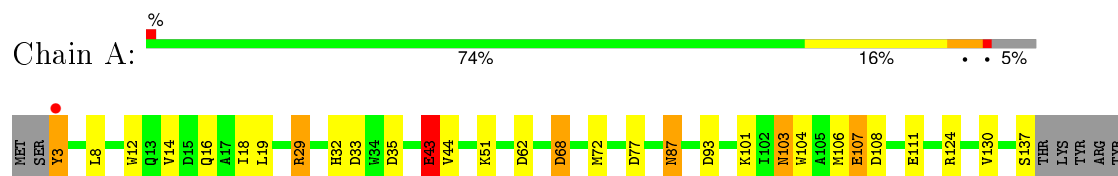
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	66	Total O 66 66	0	0
3	B	19	Total O 19 19	0	0
3	C	35	Total O 35 35	0	0
3	D	21	Total O 21 21	0	0
3	E	36	Total O 36 36	0	0
3	F	15	Total O 15 15	0	0

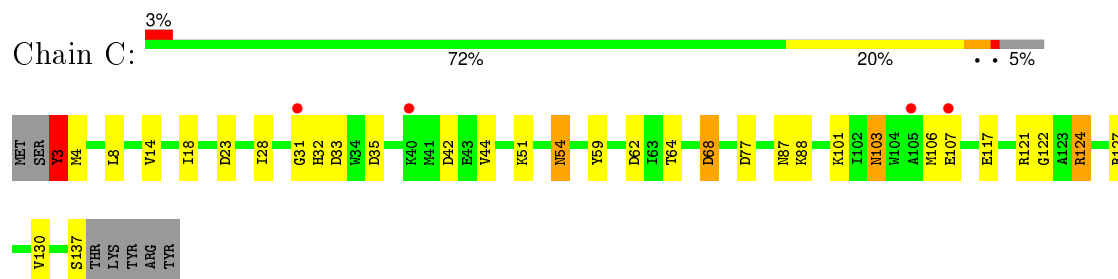
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

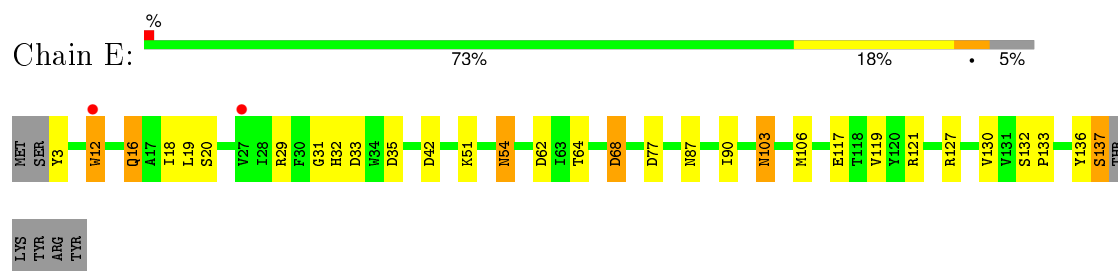
- Molecule 1: Spliceosomal U5 snRNP-specific 15 kDa protein



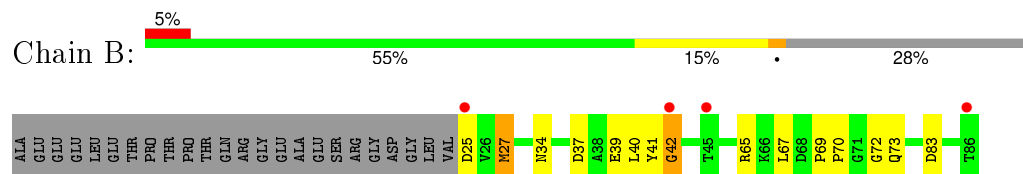
- Molecule 1: Spliceosomal U5 snRNP-specific 15 kDa protein



