



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:58 PM GMT

PDB ID : 4UF6
Title : UCH-L5 in complex with ubiquitin-propargyl bound to an activating fragment of INO80G
Authors : Sahtoe, D.D.; Van Dijk, W.J.; El Oualid, F.; Ekkebus, R.; Ovaa, H.; Sixma, T.K.
Deposited on : 2014-12-23
Resolution : 3.69 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

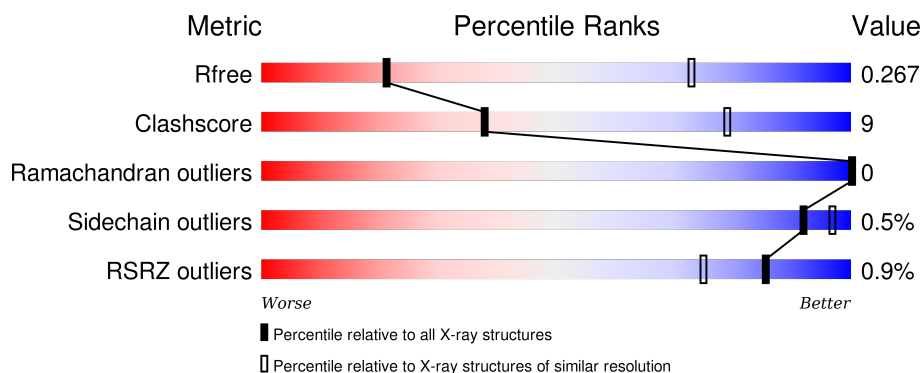
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.69 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	1101 (3.90-3.50)
Clashscore	102246	1224 (3.90-3.50)
Ramachandran outliers	100387	1172 (3.90-3.50)
Sidechain outliers	100360	1170 (3.90-3.50)
RSRZ outliers	91569	1108 (3.90-3.50)

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	331	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 66%, yellow 66%, yellow 88%, grey 88%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 66% 22% 12% </div> </div>
1	D	331	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 0%, green 69%, yellow 69%, yellow 89%, grey 89%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 69% 20% 11% </div> </div>
1	G	331	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, green 1%, green 66%, yellow 66%, yellow 87%, grey 87%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 66% 21% 13% </div> </div>
1	J	331	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 0%, green 68%, yellow 68%, yellow 89%, grey 89%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 68% 21% 11% </div> </div>
2	B	76	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 4%, green 4%, green 80%, yellow 80%, yellow 98%, grey 98%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 80% 18% . </div> </div>

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Mol	Chain	Length	Quality of chain
2	E	76	
2	H	76	
2	K	76	
3	C	65	
3	F	65	
3	I	65	
3	L	65	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13486 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	0	0
			2360	1503	405	440	12			
1	D	295	Total	C	N	O	S	0	0	0
			2382	1515	409	446	12			
1	G	287	Total	C	N	O	S	0	0	0
			2313	1475	397	430	11			
1	J	294	Total	C	N	O	S	0	0	0
			2377	1514	407	444	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q9Y5K5
A	-1	PRO	-	EXPRESSION TAG	UNP Q9Y5K5
A	0	GLY	-	EXPRESSION TAG	UNP Q9Y5K5
D	-2	GLY	-	EXPRESSION TAG	UNP Q9Y5K5
D	-1	PRO	-	EXPRESSION TAG	UNP Q9Y5K5
D	0	GLY	-	EXPRESSION TAG	UNP Q9Y5K5
G	-2	GLY	-	EXPRESSION TAG	UNP Q9Y5K5
G	-1	PRO	-	EXPRESSION TAG	UNP Q9Y5K5
G	0	GLY	-	EXPRESSION TAG	UNP Q9Y5K5
J	-2	GLY	-	EXPRESSION TAG	UNP Q9Y5K5
J	-1	PRO	-	EXPRESSION TAG	UNP Q9Y5K5
J	0	GLY	-	EXPRESSION TAG	UNP Q9Y5K5

- Molecule 2 is a protein called POLYUBIQUITIN-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	76	Total	C	N	O	S	0	0	0
			601	379	105	116	1			
2	E	76	Total	C	N	O	S	0	0	0
			601	379	105	116	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	76	Total	C	N	O	S	0	0	0
			601	379	105	116	1			
2	K	76	Total	C	N	O	S	0	0	0
			601	379	105	116	1			

- Molecule 3 is a protein called NUCLEAR FACTOR RELATED TO KAPPA-B-BINDING PROTEIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	50	Total	C	N	O	0	0	0
			417	267	64	86			
3	F	50	Total	C	N	O	0	0	0
			417	267	64	86			
3	I	49	Total	C	N	O	0	0	0
			408	262	63	83			
3	L	49	Total	C	N	O	0	0	0
			408	262	63	83			

