



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 20, 2025 – 10:39 AM JST

PDB ID : 9K7Q / pdb\_00009k7q  
Title : Parkinson disease protein 7 (DJ-1) and Alpha-synuclein (Alpha-syn) complex  
Authors : Han, C.W.; Kim, D.H.  
Deposited on : 2024-10-24  
Resolution : 4.58 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.46

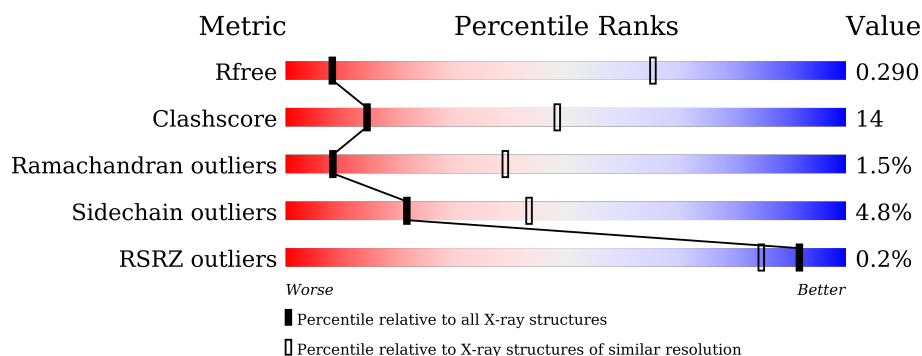
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 4.58 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))
$R_{free}$	164625	1065 (5.26-3.90)
Clashscore	180529	1120 (5.26-3.90)
Ramachandran outliers	177936	1014 (5.26-3.90)
Sidechain outliers	177891	1014 (5.26-3.88)
RSRZ outliers	164620	1061 (5.26-3.90)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	187	<div> <div>84%</div> <div>14%</div> <div>.</div> </div>
1	D	187	<div> <div>86%</div> <div>13%</div> <div>.</div> </div>
1	E	187	<div> <div>80%</div> <div>17%</div> <div>.</div> </div>
1	F	187	<div> <div>%</div> <div>80%</div> <div>18%</div> <div>.</div> </div>
1	H	187	<div> <div>%</div> <div>82%</div> <div>17%</div> <div>.</div> </div>
1	J	187	<div> <div>84%</div> <div>16%</div> </div>

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Mol	Chain	Length	Quality of chain
2	M	92	<div><div><div></div><div></div><div></div><div></div></div><div>%27%42%25%5%</div></div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8885 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 Parkinson disease protein 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	187	Total	C	N	O	S	0	0	0
			1375	865	240	263	7			
1	D	187	Total	C	N	O	S	0	0	0
			1375	865	240	263	7			
1	E	187	Total	C	N	O	S	0	0	0
			1375	865	240	263	7			
1	F	187	Total	C	N	O	S	0	0	0
			1375	865	240	263	7			
1	H	187	Total	C	N	O	S	0	0	0
			1375	865	240	263	7			
1	J	187	Total	C	N	O	S	0	0	0
			1375	865	240	263	7			


- Molecule 2 is a protein called Alpha-synuclein.

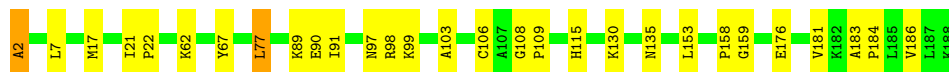
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	92	Total	C	N	O	S	0	0	0
			635	395	110	128	2			

### 3 Residue-property plots


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: Parkinson disease protein 7

Chain A: 




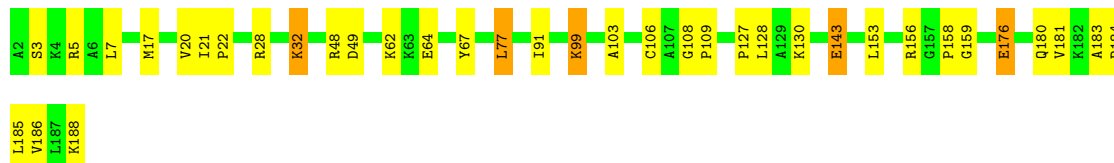
- Molecule 1: Parkinson disease protein 7

Chain D: 




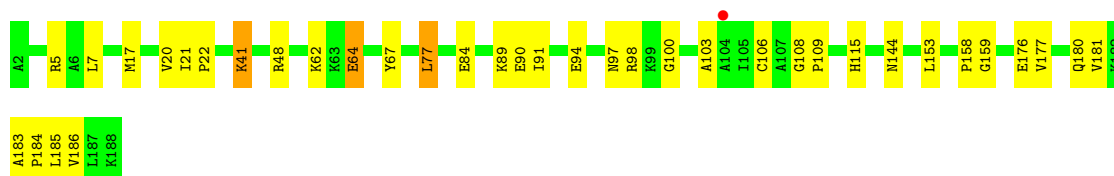
- Molecule 1: Parkinson disease protein 7