- Molecule 1: Spliceosomal U5 snRNP-specific 15 kDa protein

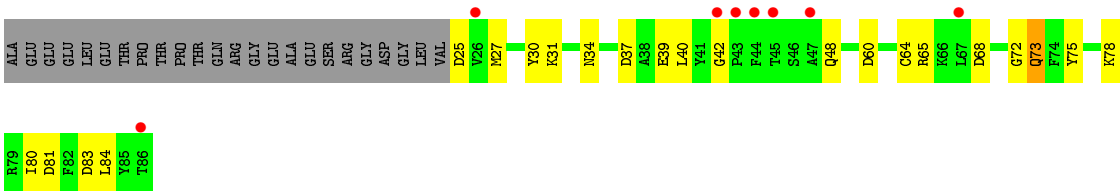


- Molecule 2: CD2 antigen cytoplasmic tail-binding protein 2

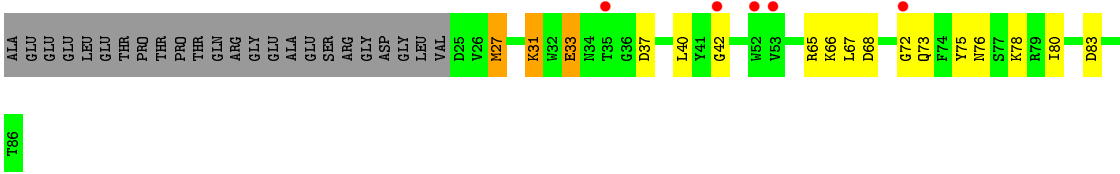


- Molecule 2: CD2 antigen cytoplasmic tail-binding protein 2





• Molecule 2: CD2 antigen cytoplasmic tail-binding protein 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	135.73 Å 75.91 Å 77.03 Å 90.00° 94.27° 90.00°	Depositor
Resolution (Å)	49.39 – 2.35 49.27 – 2.35	Depositor EDS
% Data completeness (in resolution range)	91.8 (49.39-2.35) 91.8 (49.27-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.34 Å)	Xtriage
Refinement program	REFMAC 5.1	Depositor
R, R_{free}	0.216 , 0.262 0.214 , 0.258	Depositor DCC
R_{free} test set	1486 reflections (5.18%)	DCC
Wilson B-factor (Å ²)	51.6	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 30195 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5082	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% 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, 0.375 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	2.64	6/1139 (0.5%)	2.26	18/1541 (1.2%)
1	C	0.85	5/1139 (0.4%)	1.18	15/1541 (1.0%)
1	E	2.29	7/1139 (0.6%)	2.42	18/1541 (1.2%)
2	B	0.44	0/537	0.73	0/729
2	D	0.66	1/537 (0.2%)	0.89	6/729 (0.8%)
2	F	0.86	2/537 (0.4%)	0.99	5/729 (0.7%)
All	All	1.75	21/5028 (0.4%)	1.75	62/6810 (0.9%)

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	2
1	C	0	1
1	E	0	1
2	F	0	1
All	All	0	5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	43	GLU	CD-OE2	57.18	1.88	1.25
1	E	12	TRP	CE2-CZ2	55.88	2.34	1.39
1	E	12	TRP	CG-CD2	43.27	2.17	1.43
1	A	43	GLU	CD-OE1	-43.03	0.78	1.25
1	A	3	TYR	CD2-CE2	-33.10	0.89	1.39
1	A	3	TYR	CB-CG	28.84	1.95	1.51
1	A	3	TYR	CE1-CZ	-24.38	1.06	1.38
1	E	12	TRP	CZ2-CH2	20.09	1.75	1.37
1	C	3	TYR	CB-CG	15.16	1.74	1.51
1	E	12	TRP	CD2-CE3	-13.67	1.19	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	124	ARG	NE-CZ	-13.44	1.15	1.33
2	F	33	GLU	CD-OE1	13.43	1.40	1.25
2	D	60	ASP	CB-CG	11.34	1.75	1.51
1	C	88	LYS	CD-CE	10.26	1.76	1.51
2	F	33	GLU	CD-OE2	10.03	1.36	1.25
1	E	16	GLN	CB-CG	8.59	1.75	1.52
1	E	54	ASN	CG-OD1	8.43	1.42	1.24
1	A	107	GLU	CB-CG	-7.38	1.38	1.52
1	E	12	TRP	CE3-CZ3	-7.18	1.26	1.38
1	C	54	ASN	CG-ND2	6.75	1.49	1.32
1	C	54	ASN	CG-OD1	-6.57	1.09	1.24