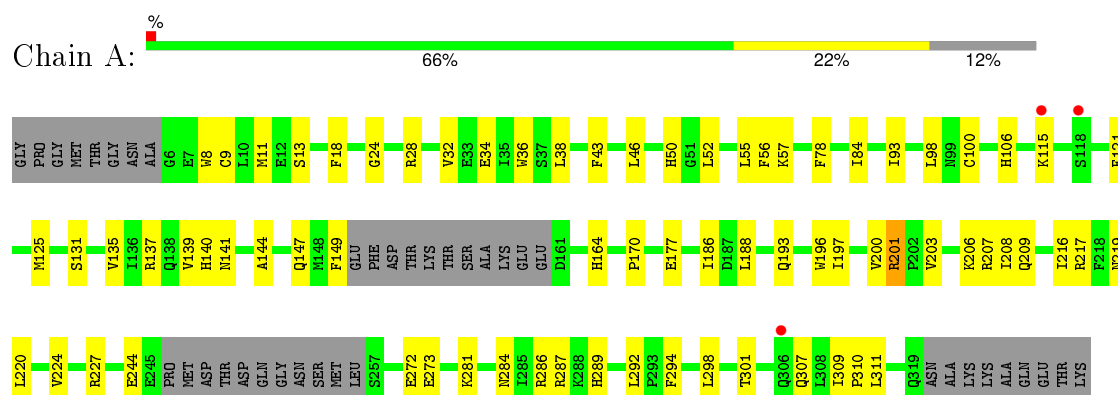
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	37	GLY	-	EXPRESSION TAG	UNP Q6P4R8
C	38	PRO	-	EXPRESSION TAG	UNP Q6P4R8
C	39	GLY	-	EXPRESSION TAG	UNP Q6P4R8
F	37	GLY	-	EXPRESSION TAG	UNP Q6P4R8
F	38	PRO	-	EXPRESSION TAG	UNP Q6P4R8
F	39	GLY	-	EXPRESSION TAG	UNP Q6P4R8
I	37	GLY	-	EXPRESSION TAG	UNP Q6P4R8
I	38	PRO	-	EXPRESSION TAG	UNP Q6P4R8
I	39	GLY	-	EXPRESSION TAG	UNP Q6P4R8
L	37	GLY	-	EXPRESSION TAG	UNP Q6P4R8
L	38	PRO	-	EXPRESSION TAG	UNP Q6P4R8
L	39	GLY	-	EXPRESSION TAG	UNP Q6P4R8

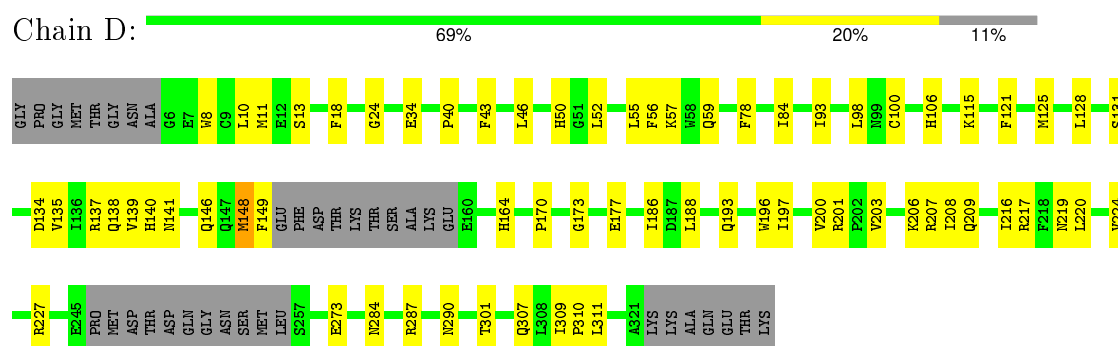
3 Residue-property plots [i](#)

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.

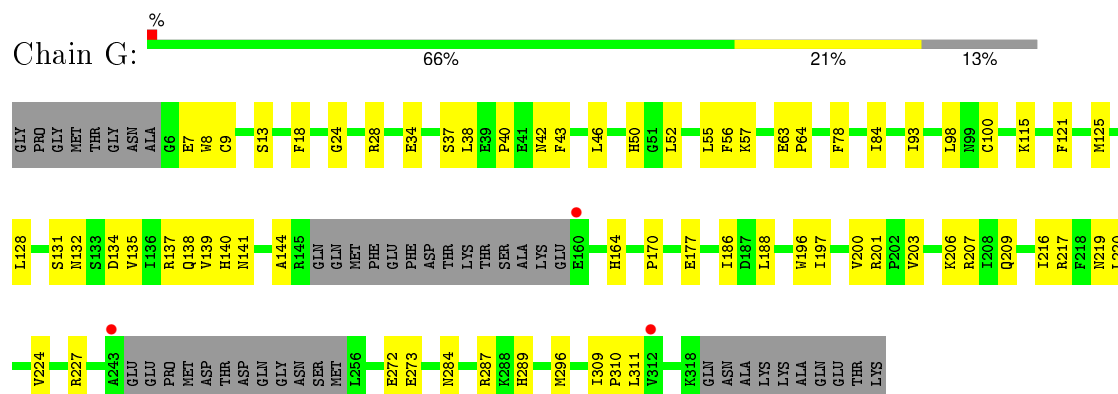
• Molecule 1:



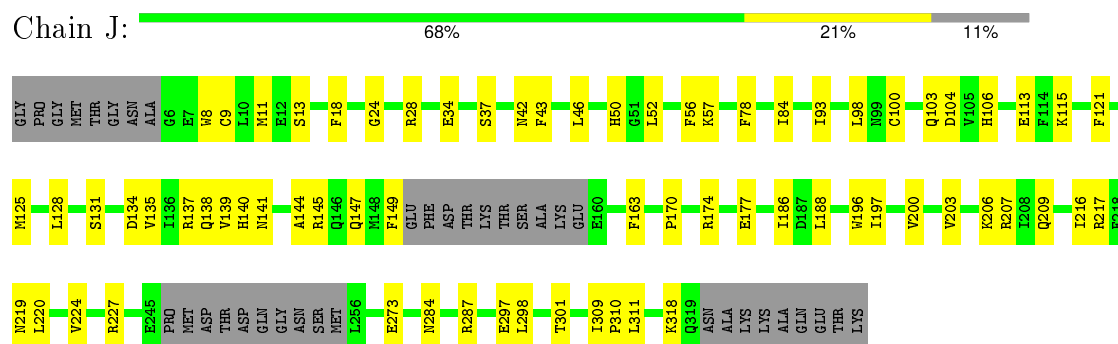
• Molecule 1:



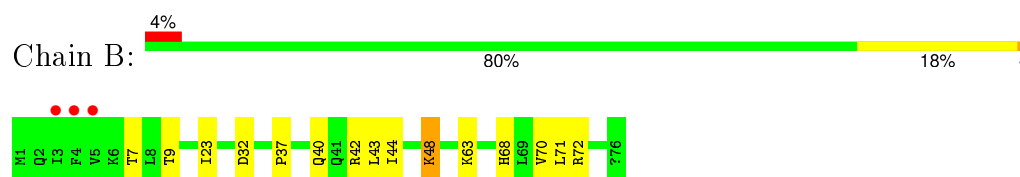
• Molecule 1:



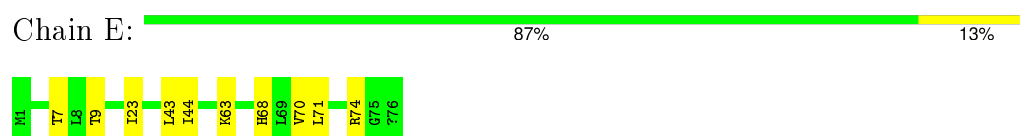
- Molecule 1:



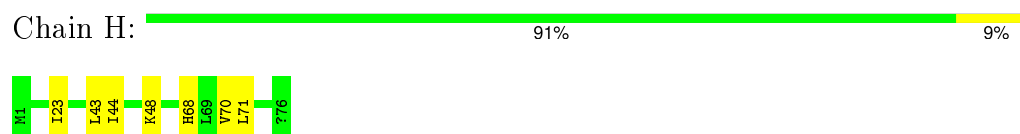
- Molecule 2: POLYUBIQUITIN-B



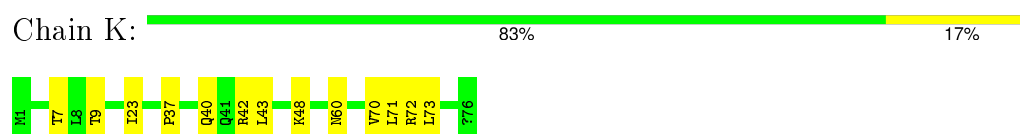
- Molecule 2: POLYUBIQUITIN-B



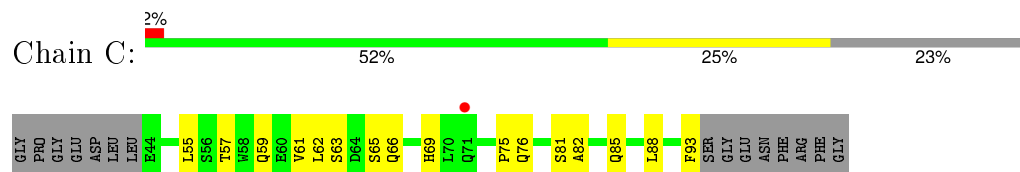
- Molecule 2: POLYUBIQUITIN-B



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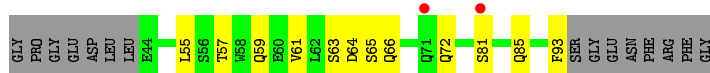


- Molecule 3: NUCLEAR FACTOR RELATED TO KAPPA-B-BINDING PROTEIN



- Molecule 3: NUCLEAR FACTOR RELATED TO KAPPA-B-BINDING PROTEIN





● Molecule 3: NUCLEAR FACTOR RELATED TO KAPPA-B-BINDING PROTEIN



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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	152.08 Å 137.79 Å 98.92 Å 90.00° 102.54° 90.00°	Depositor
Resolution (Å)	100.99 – 3.69 47.74 – 3.69	Depositor EDS
% Data completeness (in resolution range)	97.9 (100.99-3.69) 98.0 (47.74-3.69)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 3.67 Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.238 , 0.269 0.240 , 0.267	Depositor DCC
R_{free} test set	1059 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	82.8	Xtriage
Anisotropy	0.568	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 29.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 21089 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	13486	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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

Bond lengths and bond angles in the following residue types are not validated in this section: AYE

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	0.44	0/2404	0.64	2/3236 (0.1%)
1	D	0.44	0/2426	0.62	0/3266
1	G	0.41	0/2356	0.59	0/3173
1	J	0.46	0/2421	0.62	0/3259
2	B	0.34	0/603	0.57	0/811
2	E	0.39	0/603	0.61	0/811
2	H	0.35	0/603	0.57	0/811
2	K	0.35	0/603	0.57	0/811
3	C	0.43	0/427	0.63	0/581
3	F	0.45	0/427	0.68	0/581
3	I	0.43	0/418	0.63	0/569
3	L	0.46	0/418	0.63	0/569
All	All	0.43	0/13709	0.61	2/18478 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	201	ARG	NE-CZ-NH1	7.37	123.99	120.30
1	A	201	ARG	NE-CZ-NH2	-6.35	117.13	120.30