Chain E: 




- Molecule 1: Parkinson disease protein 7

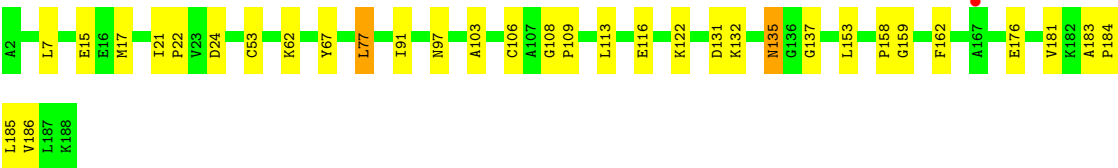
Chain F: 



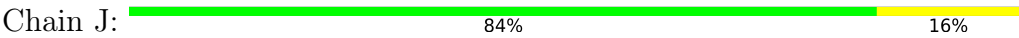
- Molecule 1: Parkinson disease protein 7

Chain H: 

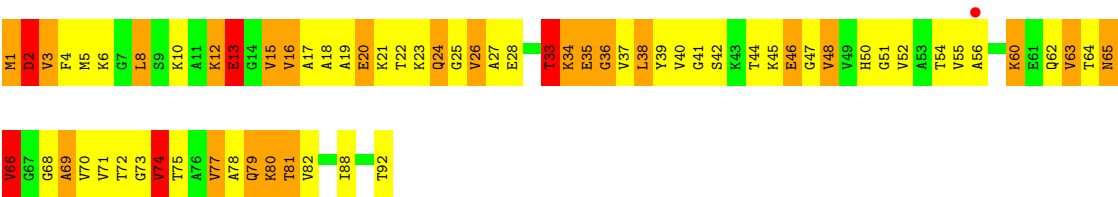




● Molecule 1: Parkinson disease protein 7



● Molecule 2: Alpha-synuclein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	75.28Å 75.27Å 75.39Å 89.89° 89.92° 60.04°	Depositor
Resolution (Å)	7.66 – 4.58 7.66 – 4.58	Depositor EDS
% Data completeness (in resolution range)	75.4 (7.66-4.58) 78.0 (7.66-4.58)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.58 (at 4.61Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.145 , 0.222 0.193 , 0.290	Depositor DCC
$R_{free}$ test set	336 reflections (4.16%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	68.4	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.23 , 44.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.55$ , $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	0.000 for h-k,h,l 0.000 for k,-h+k,l 0.250 for -k,h-k,l 0.250 for -h+k,-h,l 0.268 for -h+k,k,-l 0.309 for h,h-k,-l 0.000 for -h,-k,l 0.000 for k,h,-l 0.277 for -k,-h,-l 0.000 for h-k,-k,-l 0.000 for -h,-h+k,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	8885	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.93% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  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	0.59	0/1392	1.12	2/1880 (0.1%)
1	D	0.58	0/1392	1.10	2/1880 (0.1%)
1	E	0.56	0/1392	1.10	5/1880 (0.3%)
1	F	0.56	0/1392	1.12	2/1880 (0.1%)
1	H	0.57	0/1392	1.12	2/1880 (0.1%)
1	J	0.56	0/1392	1.12	3/1880 (0.2%)
2	M	0.62	0/637	1.59	8/856 (0.9%)
All	All	0.57	0/8989	1.15	24/12136 (0.2%)

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	1
2	M	0	6
All	All	0	7

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	176	GLU	CB-CG-CD	9.77	129.21	112.60
2	M	64	THR	N-CA-C	-7.50	104.65	113.88
2	M	2	ASP	CB-CA-C	-7.03	98.94	109.90
1	J	76	ASN	CB-CA-C	6.85	121.63	110.88
1	A	176	GLU	CB-CG-CD	6.71	124.01	112.60
1	E	143	GLU	CB-CG-CD	6.69	123.97	112.60
1	E	99	LYS	CB-CG-CD	6.47	126.19	111.30
2	M	24	GLN	CB-CA-C	-6.47	100.64	110.92
1	D	64	GLU	CB-CG-CD	-6.26	101.96	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	79	GLN	N-CA-C	-6.19	104.50	114.09
1	H	24	ASP	CA-CB-CG	6.13	118.73	112.60
1	F	176	GLU	CB-CG-CD	6.12	123.01	112.60
1	A	90	GLU	CB-CG-CD	-6.05	102.31	112.60
1	D	170	GLU	CB-CG-CD	5.83	122.52	112.60
2	M	70	VAL	N-CA-C	-5.76	107.51	113.10
1	J	94	GLU	CB-CG-CD	5.68	122.26	112.60
2	M	33	THR	CA-CB-OG1	5.67	118.10	109.60
1	E	176	GLU	CB-CG-CD	5.63	122.17	112.60
1	F	144	ASN	CA-CB-CG	-5.52	107.08	112.60
2	M	4	PHE	CA-CB-CG	-5.33	108.47	113.80
1	E	130	LYS	CA-C-N	5.29	127.68	120.54
1	E	130	LYS	C-N-CA	5.29	127.68	120.54
1	J	144	ASN	CA-CB-CG	-5.23	107.37	112.60
2	M	39	TYR	CB-CA-C	5.16	119.16	109.71