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	12	TRP	CH2-CZ2-CE2	-48.96	68.44	117.40
1	A	3	TYR	CB-CG-CD1	-43.77	94.74	121.00
1	A	3	TYR	CB-CG-CD2	43.49	147.09	121.00
1	E	12	TRP	CD1-CG-CD2	-39.97	74.32	106.30
1	A	3	TYR	CD1-CE1-CZ	-36.66	86.81	119.80
1	E	12	TRP	CZ3-CH2-CZ2	27.37	154.45	121.60
1	E	12	TRP	CD2-CE3-CZ3	27.22	154.18	118.80
1	E	12	TRP	CG-CD1-NE1	27.22	137.32	110.10
1	A	3	TYR	CE1-CZ-CE2	20.01	151.82	119.80
1	E	12	TRP	CB-CG-CD1	19.36	152.16	127.00
1	A	43	GLU	CG-CD-OE2	-18.86	80.59	118.30
1	E	12	TRP	CB-CG-CD2	-18.50	102.55	126.60
1	E	12	TRP	CG-CD2-CE3	17.77	149.90	133.90
1	E	12	TRP	CE3-CZ3-CH2	-16.78	102.75	121.20
1	A	3	TYR	CE1-CZ-OH	-16.68	75.08	120.10
1	A	43	GLU	OE1-CD-OE2	-16.68	103.29	123.30
1	C	124	ARG	NE-CZ-NH1	-16.42	112.09	120.30
1	C	124	ARG	NE-CZ-NH2	15.38	127.99	120.30
1	C	3	TYR	CB-CG-CD1	-14.42	112.35	121.00
1	E	12	TRP	CE2-CD2-CE3	-14.25	101.60	118.70
2	F	33	GLU	OE1-CD-OE2	-14.23	106.22	123.30
1	C	3	TYR	CA-CB-CG	-14.16	86.50	113.40
1	A	43	GLU	CG-CD-OE1	12.72	143.75	118.30
1	E	12	TRP	NE1-CE2-CZ2	-11.55	117.69	130.40
1	E	54	ASN	OD1-CG-ND2	-11.50	95.45	121.90
1	C	3	TYR	CB-CG-CD2	10.55	127.33	121.00
1	C	3	TYR	N-CA-CB	10.18	128.93	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	TYR	CA-CB-CG	-9.86	94.66	113.40
1	E	12	TRP	CD2-CE2-CZ2	9.57	133.78	122.30
2	D	25	ASP	N-CA-CB	9.34	127.40	110.60
2	D	25	ASP	CB-CA-C	-8.71	92.99	110.40
1	C	4	MET	CA-CB-CG	8.57	127.87	113.30
1	A	33	ASP	CB-CG-OD2	8.43	125.89	118.30
2	D	25	ASP	CA-CB-CG	-7.42	97.07	113.40
2	F	37	ASP	CB-CG-OD1	-6.88	112.11	118.30
2	F	37	ASP	CB-CG-OD2	6.78	124.40	118.30
1	E	68	ASP	CB-CG-OD2	6.47	124.12	118.30
1	A	3	TYR	CG-CD2-CE2	-6.46	116.14	121.30
1	A	124	ARG	NE-CZ-NH2	-6.35	117.12	120.30
1	A	68	ASP	CB-CG-OD2	6.30	123.97	118.30
1	A	3	TYR	CZ-CE2-CD2	-6.30	114.13	119.80
1	E	77	ASP	CB-CG-OD2	6.29	123.96	118.30
1	A	77	ASP	CB-CG-OD2	5.91	123.62	118.30
1	C	33	ASP	CB-CG-OD2	5.88	123.60	118.30
1	E	35	ASP	CB-CG-OD2	5.71	123.44	118.30
1	C	68	ASP	CB-CG-OD2	5.67	123.40	118.30
1	C	23	ASP	CB-CG-OD2	5.61	123.35	118.30
1	C	42	ASP	CB-CG-OD2	5.47	123.22	118.30
1	C	77	ASP	CB-CG-OD2	5.44	123.20	118.30
1	E	42	ASP	CB-CG-OD2	5.41	123.17	118.30
1	A	62	ASP	CB-CG-OD2	5.41	123.17	118.30
2	D	37	ASP	CB-CG-OD2	5.32	123.08	118.30
2	F	33	GLU	CG-CD-OE2	-5.28	107.74	118.30
1	A	93	ASP	CB-CG-OD2	5.26	123.03	118.30
1	E	33	ASP	CB-CG-OD2	5.22	123.00	118.30
1	C	35	ASP	CB-CG-OD2	5.20	122.98	118.30
1	C	3	TYR	N-CA-C	-5.15	97.10	111.00
2	D	68	ASP	CB-CG-OD2	5.12	122.91	118.30
2	F	68	ASP	CB-CG-OD2	5.12	122.91	118.30
1	C	124	ARG	CD-NE-CZ	5.11	130.75	123.60
1	A	35	ASP	CB-CG-OD2	5.07	122.86	118.30
2	D	81	ASP	CB-CG-OD2	5.02	122.81	118.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	3	TYR	Sidechain
1	A	43	GLU	Sidechain