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	2360	0	2348	62	1
1	D	2382	0	2365	54	0
1	G	2313	0	2311	53	0
1	J	2377	0	2365	50	1
2	B	601	0	630	12	2
2	E	601	0	630	10	1
2	H	601	0	630	6	0
2	K	601	0	630	8	1
3	C	417	0	383	17	0
3	F	417	0	383	11	0
3	I	408	0	377	7	1
3	L	408	0	377	11	1
All	All	13486	0	13429	247	4

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 9.

All (247) 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:G:40:PRO:HB3	1:G:201:ARG:HH22	1.20	1.02
1:D:135:VAL:O	1:D:139:VAL:HG23	1.63	0.99
1:A:135:VAL:O	1:A:139:VAL:HG23	1.62	0.98
1:J:135:VAL:O	1:J:139:VAL:HG23	1.62	0.97
1:G:135:VAL:O	1:G:139:VAL:HG23	1.62	0.97
1:G:40:PRO:HB3	1:G:201:ARG:NH2	1.99	0.78
1:G:38:LEU:O	1:G:201:ARG:HD3	1.90	0.71
1:A:43:PHE:CG	1:A:197:ILE:HG21	2.27	0.70
1:D:43:PHE:CG	1:D:197:ILE:HG21	2.27	0.69
1:J:43:PHE:CG	1:J:197:ILE:HG21	2.27	0.69
1:G:43:PHE:CG	1:G:197:ILE:HG21	2.28	0.68
1:D:121:PHE:HB3	1:D:125:MET:HB2	1.75	0.68
1:G:121:PHE:HB3	1:G:125:MET:HB2	1.76	0.67
1:J:121:PHE:HB3	1:J:125:MET:HB2	1.76	0.66
1:A:34:GLU:CD	2:B:70:VAL:HG11	2.16	0.66
1:A:121:PHE:HB3	1:A:125:MET:HB2	1.76	0.66
1:J:34:GLU:CD	2:K:70:VAL:HG11	2.16	0.65
1:D:311:LEU:HD13	3:F:57:THR:HG21	1.78	0.64
1:D:40:PRO:HD3	1:D:201:ARG:HH12	1.62	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:301:THR:HG21	3:L:65:SER:HB2	1.81	0.63
1:D:34:GLU:CD	2:E:70:VAL:HG11	2.21	0.61
1:G:40:PRO:CB	1:G:201:ARG:HH22	2.03	0.61
1:G:40:PRO:HD3	1:G:201:ARG:HH12	1.65	0.61
3:F:55:LEU:O	3:F:59:GLN:NE2	2.34	0.60
3:C:55:LEU:O	3:C:59:GLN:NE2	2.34	0.60
1:D:106:HIS:CD2	1:G:28:ARG:HB3	2.37	0.60
1:J:217:ARG:HA	2:K:71:LEU:CD1	2.31	0.59
1:J:217:ARG:HA	2:K:71:LEU:HD12	1.83	0.59
3:L:55:LEU:O	3:L:59:GLN:NE2	2.35	0.59
1:G:34:GLU:CD	2:H:70:VAL:HG11	2.23	0.59
1:D:217:ARG:HA	2:E:71:LEU:CD1	2.34	0.58
3:I:55:LEU:O	3:I:59:GLN:NE2	2.36	0.58
1:D:307:GLN:OE1	3:F:61:VAL:HG11	2.04	0.57
1:G:272:GLU:OE1	2:H:48:LYS:NZ	2.37	0.57
1:D:98:LEU:HD22	1:D:115:LYS:HB2	1.87	0.57
1:A:36:TRP:CE3	2:B:44:ILE:HD13	2.39	0.57
1:A:272:GLU:CD	2:B:48:LYS:HE3	2.26	0.56
1:J:301:THR:HG21	3:L:65:SER:CB	2.36	0.56
1:G:24:GLY:HA3	1:G:139:VAL:HG11	1.88	0.56
1:G:98:LEU:HD22	1:G:115:LYS:HB2	1.87	0.56
1:D:50:HIS:HB3	1:D:100:CYS:SG	2.46	0.55
1:A:98:LEU:HD22	1:A:115:LYS:HB2	1.87	0.55
1:D:24:GLY:HA3	1:D:139:VAL:HG11	1.87	0.55
1:A:24:GLY:HA3	1:A:139:VAL:HG11	1.89	0.55
1:J:98:LEU:HD22	1:J:115:LYS:HB2	1.87	0.55
1:D:217:ARG:HA	2:E:71:LEU:HD12	1.89	0.54
1:J:24:GLY:HA3	1:J:139:VAL:HG11	1.88	0.54
1:A:38:LEU:O	1:A:201:ARG:HD2	2.08	0.54
1:A:50:HIS:HB3	1:A:100:CYS:SG	2.48	0.54
1:G:50:HIS:HB3	1:G:100:CYS:SG	2.48	0.54
1:J:50:HIS:HB3	1:J:100:CYS:SG	2.48	0.53
1:J:147:GLN:OE1	1:J:149:PHE:CE1	2.62	0.53
1:G:63:GLU:HG2	1:G:64:PRO:HD2	1.91	0.53
1:D:84:ILE:HD11	1:D:131:SER:HB3	1.91	0.53
3:C:82:ALA:HB2	3:F:64:ASP:OD2	2.10	0.52
1:J:216:ILE:O	1:J:216:ILE:HG12	2.10	0.52
1:A:84:ILE:HD11	1:A:131:SER:HB3	1.92	0.52
1:D:106:HIS:NE2	1:G:28:ARG:HB3	2.24	0.52
1:J:103:GLN:HG2	1:J:104:ASP:N	2.26	0.51
1:D:146:GLN:HE22	1:D:290:ASN:HA	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:216:ILE:HG12	1:D:216:ILE:O	2.10	0.51
1:G:84:ILE:HD11	1:G:131:SER:HB3	1.93	0.51
1:A:170:PRO:C	1:A:193:GLN:HE22	2.13	0.51
1:G:311:LEU:HD13	3:I:57:THR:HG21	1.91	0.51
1:G:216:ILE:HG12	1:G:216:ILE:O	2.10	0.51
1:D:56:PHE:CZ	1:D:207:ARG:NH2	2.79	0.50
1:J:84:ILE:HD11	1:J:131:SER:HB3	1.93	0.50
1:A:298:LEU:HA	3:C:69:HIS:CE1	2.47	0.50
1:A:216:ILE:HG12	1:A:216:ILE:O	2.11	0.50
1:G:46:LEU:HB3	1:G:224:VAL:HG11	1.94	0.49
1:A:38:LEU:O	1:A:201:ARG:CD	2.60	0.49
1:D:309:ILE:HB	1:D:310:PRO:HD3	1.93	0.49
1:D:146:GLN:NE2	1:D:290:ASN:HA	2.27	0.49
1:J:46:LEU:HB3	1:J:224:VAL:HG11	1.94	0.49
1:A:311:LEU:HD13	3:C:57:THR:HG21	1.94	0.49