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	2	ALA	Peptide
2	M	13	GLU	Peptide
2	M	54	THR	Peptide
2	M	65	ASN	Peptide
2	M	66	VAL	Peptide
2	M	81	THR	Peptide
2	M	88	ILE	Peptide

## 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1375	0	1431	25	1
1	D	1375	0	1431	20	0
1	E	1375	0	1431	47	0
1	F	1375	0	1431	42	0
1	H	1375	0	1431	52	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	J	1375	0	1431	23	1
2	M	635	0	673	122	2
All	All	8885	0	9259	250	2

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 14.

All (250) 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:A:7:LEU:HB2	1:A:67:TYR:CD2	1.67	1.30
1:D:7:LEU:HB2	1:D:67:TYR:CD2	1.67	1.29
1:H:7:LEU:HB2	1:H:67:TYR:CD2	1.68	1.28
1:F:7:LEU:HB2	1:F:67:TYR:CD2	1.68	1.27
1:J:7:LEU:HB2	1:J:67:TYR:CD2	1.67	1.27
1:E:7:LEU:HB2	1:E:67:TYR:CD2	1.68	1.26
1:F:180:GLN:NE2	2:M:28:GLU:HB3	1.60	1.15
2:M:56:ALA:O	2:M:60:LYS:HB2	1.50	1.10
1:F:180:GLN:HE22	2:M:28:GLU:CB	1.66	1.07
1:D:7:LEU:HB2	1:D:67:TYR:HD2	1.01	1.03
1:J:7:LEU:HB2	1:J:67:TYR:HD2	1.01	1.00
1:E:7:LEU:HB2	1:E:67:TYR:HD2	1.03	0.99
1:F:180:GLN:NE2	2:M:28:GLU:CB	2.26	0.97
1:F:180:GLN:CD	2:M:28:GLU:HB3	1.88	0.97
1:E:128:LEU:HA	2:M:38:LEU:HD21	1.46	0.96
1:A:7:LEU:HB2	1:A:67:TYR:HD2	1.03	0.96
2:M:46:GLU:O	2:M:50:HIS:HB3	1.66	0.95
1:F:7:LEU:HB2	1:F:67:TYR:HD2	1.02	0.94
2:M:1:MET:H1	2:M:6:LYS:HG2	1.30	0.94
2:M:13:GLU:O	2:M:17:ALA:HB3	1.68	0.93
1:F:180:GLN:HE22	2:M:28:GLU:CD	1.78	0.92
1:H:7:LEU:HB2	1:H:67:TYR:HD2	1.05	0.91
2:M:1:MET:N	2:M:6:LYS:HG2	1.84	0.91
1:H:77:LEU:N	1:H:77:LEU:HD23	1.86	0.90
1:H:77:LEU:HD23	1:H:77:LEU:H	1.36	0.90
1:J:7:LEU:CB	1:J:67:TYR:HD2	1.84	0.89
1:H:113:LEU:HD11	2:M:34:LYS:NZ	1.88	0.89
1:E:156:ARG:HH22	2:M:40:VAL:HA	1.35	0.88
1:F:7:LEU:CB	1:F:67:TYR:HD2	1.86	0.87
1:E:7:LEU:CB	1:E:67:TYR:HD2	1.87	0.86
1:D:7:LEU:CB	1:D:67:TYR:HD2	1.85	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:41:LYS:HD2	1:F:41:LYS:O	1.77	0.84
1:A:7:LEU:CB	1:A:67:TYR:HD2	1.86	0.84
1:H:7:LEU:CB	1:H:67:TYR:HD2	1.88	0.84
1:H:135:ASN:HB2	2:M:28:GLU:HG2	1.61	0.83
1:H:113:LEU:CD1	2:M:34:LYS:NZ	2.43	0.82
1:E:127:PRO:CG	2:M:40:VAL:HG22	2.11	0.81
2:M:47:GLY:HA2	2:M:51:GLY:H	1.46	0.81