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Mol	Chain	Res	Type	Group
1	C	3	TYR	Sidechain
1	E	54	ASN	Sidechain
2	F	33	GLU	Sidechain

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	1112	0	1084	10	0
1	C	1112	0	1085	16	0
1	E	1112	0	1085	14	0
2	B	518	0	461	11	0
2	D	518	0	461	11	0
2	F	518	0	461	10	0
3	A	66	0	0	1	0
3	B	19	0	0	2	0
3	C	35	0	0	2	0
3	D	21	0	0	1	0
3	E	36	0	0	2	0
3	F	15	0	0	0	0
All	All	5082	0	4637	69	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 7.

All (69) 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:B:39:GLU:OE1	3:B:87:HOH:O	1.71	1.09
2:D:39:GLU:OE1	3:D:89:HOH:O	1.89	0.90
2:F:65:ARG:HD2	2:F:72:GLY:HA2	1.52	0.89
2:F:27:MET:HB3	2:F:67:LEU:HD12	1.58	0.84
2:B:40:LEU:HB2	2:D:39:GLU:HG3	1.59	0.83
1:C:103:ASN:H	1:C:103:ASN:HD22	1.26	0.83
1:A:32:HIS:HD2	3:A:153:HOH:O	1.65	0.80
2:B:65:ARG:HH11	2:B:72:GLY:HA3	1.53	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:65:ARG:HD2	2:D:72:GLY:HA2	1.70	0.72
2:F:65:ARG:HD2	2:F:72:GLY:CA	2.20	0.71
2:B:41:TYR:O	2:B:42:GLY:O	2.08	0.71
1:C:44:VAL:HG21	1:C:107:GLU:HA	1.74	0.67
2:F:76:ASN:OD1	2:F:78:LYS:HB2	1.97	0.64
1:E:3:TYR:N	3:E:175:HOH:O	2.31	0.64
2:F:72:GLY:O	2:F:73:GLN:HB3	1.98	0.63
2:D:65:ARG:HH11	2:D:72:GLY:HA3	1.64	0.63
2:B:70:PRO:HB3	2:D:30:TYR:OH	1.98	0.63
1:E:103:ASN:H	1:E:103:ASN:HD22	1.47	0.63
1:A:108:ASP:HB3	1:A:111:GLU:HB2	1.80	0.63
1:C:103:ASN:HD22	1:C:103:ASN:N	1.94	0.61
2:D:72:GLY:O	2:D:73:GLN:HB3	2.01	0.60
1:E:32:HIS:HD2	3:E:143:HOH:O	1.85	0.60
1:C:18:ILE:HG22	1:C:87:ASN:HD22	1.69	0.57
2:D:64:CYS:SG	2:D:80:ILE:HD12	2.46	0.56
1:C:106:MET:O	1:C:106:MET:HG3	2.07	0.55
2:F:65:ARG:HH11	2:F:72:GLY:HA3	1.71	0.54
1:C:103:ASN:H	1:C:103:ASN:ND2	2.01	0.54
2:B:65:ARG:HD2	2:B:72:GLY:HA2	1.90	0.54
2:D:31:LYS:HG3	2:D:40:LEU:HD23	1.89	0.54
1:C:8:LEU:HD13	1:C:14:VAL:HA	1.89	0.54
2:B:72:GLY:O	2:B:73:GLN:HB3	2.09	0.52