1:D:170:PRO:HB2	1:D:193:GLN:OE1	2.13	0.49
1:A:56:PHE:CZ	1:A:207:ARG:NH2	2.81	0.49
1:A:309:ILE:HB	1:A:310:PRO:HD3	1.94	0.49
3:I:81:SER:O	3:I:85:GLN:HG2	2.13	0.49
1:G:78:PHE:CE2	1:G:98:LEU:HD12	2.47	0.48
1:J:298:LEU:HA	3:L:69:HIS:CE1	2.47	0.48
1:D:137:ARG:O	1:D:141:ASN:ND2	2.46	0.48
3:L:81:SER:O	3:L:85:GLN:HG2	2.14	0.48
3:C:81:SER:O	3:C:85:GLN:HG2	2.13	0.48
1:J:188:LEU:HD23	1:J:203:VAL:HG21	1.95	0.48
1:J:56:PHE:CZ	1:J:207:ARG:NH2	2.81	0.48
1:A:307:GLN:OE1	3:C:61:VAL:HG11	2.14	0.48
1:J:309:ILE:HB	1:J:310:PRO:HD3	1.93	0.48
1:D:46:LEU:HB3	1:D:224:VAL:HG11	1.96	0.48
1:G:137:ARG:O	1:G:141:ASN:ND2	2.47	0.48
1:J:318:LYS:NZ	3:L:52:VAL:HG13	2.28	0.48
1:A:121:PHE:CZ	1:G:296:MET:HE3	2.49	0.48
1:A:177:GLU:HB3	1:A:186:ILE:HB	1.96	0.48
1:A:188:LEU:HD23	1:A:203:VAL:HG21	1.96	0.48
1:D:301:THR:HG21	3:F:65:SER:HB2	1.94	0.48
1:A:46:LEU:HB3	1:A:224:VAL:HG11	1.95	0.48
1:A:137:ARG:O	1:A:141:ASN:ND2	2.46	0.48
1:J:177:GLU:HB3	1:J:186:ILE:HB	1.96	0.47
1:A:78:PHE:CE1	1:A:98:LEU:HD12	2.49	0.47
1:G:56:PHE:CZ	1:G:207:ARG:NH2	2.83	0.47
1:D:188:LEU:HD23	1:D:203:VAL:HG21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:309:ILE:HB	1:G:310:PRO:HD3	1.95	0.47
1:G:177:GLU:HB3	1:G:186:ILE:HB	1.97	0.47
1:D:173:GLY:H	1:D:193:GLN:HG2	1.79	0.47
1:D:177:GLU:HB3	1:D:186:ILE:HB	1.96	0.47
3:F:81:SER:O	3:F:85:GLN:HG2	2.14	0.47
1:G:18:PHE:CD2	1:G:93:ILE:HD11	2.50	0.47
1:A:289:HIS:CG	3:C:75:PRO:HB3	2.50	0.47
1:J:18:PHE:CD2	1:J:93:ILE:HD11	2.50	0.47
1:A:18:PHE:CD2	1:A:93:ILE:HD11	2.50	0.46
3:C:63:SER:O	3:C:66:GLN:HG2	2.16	0.46
1:G:188:LEU:HD23	1:G:203:VAL:HG21	1.96	0.46
1:J:137:ARG:O	1:J:141:ASN:ND2	2.48	0.46
1:J:78:PHE:CE1	1:J:98:LEU:HD12	2.50	0.46
3:F:63:SER:O	3:F:66:GLN:HG2	2.16	0.46
2:H:23:ILE:HG23	2:H:43:LEU:HD12	1.98	0.46
1:J:297:GLU:HG3	3:L:73:PHE:CZ	2.50	0.46
1:J:52:LEU:HD12	1:J:200:VAL:HG11	1.98	0.46
3:F:55:LEU:HA	3:F:93:PHE:HB3	1.97	0.46
3:L:63:SER:O	3:L:66:GLN:HG2	2.16	0.46
1:D:52:LEU:HB3	1:D:220:LEU:HD11	1.98	0.46
3:C:55:LEU:HA	3:C:93:PHE:HB3	1.97	0.45
1:G:52:LEU:HB3	1:G:220:LEU:HD11	1.98	0.45
1:D:52:LEU:HD12	1:D:200:VAL:HG11	1.99	0.45
1:A:106:HIS:HE1	1:J:28:ARG:HB3	1.81	0.45
1:A:28:ARG:HD2	1:J:106:HIS:HB3	1.99	0.45
1:G:52:LEU:HD12	1:G:200:VAL:HG11	1.99	0.45
1:D:78:PHE:CE1	1:D:98:LEU:HD12	2.52	0.45
1:A:217:ARG:HA	2:B:71:LEU:CD1	2.46	0.45
1:J:163:PHE:CZ	2:K:73:LEU:HD22	2.52	0.45
1:A:170:PRO:HG3	1:A:196:TRP:CD1	2.52	0.45
1:A:52:LEU:HD12	1:A:200:VAL:HG11	1.98	0.44
2:K:23:ILE:HG23	2:K:43:LEU:HD12	1.99	0.44
1:J:170:PRO:HG3	1:J:196:TRP:CD1	2.53	0.44
2:B:23:ILE:HG23	2:B:43:LEU:HD12	1.98	0.44
1:D:148:MET:N	1:D:148:MET:SD	2.91	0.44
3:I:63:SER:O	3:I:66:GLN:HG2	2.16	0.44
3:L:80:ASP:O	3:L:84:GLN:HG2	2.18	0.44
1:D:13:SER:OG	1:D:219:ASN:ND2	2.50	0.44
3:F:63:SER:HB2	3:F:65:SER:OG	2.18	0.44
3:I:63:SER:HB2	3:I:65:SER:OG	2.18	0.44
1:G:8:TRP:HE3	1:G:140:HIS:ND1	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:LEU:HB3	1:A:220:LEU:HD11	1.98	0.44
1:J:52:LEU:HB3	1:J:220:LEU:HD11	1.99	0.44
1:A:147:GLN:OE1	1:G:7:GLU:HG2	2.18	0.44
2:E:23:ILE:HG23	2:E:43:LEU:HD12	1.98	0.44
1:A:281:LYS:NZ	1:J:113:GLU:OE2	2.47	0.44
2:K:7:THR:OG1	2:K:9:THR:OG1	2.29	0.44
1:D:196:TRP:HZ3	1:D:200:VAL:HG21	1.83	0.43
3:C:63:SER:HB2	3:C:65:SER:OG	2.18	0.43
1:D:8:TRP:HE3	1:D:140:HIS:ND1	2.16	0.43
1:D:18:PHE:CD2	1:D:93:ILE:HD11	2.52	0.43
1:J:9:CYS:HB3	1:J:144:ALA:HB2	2.00	0.43
1:J:196:TRP:HZ3	1:J:200:VAL:HG21	1.83	0.43
3:I:80:ASP:O	3:I:84:GLN:HG2	2.18	0.43
1:A:8:TRP:HE3	1:A:140:HIS:ND1	2.16	0.43
1:J:219:ASN:OD1	1:J:219:ASN:C	2.57	0.43
1:A:196:TRP:HZ3	1:A:200:VAL:HG21	1.84	0.43
3:L:63:SER:HB2	3:L:65:SER:OG	2.18	0.43
1:A:84:ILE:HD11	1:A:131:SER:CB	2.49	0.43