1:H:131:ASP:OD2	2:M:23:LYS:HG2	1.81	0.80
2:M:63:VAL:HG13	2:M:66:VAL:O	1.82	0.79
2:M:78:ALA:CA	2:M:81:THR:HG22	2.15	0.75
2:M:36:GLY:C	2:M:38:LEU:H	1.95	0.75
1:E:5:ARG:NH1	1:E:64:GLU:O	2.20	0.75
2:M:5:MET:CE	2:M:8:LEU:HD23	2.16	0.75
2:M:63:VAL:CG1	2:M:66:VAL:O	2.36	0.74
2:M:44:THR:O	2:M:46:GLU:N	2.16	0.74
1:H:7:LEU:HB2	1:H:67:TYR:CE2	2.24	0.73
1:F:180:GLN:HE22	2:M:28:GLU:CG	2.02	0.73
1:A:98:ARG:HG3	1:A:98:ARG:HH11	1.54	0.73
1:A:7:LEU:HB2	1:A:67:TYR:CE2	2.23	0.72
1:F:7:LEU:HB2	1:F:67:TYR:CE2	2.24	0.72
1:H:135:ASN:HB2	2:M:28:GLU:HA	1.70	0.72
1:E:127:PRO:C	2:M:38:LEU:HD23	2.14	0.72
2:M:78:ALA:HB1	2:M:81:THR:CG2	2.20	0.72
1:A:130:LYS:NZ	1:E:3:SER:OG	2.23	0.71
1:D:7:LEU:HB2	1:D:67:TYR:CE2	2.23	0.71
1:F:180:GLN:CD	2:M:28:GLU:CB	2.62	0.71
1:H:113:LEU:HD11	2:M:34:LYS:HZ3	1.53	0.70
1:H:113:LEU:CD1	2:M:34:LYS:HZ1	2.04	0.70
2:M:1:MET:HG3	2:M:6:LYS:HZ3	1.57	0.70
2:M:13:GLU:O	2:M:17:ALA:CB	2.39	0.70
1:J:7:LEU:HB2	1:J:67:TYR:CE2	2.24	0.70
2:M:78:ALA:HA	2:M:81:THR:HG22	1.74	0.70
1:A:159:GLY:N	1:D:185:LEU:O	2.24	0.70
1:J:5:ARG:NH1	1:J:64:GLU:O	2.24	0.70
1:E:7:LEU:HB2	1:E:67:TYR:CE2	2.25	0.69
1:F:7:LEU:HD13	1:F:67:TYR:HE2	1.57	0.69
2:M:78:ALA:HB1	2:M:81:THR:HG22	1.75	0.69
2:M:5:MET:HA	2:M:8:LEU:HB2	1.74	0.68
2:M:56:ALA:O	2:M:60:LYS:CB	2.37	0.68
1:F:180:GLN:OE1	2:M:28:GLU:C	2.36	0.68
1:H:7:LEU:HD13	1:H:67:TYR:HE2	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:7:LEU:HD13	1:J:67:TYR:HE2	1.58	0.68
1:A:7:LEU:HD13	1:A:67:TYR:HE2	1.56	0.68
1:F:180:GLN:NE2	2:M:28:GLU:CD	2.49	0.67
2:M:78:ALA:CB	2:M:81:THR:HG22	2.25	0.67
1:D:7:LEU:HD13	1:D:67:TYR:HE2	1.58	0.67
1:F:180:GLN:OE1	2:M:28:GLU:CB	2.43	0.67
1:H:135:ASN:HB2	2:M:28:GLU:CG	2.25	0.67
1:H:113:LEU:HD11	2:M:34:LYS:HZ1	1.58	0.66
2:M:72:THR:HA	2:M:75:THR:OG1	1.96	0.66
1:E:7:LEU:HD13	1:E:67:TYR:HE2	1.59	0.66
1:E:127:PRO:C	2:M:38:LEU:CD2	2.70	0.65
2:M:48:VAL:HA	2:M:52:VAL:HG23	1.78	0.65
1:E:128:LEU:CA	2:M:38:LEU:HD21	2.25	0.65
2:M:81:THR:HG23	2:M:81:THR:O	1.97	0.65
2:M:62:GLN:O	2:M:66:VAL:HG23	1.98	0.64
1:E:62:LYS:HE2	1:E:91:ILE:HD11	1.79	0.64
1:F:62:LYS:HE2	1:F:91:ILE:HD11	1.81	0.63
1:E:127:PRO:O	2:M:38:LEU:CD2	2.47	0.63
1:E:127:PRO:CB	2:M:40:VAL:HG22	2.29	0.62
1:H:62:LYS:HE2	1:H:91:ILE:HD11	1.81	0.62