1:E:117:GLU:OE2	1:E:121:ARG:NH2	2.43	0.51
1:C:28:ILE:HD13	1:C:59:TYR:HB2	1.92	0.51
1:C:44:VAL:CG2	1:C:107:GLU:HA	2.42	0.50
1:C:32:HIS:HE1	1:C:64:THR:OG1	1.94	0.50
1:A:103:ASN:HD22	1:A:103:ASN:H	1.60	0.49
1:C:122:GLY:HA2	2:D:84:LEU:HD22	1.94	0.49
1:A:18:ILE:HG22	1:A:87:ASN:HD22	1.77	0.49
2:F:75:TYR:HB2	2:F:80:ILE:HD11	1.95	0.49
2:B:27:MET:HB3	2:B:67:LEU:HD12	1.94	0.49
1:C:103:ASN:ND2	1:C:103:ASN:N	2.61	0.48
1:E:90:ILE:HG21	1:E:119:VAL:HG13	1.95	0.48
1:E:136:TYR:O	1:E:137:SER:HB3	2.13	0.48
1:C:54:ASN:HB2	3:C:152:HOH:O	2.12	0.48
1:E:32:HIS:HE1	1:E:64:THR:OG1	1.96	0.48
2:D:75:TYR:HB2	2:D:80:ILE:HD11	1.97	0.47
1:A:29:ARG:HD2	1:A:29:ARG:C	2.34	0.47
2:B:69:PRO:HB2	2:B:70:PRO:HD2	1.96	0.47
1:E:18:ILE:HG22	1:E:87:ASN:HD22	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:VAL:HG21	1:A:107:GLU:HA	1.98	0.46
1:C:3:TYR:HB3	3:C:174:HOH:O	2.15	0.46
1:C:117:GLU:OE2	1:C:121:ARG:NH2	2.50	0.45
2:F:65:ARG:NH1	2:F:72:GLY:HA3	2.32	0.44
2:F:66:LYS:HE3	2:F:66:LYS:HB2	1.73	0.44
1:C:31:GLY:O	1:C:62:ASP:HA	2.18	0.43
1:E:32:HIS:CE1	1:E:64:THR:OG1	2.72	0.43
1:E:29:ARG:HD2	1:E:29:ARG:C	2.39	0.42
1:E:16:GLN:O	1:E:20:SER:HB3	2.19	0.42
2:D:65:ARG:NH1	2:D:72:GLY:HA3	2.32	0.42
2:F:31:LYS:HG3	2:F:40:LEU:HD23	2.02	0.42
1:E:106:MET:O	1:E:106:MET:HG3	2.18	0.42
1:A:12:TRP:O	1:A:16:GLN:HG2	2.19	0.42
2:B:65:ARG:NH1	2:B:72:GLY:HA3	2.29	0.42
1:E:31:GLY:O	1:E:62:ASP:HA	2.19	0.41
1:A:106:MET:HG3	1:A:106:MET:O	2.20	0.41
1:A:103:ASN:HD22	1:A:104:TRP:HE3	1.68	0.41
2:B:39:GLU:HG2	3:B:95:HOH:O	2.21	0.41
1:A:8:LEU:HD13	1:A:14:VAL:HA	2.02	0.40
1:E:132:SER:HA	1:E:133:PRO:HD3	1.99	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	133/142 (94%)	131 (98%)	2 (2%)	0	100	100
1	C	133/142 (94%)	132 (99%)	1 (1%)	0	100	100
1	E	133/142 (94%)	133 (100%)	0	0	100	100
2	B	60/86 (70%)	57 (95%)	2 (3%)	1 (2%)	11	8