1:G:196:TRP:HZ3	1:G:200:VAL:HG21	1.83	0.43
1:J:13:SER:OG	1:J:219:ASN:ND2	2.51	0.43
1:G:284:ASN:OD1	1:G:287:ARG:NH1	2.51	0.43
1:A:286:ARG:HD3	3:C:88:LEU:HD13	2.00	0.43
2:H:44:ILE:HB	2:H:68:HIS:HB2	2.01	0.43
1:D:227:ARG:HB2	1:D:273:GLU:CD	2.39	0.43
1:G:13:SER:OG	1:G:219:ASN:ND2	2.51	0.43
1:A:311:LEU:HD22	3:C:57:THR:HG22	2.00	0.43
1:A:284:ASN:OD1	1:A:287:ARG:NH1	2.52	0.43
1:A:292:LEU:HD23	1:G:132:ASN:HA	2.01	0.43
1:D:284:ASN:OD1	1:D:287:ARG:NH1	2.51	0.43
2:B:44:ILE:HB	2:B:68:HIS:HB2	2.00	0.42
1:D:84:ILE:HD11	1:D:131:SER:CB	2.49	0.42
2:E:44:ILE:HB	2:E:68:HIS:HB2	1.99	0.42
1:A:140:HIS:CD2	1:A:140:HIS:C	2.93	0.42
1:G:227:ARG:HB2	1:G:273:GLU:CD	2.39	0.42
1:D:301:THR:HG21	3:F:65:SER:CB	2.50	0.42
1:G:170:PRO:HG3	1:G:196:TRP:CD1	2.54	0.42
1:A:9:CYS:HB3	1:A:144:ALA:HB2	2.01	0.42
1:D:10:LEU:CB	2:E:74:ARG:HG2	2.49	0.42
1:D:208:ILE:HG21	2:E:9:THR:HG22	2.00	0.42
1:J:206:LYS:HA	1:J:209:GLN:HG2	2.00	0.42
1:J:284:ASN:OD1	1:J:287:ARG:NH1	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:MET:CE	1:A:140:HIS:CD2	3.03	0.42
1:J:128:LEU:O	1:J:131:SER:N	2.53	0.42
1:A:311:LEU:HD22	3:C:57:THR:CG2	2.49	0.42
1:D:10:LEU:CD1	2:E:74:ARG:HG2	2.49	0.42
1:D:219:ASN:OD1	1:D:219:ASN:C	2.58	0.42
1:A:227:ARG:HB2	1:A:273:GLU:CD	2.39	0.42
1:G:57:LYS:HD3	1:G:217:ARG:HH21	1.85	0.42
1:D:206:LYS:HA	1:D:209:GLN:HG2	2.01	0.42
1:D:170:PRO:HG3	1:D:196:TRP:CD1	2.54	0.42
1:A:149:PHE:N	1:A:149:PHE:CD1	2.88	0.42
1:A:28:ARG:HB3	1:J:106:HIS:CG	2.55	0.41
1:G:206:LYS:HA	1:G:209:GLN:HG2	2.01	0.41
1:A:36:TRP:CE3	2:B:44:ILE:CD1	3.03	0.41
1:G:84:ILE:HD11	1:G:131:SER:CB	2.49	0.41
1:A:217:ARG:HA	2:B:71:LEU:HD12	2.02	0.41
1:G:9:CYS:HB3	1:G:144:ALA:HB2	2.02	0.41
1:A:206:LYS:HA	1:A:209:GLN:HG2	2.02	0.41
2:B:7:THR:OG1	2:B:9:THR:OG1	2.29	0.41
1:D:149:PHE:N	1:D:149:PHE:CD1	2.88	0.41
1:J:311:LEU:HD13	3:L:57:THR:HG21	2.02	0.41
1:J:227:ARG:HB2	1:J:273:GLU:CD	2.40	0.41
1:J:84:ILE:HD11	1:J:131:SER:CB	2.50	0.41
1:G:128:LEU:O	1:G:131:SER:N	2.54	0.41
1:A:57:LYS:HD3	1:A:217:ARG:HH21	1.85	0.41
1:D:11:MET:CE	1:D:140:HIS:CD2	3.04	0.41
1:G:217:ARG:HA	2:H:71:LEU:CD1	2.51	0.41
1:A:208:ILE:HG21	2:B:9:THR:HG22	2.01	0.41
3:C:76:GLN:OE1	3:F:72:GLN:HG3	2.21	0.41
1:G:40:PRO:CB	1:G:201:ARG:NH2	2.76	0.41
1:D:55:LEU:HA	1:D:164:HIS:O	2.21	0.41
1:D:24:GLY:HA3	1:D:139:VAL:CG1	2.50	0.41
1:D:57:LYS:HD3	1:D:217:ARG:HH21	1.86	0.41
1:A:8:TRP:CE2	1:A:84:ILE:HD13	2.55	0.41
1:J:11:MET:CE	1:J:140:HIS:CD2	3.03	0.41
1:J:8:TRP:HE3	1:J:140:HIS:ND1	2.18	0.41
2:B:37:PRO:HD2	2:B:40:GLN:OE1	2.21	0.41
2:B:42:ARG:CZ	2:B:72:ARG:HG3	2.51	0.41
1:J:134:ASP:O	1:J:138:GLN:HG3	2.21	0.41
1:D:40:PRO:CD	1:D:201:ARG:HH12	2.31	0.41
1:G:8:TRP:CE2	1:G:84:ILE:HD13	2.56	0.41
1:A:301:THR:HG21	3:C:65:SER:CB	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:208:ILE:CG2	2:E:9:THR:HG22	2.51	0.41
1:A:55:LEU:HA	1:A:164:HIS:O	2.21	0.41
1:G:217:ARG:HA	2:H:71:LEU:HD12	2.02	0.41
1:G:289:HIS:CG	3:I:75:PRO:HB3	2.56	0.41
2:K:42:ARG:CZ	2:K:72:ARG:HG3	2.50	0.41
1:A:18:PHE:HB2	1:A:32:VAL:HG21	2.04	0.40
1:A:294:PHE:CE1	3:C:69:HIS:HB3	2.56	0.40
1:G:134:ASP:O	1:G:138:GLN:HG3	2.22	0.40
1:A:219:ASN:OD1	1:A:219:ASN:C	2.60	0.40
1:G:37:SER:O	1:G:42:ASN:OD1	2.40	0.40
1:J:57:LYS:HD3	1:J:217:ARG:HH21	1.85	0.40
1:G:63:GLU:HG2	1:G:64:PRO:CD	2.51	0.40
1:G:140:HIS:CD2	1:G:140:HIS:C	2.94	0.40
1:J:37:SER:O	1:J:42:ASN:OD1	2.40	0.40
1:A:43:PHE:CD2	1:A:197:ILE:HG21	2.56	0.40
1:J:8:TRP:CE2	1:J:84:ILE:HD13	2.56	0.40
3:C:62:LEU:HB3	3:C:66:GLN:NE2	2.36	0.40
2:E:7:THR:OG1	2:E:9:THR:OG1	2.29	0.40
1:A:13:SER:OG	1:A:219:ASN:ND2	2.54	0.40
1:D:134:ASP:O	1:D:138:GLN:HG3	2.21	0.40
1:D:128:LEU:O	1:D:131:SER:N	2.55	0.40
2:K:37:PRO:HD2	2:K:40:GLN:OE1	2.22	0.40
1:G:55:LEU:HA	1:G:164:HIS:O	2.21	0.40