1:E:128:LEU:HA	2:M:38:LEU:CD2	2.26	0.62
1:D:62:LYS:HE2	1:D:91:ILE:HD11	1.80	0.62
1:J:62:LYS:HE2	1:J:91:ILE:HD11	1.81	0.61
1:H:185:LEU:O	1:J:159:GLY:N	2.33	0.61
1:A:62:LYS:HE2	1:A:91:ILE:HD11	1.81	0.61
1:H:77:LEU:N	1:H:77:LEU:CD2	2.57	0.61
2:M:47:GLY:HA2	2:M:51:GLY:N	2.15	0.60
1:H:132:LYS:N	2:M:27:ALA:HB1	2.15	0.60
1:H:135:ASN:CB	2:M:28:GLU:HA	2.31	0.60
1:A:98:ARG:HG3	1:A:98:ARG:NH1	2.16	0.60
2:M:33:THR:O	2:M:36:GLY:N	2.35	0.59
1:H:135:ASN:CG	2:M:28:GLU:HA	2.27	0.59
1:E:127:PRO:HB3	2:M:40:VAL:HG22	1.84	0.59
1:E:127:PRO:HG3	2:M:40:VAL:HG22	1.85	0.58
1:F:17:MET:HE3	1:F:158:PRO:HB3	1.86	0.57
2:M:36:GLY:O	2:M:38:LEU:N	2.35	0.57
1:E:49:ASP:OD2	1:H:122:LYS:NZ	2.36	0.57
1:H:135:ASN:HB2	2:M:28:GLU:CA	2.36	0.56
1:E:49:ASP:CG	1:H:122:LYS:NZ	2.64	0.56
2:M:66:VAL:HG13	2:M:69:ALA:HB3	1.88	0.56
1:F:90:GLU:HG2	1:F:94:GLU:OE2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:180:GLN:OE1	2:M:28:GLU:HB3	2.01	0.56
1:A:135:ASN:OD1	1:E:180:GLN:CD	2.50	0.55
1:A:7:LEU:CB	1:A:67:TYR:CD2	2.62	0.55
2:M:17:ALA:C	2:M:19:ALA:H	2.14	0.55
2:M:35:GLU:O	2:M:38:LEU:HB2	2.07	0.55
1:H:116:GLU:OE1	2:M:34:LYS:HA	2.08	0.54
1:D:17:MET:HE3	1:D:158:PRO:HB3	1.90	0.53
1:H:135:ASN:CB	2:M:28:GLU:HG2	2.35	0.53
2:M:22:THR:O	2:M:26:VAL:HG22	2.09	0.53
1:F:41:LYS:O	1:F:41:LYS:CD	2.53	0.53
2:M:12:LYS:HG3	2:M:12:LYS:O	2.09	0.53
1:H:53:CYS:SG	1:J:53:CYS:SG	3.07	0.52
2:M:1:MET:H1	2:M:6:LYS:CG	2.12	0.52
1:H:17:MET:HE3	1:H:158:PRO:HB3	1.93	0.51
1:J:7:LEU:CB	1:J:67:TYR:CD2	2.62	0.51
1:E:7:LEU:CB	1:E:67:TYR:CD2	2.63	0.51
1:D:7:LEU:CB	1:D:67:TYR:CD2	2.62	0.51
1:E:156:ARG:NH2	2:M:40:VAL:HG13	2.26	0.51
1:E:185:LEU:O	1:F:159:GLY:N	2.44	0.51
2:M:17:ALA:O	2:M:20:GLU:HG3	2.10	0.51
1:F:5:ARG:NH1	1:F:64:GLU:O	2.36	0.51
1:E:127:PRO:O	2:M:38:LEU:HD21	2.10	0.50
1:H:21:ILE:HB	1:H:22:PRO:HD3	1.94	0.50
2:M:24:GLN:HB3	2:M:28:GLU:OE1	2.12	0.50
1:A:21:ILE:HB	1:A:22:PRO:HD3	1.94	0.50
1:H:113:LEU:HD12	2:M:34:LYS:NZ	2.24	0.50
1:H:159:GLY:HA3	1:J:185:LEU:O	2.11	0.50
1:D:181:VAL:O	1:D:184:PRO:HD2	2.12	0.50
1:F:21:ILE:HB	1:F:22:PRO:HD3	1.94	0.50
1:F:181:VAL:O	1:F:184:PRO:HD2	2.12	0.50
1:H:132:LYS:CA	2:M:27:ALA:HB1	2.42	0.49
2:M:68:GLY:O	2:M:69:ALA:CB	2.59	0.49