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	60/86 (70%)	57 (95%)	2 (3%)	1 (2%)	11	8
2	F	60/86 (70%)	58 (97%)	1 (2%)	1 (2%)	11	8
All	All	579/684 (85%)	568 (98%)	8 (1%)	3 (0%)	34	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	42	GLY
2	D	42	GLY
2	F	42	GLY

5.3.2 Protein sidechains ⓘ

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	123/130 (95%)	112 (91%)	11 (9%)	12	11
1	C	123/130 (95%)	115 (94%)	8 (6%)	21	24
1	E	123/130 (95%)	115 (94%)	8 (6%)	21	24
2	B	54/73 (74%)	49 (91%)	5 (9%)	11	10
2	D	54/73 (74%)	48 (89%)	6 (11%)	8	7
2	F	54/73 (74%)	51 (94%)	3 (6%)	26	31
All	All	531/609 (87%)	490 (92%)	41 (8%)	16	17

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	29	ARG
1	A	43	GLU
1	A	51	LYS
1	A	68	ASP
1	A	72	MET
1	A	87	ASN

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Mol	Chain	Res	Type
1	A	101	LYS
1	A	103	ASN
1	A	130	VAL
1	A	137	SER
1	C	51	LYS
1	C	68	ASP
1	C	101	LYS
1	C	103	ASN
1	C	124	ARG
1	C	127	ARG
1	C	130	VAL
1	C	137	SER
1	E	12	TRP
1	E	19	LEU
1	E	51	LYS
1	E	68	ASP
1	E	103	ASN
1	E	127	ARG
1	E	130	VAL
1	E	137	SER
2	B	25	ASP
2	B	27	MET
2	B	34	ASN
2	B	37	ASP
2	B	83	ASP
2	D	27	MET
2	D	34	ASN
2	D	48	GLN
2	D	73	GLN
2	D	78	LYS
2	D	83	ASP
2	F	27	MET
2	F	31	LYS
2	F	83	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	HIS
1	A	87	ASN
1	A	89	HIS
1	A	103	ASN

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Mol	Chain	Res	Type
1	C	32	HIS
1	C	87	ASN
1	C	89	HIS
1	C	103	ASN
1	E	32	HIS
1	E	87	ASN
1	E	89	HIS
1	E	103	ASN
2	B	34	ASN
2	D	34	ASN
2	D	50	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	135/142 (95%)	0.45	1 (0%) 89 93	34, 39, 46, 58	5 (3%)
1	C	135/142 (95%)	0.39	4 (2%) 54 64	34, 40, 47, 58	5 (3%)
1	E	135/142 (95%)	0.54	2 (1%) 76 84	35, 40, 45, 56	3 (2%)
2	B	62/86 (72%)	0.57	4 (6%) 22 33	36, 40, 43, 44	0
2	D	62/86 (72%)	0.90	8 (12%) 5 8	37, 41, 43, 44	2 (3%)
2	F	62/86 (72%)	0.81	5 (8%) 15 23	35, 39, 44, 45	2 (3%)
All	All	591/684 (86%)	0.56	24 (4%) 41 53	34, 40, 45, 58	17 (2%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	86	THR	4.1
2	D	42	GLY	3.7
2	F	72	GLY	3.5
2	F	35	THR	3.4
2	D	86	THR	3.2
1	A	3	TYR	3.0
2	B	42	GLY	2.9
1	C	107	GLU	2.9
2	D	47	ALA	2.9
2	B	25	ASP	2.9
2	D	67	LEU	2.8
2	D	44	PHE	2.8
2	F	42	GLY	2.8
1	E	12	TRP	2.8
2	D	43	PRO	2.6
2	F	53	VAL	2.6
2	D	45	THR	2.5
1	C	31	GLY	2.4
1	C	40	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	45	THR	2.3
2	D	26	VAL	2.2
1	C	105	ALA	2.2
2	F	52	TRP	2.2
1	E	27	VAL	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 carbohydrates 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.