All (4) 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:244:GLU:O	1:J:174:ARG:NH1[2_755]	1.99	0.21
2:B:32:ASP:O	2:K:60:ASN:ND2[4_655]	2.05	0.15
2:E:63:LYS:NZ	3:I:82:ALA:O[4_656]	2.14	0.06
2:B:63:LYS:NZ	3:L:79:GLU:OE1[4_655]	2.18	0.02

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	286/331 (86%)	273 (96%)	13 (4%)	0	100	100
1	D	289/331 (87%)	276 (96%)	13 (4%)	0	100	100
1	G	281/331 (85%)	269 (96%)	12 (4%)	0	100	100
1	J	288/331 (87%)	276 (96%)	12 (4%)	0	100	100
2	B	73/76 (96%)	71 (97%)	2 (3%)	0	100	100
2	E	73/76 (96%)	70 (96%)	3 (4%)	0	100	100
2	H	73/76 (96%)	71 (97%)	2 (3%)	0	100	100
2	K	73/76 (96%)	71 (97%)	2 (3%)	0	100	100
3	C	48/65 (74%)	41 (85%)	7 (15%)	0	100	100
3	F	48/65 (74%)	41 (85%)	7 (15%)	0	100	100
3	I	47/65 (72%)	39 (83%)	8 (17%)	0	100	100
3	L	47/65 (72%)	39 (83%)	8 (17%)	0	100	100
All	All	1626/1888 (86%)	1537 (94%)	89 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	257/288 (89%)	257 (100%)	0	100	100
1	D	259/288 (90%)	257 (99%)	2 (1%)	86	94
1	G	252/288 (88%)	252 (100%)	0	100	100
1	J	259/288 (90%)	258 (100%)	1 (0%)	93	98
2	B	68/68 (100%)	67 (98%)	1 (2%)	72	90
2	E	68/68 (100%)	68 (100%)	0	100	100
2	H	68/68 (100%)	68 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	K	68/68 (100%)	67 (98%)	1 (2%)	72	90
3	C	48/59 (81%)	48 (100%)	0	100	100
3	F	48/59 (81%)	48 (100%)	0	100	100
3	I	47/59 (80%)	46 (98%)	1 (2%)	61	86
3	L	47/59 (80%)	46 (98%)	1 (2%)	61	86
All	All	1489/1660 (90%)	1482 (100%)	7 (0%)	92	97