1:E:17:MET:HE3	1:E:158:PRO:HB3	1.94	0.49
2:M:13:GLU:HA	2:M:16:VAL:H	1.78	0.49
1:F:177:VAL:HG22	2:M:28:GLU:OE1	2.12	0.49
1:J:21:ILE:HB	1:J:22:PRO:HD3	1.95	0.49
2:M:68:GLY:O	2:M:69:ALA:HB2	2.11	0.49
1:D:21:ILE:HB	1:D:22:PRO:HD3	1.95	0.49
1:H:181:VAL:O	1:H:184:PRO:HD2	2.13	0.49
1:E:21:ILE:HB	1:E:22:PRO:HD3	1.95	0.48
1:E:181:VAL:O	1:E:184:PRO:HD2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:5:MET:HE1	2:M:8:LEU:HD23	1.95	0.48
1:A:181:VAL:O	1:A:184:PRO:HD2	2.13	0.48
1:A:17:MET:HE3	1:A:158:PRO:HB3	1.95	0.48
2:M:2:ASP:O	2:M:3:VAL:C	2.57	0.47
2:M:5:MET:HE2	2:M:8:LEU:HD23	1.94	0.47
1:H:135:ASN:OD1	2:M:28:GLU:O	2.33	0.47
2:M:63:VAL:HG12	2:M:66:VAL:O	2.11	0.47
2:M:71:VAL:HA	2:M:74:VAL:HG23	1.96	0.47
2:M:36:GLY:C	2:M:38:LEU:N	2.65	0.47
1:H:131:ASP:OD2	2:M:23:LYS:CG	2.58	0.47
2:M:66:VAL:CG1	2:M:69:ALA:HB3	2.45	0.47
1:F:77:LEU:H	1:F:77:LEU:HG	1.56	0.47
1:H:185:LEU:O	1:J:159:GLY:HA3	2.15	0.47
1:J:181:VAL:O	1:J:184:PRO:HD2	2.14	0.47
1:H:135:ASN:HB2	2:M:28:GLU:CB	2.46	0.46
2:M:19:ALA:C	2:M:21:LYS:H	2.24	0.46
2:M:79:GLN:HA	2:M:82:VAL:HG22	1.96	0.46
1:F:98:ARG:NH1	1:F:100:GLY:HA3	2.31	0.46
2:M:21:LYS:O	2:M:24:GLN:OE1	2.33	0.46
1:A:159:GLY:HA3	1:D:185:LEU:O	2.16	0.46
1:E:143:GLU:HB3	2:M:41:GLY:CA	2.46	0.46
1:F:180:GLN:OE1	2:M:28:GLU:O	2.34	0.45
1:A:99:LYS:NZ	1:F:84:GLU:OE2	2.33	0.45
2:M:13:GLU:C	2:M:15:VAL:H	2.23	0.45
1:A:103:ALA:HA	1:A:153:LEU:O	2.17	0.45
2:M:40:VAL:HG12	2:M:42:SER:OG	2.17	0.45
2:M:47:GLY:HA2	2:M:51:GLY:HA3	1.98	0.45
1:A:159:GLY:CA	1:D:185:LEU:O	2.64	0.45
2:M:35:GLU:O	2:M:38:LEU:CB	2.65	0.45
1:A:77:LEU:H	1:A:77:LEU:HG	1.27	0.45
1:E:28:ARG:HG2	1:F:48:ARG:HH21	1.82	0.45
1:J:98:ARG:NH1	1:J:100:GLY:HA3	2.32	0.45
2:M:78:ALA:CB	2:M:81:THR:CG2	2.89	0.45
1:E:159:GLY:N	1:F:185:LEU:O	2.48	0.45
1:E:28:ARG:CZ	1:F:48:ARG:NH2	2.80	0.45
2:M:1:MET:H2	2:M:6:LYS:HG2	1.74	0.45
1:E:103:ALA:HA	1:E:153:LEU:O	2.17	0.44
1:E:159:GLY:HA3	1:F:185:LEU:O	2.17	0.44
1:H:162:PHE:CE1	1:J:17:MET:SD	3.11	0.44
1:J:103:ALA:HA	1:J:153:LEU:O	2.18	0.44
1:H:7:LEU:CB	1:H:67:TYR:CD2	2.63	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:77:LEU:H	1:D:77:LEU:HG	1.35	0.44
1:D:103:ALA:HA	1:D:153:LEU:O	2.18	0.44