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	48	LYS
1	D	59	GLN
1	D	148	MET
3	I	46	PRO
1	J	145	ARG
2	K	48	LYS
3	L	46	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	HIS
1	D	42	ASN
1	D	233	GLN

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, 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	292/331 (88%)	-0.06	3 (1%) 84 72	53, 79, 143, 181	0
1	D	295/331 (89%)	-0.10	0 100 100	48, 66, 125, 169	0
1	G	287/331 (86%)	-0.09	3 (1%) 84 72	50, 77, 146, 185	0
1	J	294/331 (88%)	-0.14	0 100 100	42, 63, 138, 170	0
2	B	75/76 (98%)	0.30	3 (4%) 42 28	107, 140, 187, 199	0
2	E	75/76 (98%)	-0.01	0 100 100	69, 86, 103, 106	0
2	H	75/76 (98%)	0.10	0 100 100	76, 90, 100, 102	0
2	K	75/76 (98%)	0.06	0 100 100	58, 76, 92, 97	0
3	C	50/65 (76%)	0.11	1 (2%) 68 53	102, 114, 126, 132	0
3	F	50/65 (76%)	0.23	2 (4%) 42 28	99, 113, 127, 133	0
3	I	49/65 (75%)	0.26	1 (2%) 68 53	106, 124, 144, 154	0
3	L	49/65 (75%)	0.36	2 (4%) 41 27	91, 115, 149, 158	0
All	All	1666/1888 (88%)	-0.02	15 (0%) 85 75	42, 80, 146, 199	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	71	GLN	4.0
3	F	81	SER	3.9
1	G	243	ALA	3.8
3	C	71	GLN	3.0
3	L	71	GLN	2.7
3	I	48	ILE	2.6
2	B	4	PHE	2.5
1	G	160	GLU	2.5
1	A	118	SER	2.4
1	G	312	VAL	2.3
2	B	5	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	3	ILE	2.2
1	A	306	GLN	2.0
1	A	115	LYS	2.0
3	L	83	GLU	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 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.