2:M:47:GLY:HA2	2:M:51:GLY:CA	2.48	0.44
1:D:90:GLU:HG2	1:D:94:GLU:OE2	2.17	0.44
1:H:103:ALA:HA	1:H:153:LEU:O	2.17	0.43
1:H:185:LEU:O	1:J:159:GLY:CA	2.66	0.43
1:E:143:GLU:HB3	2:M:41:GLY:HA3	1.99	0.43
2:M:33:THR:O	2:M:34:LYS:C	2.61	0.43
2:M:46:GLU:O	2:M:50:HIS:CB	2.53	0.43
1:A:183:ALA:HB3	1:A:184:PRO:HD3	2.00	0.43
1:F:103:ALA:HA	1:F:153:LEU:O	2.18	0.43
1:J:17:MET:HE3	1:J:158:PRO:HB3	2.00	0.43
1:D:108:GLY:N	1:D:109:PRO:CD	2.81	0.43
1:E:20:VAL:HG21	1:F:20:VAL:HG21	1.99	0.43
2:M:77:VAL:O	2:M:80:LYS:HB2	2.19	0.43
2:M:23:LYS:O	2:M:27:ALA:CB	2.67	0.42
1:F:7:LEU:CB	1:F:67:TYR:CD2	2.63	0.42
2:M:44:THR:C	2:M:46:GLU:H	2.18	0.42
1:D:64:GLU:OE1	1:D:64:GLU:HA	2.16	0.42
1:E:49:ASP:OD1	1:H:122:LYS:NZ	2.52	0.42
1:H:183:ALA:HB3	1:H:184:PRO:HD3	2.00	0.42
1:F:89:LYS:HG3	1:F:115:HIS:HB3	2.02	0.42
1:H:17:MET:SD	1:J:21:ILE:HA	2.59	0.42
1:A:108:GLY:N	1:A:109:PRO:CD	2.82	0.42
1:H:135:ASN:CA	2:M:28:GLU:HG2	2.50	0.42
2:M:10:LYS:NZ	2:M:13:GLU:OE2	2.45	0.42
2:M:17:ALA:C	2:M:19:ALA:N	2.77	0.42
1:E:108:GLY:N	1:E:109:PRO:CD	2.83	0.42
1:E:127:PRO:C	2:M:38:LEU:HD21	2.43	0.42
1:E:183:ALA:HB3	1:E:184:PRO:HD3	2.02	0.42
1:E:32:LYS:HB3	1:E:32:LYS:HE2	1.48	0.42
1:A:89:LYS:HG3	1:A:115:HIS:HB3	2.02	0.41
1:F:108:GLY:N	1:F:109:PRO:CD	2.82	0.41
1:H:108:GLY:N	1:H:109:PRO:CD	2.82	0.41
1:H:113:LEU:HD12	2:M:34:LYS:HZ1	1.79	0.41
1:H:135:ASN:HA	2:M:28:GLU:HG2	2.02	0.41
2:M:17:ALA:O	2:M:19:ALA:N	2.52	0.41
1:J:98:ARG:NH1	1:J:100:GLY:CA	2.84	0.41
1:D:183:ALA:HB3	1:D:184:PRO:HD3	2.02	0.41
1:H:132:LYS:HA	2:M:27:ALA:HB1	2.01	0.41
1:F:180:GLN:OE1	2:M:28:GLU:HB2	2.17	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:73:GLY:O	2:M:77:VAL:HG22	2.20	0.41
1:F:183:ALA:HB3	1:F:184:PRO:HD3	2.02	0.41
1:D:89:LYS:HG3	1:D:115:HIS:HB3	2.02	0.41
2:M:23:LYS:O	2:M:27:ALA:HB2	2.21	0.41
2:M:56:ALA:O	2:M:60:LYS:N	2.44	0.41
1:J:108:GLY:N	1:J:109:PRO:CD	2.84	0.41
1:E:77:LEU:HD22	1:E:77:LEU:HA	1.89	0.41
1:E:143:GLU:O	2:M:41:GLY:HA3	2.21	0.41
1:A:135:ASN:OD1	1:E:180:GLN:NE2	2.54	0.41
2:M:20:GLU:HG3	2:M:20:GLU:H	1.47	0.40
1:A:7:LEU:HD13	1:A:67:TYR:CE2	2.47	0.40
1:E:48:ARG:NH2	1:H:137:GLY:O	2.55	0.40
1:H:15:GLU:OE1	1:J:24:ASP:OD2	2.39	0.40
1:E:185:LEU:O	1:F:159:GLY:HA3	2.22	0.40

All (2) 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:2:ALA:N	2:M:36:GLY:O[1_545]	1.95	0.25
1:J:39:ALA:O	2:M:71:VAL:CG1[1_655]	2.18	0.02

## 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	185/187 (99%)	180 (97%)	4 (2%)	1 (0%)	25 64
1	D	185/187 (99%)	181 (98%)	3 (2%)	1 (0%)	25 64
1	E	185/187 (99%)	182 (98%)	2 (1%)	1 (0%)	25 64
1	F	185/187 (99%)	182 (98%)	2 (1%)	1 (0%)	25 64

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	185/187 (99%)	181 (98%)	3 (2%)	1 (0%)	25	64
1	J	185/187 (99%)	182 (98%)	2 (1%)	1 (0%)	25	64
2	M	90/92 (98%)	55 (61%)	23 (26%)	12 (13%)	0	4
All	All	1200/1214 (99%)	1143 (95%)	39 (3%)	18 (2%)	8	39

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	M	37	VAL
2	M	45	LYS
2	M	48	VAL
2	M	69	ALA
2	M	36	GLY
2	M	18	ALA
1	A	106	CYS
1	D	106	CYS
1	E	106	CYS
1	F	106	CYS
1	H	106	CYS
1	J	106	CYS
2	M	3	VAL
2	M	33	THR
2	M	34	LYS
2	M	74	VAL
2	M	25	GLY
2	M	66	VAL

### 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	145/145 (100%)	142 (98%)	3 (2%)	48	67
1	D	145/145 (100%)	141 (97%)	4 (3%)	38	58
1	E	145/145 (100%)	139 (96%)	6 (4%)	26	48

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	145/145 (100%)	140 (97%)	5 (3%)	32	53
1	H	145/145 (100%)	141 (97%)	4 (3%)	38	58
1	J	145/145 (100%)	143 (99%)	2 (1%)	62	75
2	M	63/63 (100%)	42 (67%)	21 (33%)	0	2
All	All	933/933 (100%)	888 (95%)	45 (5%)	21	43

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	LEU
1	A	97	ASN
1	A	186	VAL
1	D	41	LYS
1	D	77	LEU
1	D	175	LYS
1	D	186	VAL
1	E	32	LYS
1	E	77	LEU
1	E	99	LYS
1	E	176	GLU
1	E	186	VAL
1	E	188	LYS
1	F	41	LYS
1	F	64	GLU
1	F	77	LEU
1	F	97	ASN
1	F	186	VAL
1	H	77	LEU
1	H	97	ASN
1	H	135	ASN
1	H	186	VAL
1	J	97	ASN
1	J	186	VAL
2	M	1	MET
2	M	2	ASP
2	M	8	LEU
2	M	12	LYS
2	M	13	GLU
2	M	15	VAL
2	M	16	VAL
2	M	20	GLU

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Mol	Chain	Res	Type
2	M	26	VAL
2	M	35	GLU
2	M	38	LEU
2	M	46	GLU
2	M	55	VAL
2	M	60	LYS
2	M	63	VAL
2	M	65	ASN
2	M	66	VAL
2	M	74	VAL
2	M	77	VAL
2	M	80	LYS
2	M	92	THR

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

Mol	Chain	Res	Type
1	A	138	HIS
1	A	180	GLN
1	D	135	ASN
1	D	138	HIS
1	D	144	ASN
1	D	180	GLN
1	E	138	HIS
1	E	180	GLN
1	F	135	ASN
1	F	138	HIS
1	F	144	ASN
1	F	180	GLN
1	H	80	GLN
1	J	138	HIS
1	J	180	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 [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	187/187 (100%)	-0.58	0 <a href="#">100</a> <a href="#">100</a>	23, 51, 96, 124	0
1	D	187/187 (100%)	-0.53	0 <a href="#">100</a> <a href="#">100</a>	19, 57, 96, 120	0
1	E	187/187 (100%)	-0.56	0 <a href="#">100</a> <a href="#">100</a>	18, 58, 101, 189	0
1	F	187/187 (100%)	-0.59	1 (0%) <a href="#">87</a> <a href="#">76</a>	23, 57, 90, 161	0
1	H	187/187 (100%)	-0.58	1 (0%) <a href="#">87</a> <a href="#">76</a>	12, 52, 95, 145	0
1	J	187/187 (100%)	-0.56	0 <a href="#">100</a> <a href="#">100</a>	25, 58, 117, 199	0
2	M	92/92 (100%)	-0.49	1 (1%) <a href="#">77</a> <a href="#">63</a>	53, 120, 199, 233	0
All	All	1214/1214 (100%)	-0.56	3 (0%) <a href="#">92</a> <a href="#">85</a>	12, 58, 121, 233	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	167	ALA	4.6
1	F	104	ALA	2.1
2	M	56	ALA	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 oligosaccharides 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.