



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:00 PM GMT

PDB ID : 5C6V  
Title : Crystal structure of the rice Topless related protein 2 (TPR2) N-terminal domain (1-209) in complex with Arabidopsis NINJA peptide  
Authors : Ke, J.; Ma, H.; Gu, X.; Brunzelle, J.S.; Xu, H.E.; Melcher, K.  
Deposited on : 2015-06-23  
Resolution : 3.10 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
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) : rb-20026982

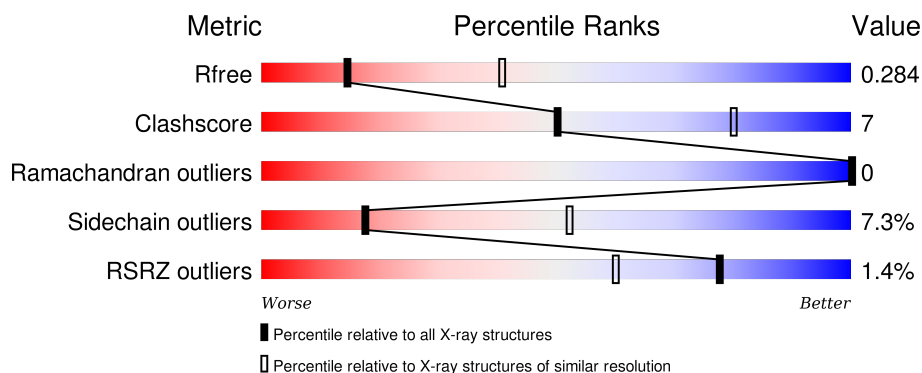
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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	1114 (3.14-3.06)
Clashscore	102246	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)

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  $\geq 3$ , 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	209	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>22%</div> <div>• •</div> </div> </div>
1	B	209	<div> <div>%</div> <div> <div></div> <div>74%</div> <div>21%</div> <div>• •</div> </div> </div>
1	C	209	<div> <div></div> <div> <div></div> <div>74%</div> <div>20%</div> <div>• 5%</div> </div> </div>
1	D	209	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>16%</div> <div>• •</div> </div> </div>
2	E	11	<div> <div>9%</div> <div> <div></div> <div>55%</div> <div>9%</div> <div>36%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	F	11	<div><div></div><div></div><div></div><div>36%27%36%</div></div>
2	G	11	<div><div></div><div></div><div></div><div>55%9%36%</div></div>
2	H	11	<div><div></div><div></div><div></div><div>9%36%18%9%36%</div></div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6914 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 ASPR2 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	202	Total	C	N	O	S	0	0	0
			1689	1091	281	313	4			
1	B	202	Total	C	N	O	S	0	0	0
			1693	1093	281	313	6			
1	C	198	Total	C	N	O	S	0	0	0
			1645	1061	273	306	5			
1	D	200	Total	C	N	O	S	0	0	0
			1665	1076	275	308	6			

- Molecule 2 is a protein called AFP homolog 2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	7	Total	C	N	O	0	0	0
			47	30	7	10			
2	F	7	Total	C	N	O	0	0	0
			47	30	7	10			
2	G	7	Total	C	N	O	0	0	0
			47	30	7	10			
2	H	7	Total	C	N	O	0	0	0
			47	30	7	10			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	14	Total	O	0	0
			14	14		
3	B	7	Total	O	0	0
			7	7		
3	C	6	Total	O	0	0
			6	6		
3	D	6	Total	O	0	0
			6	6		

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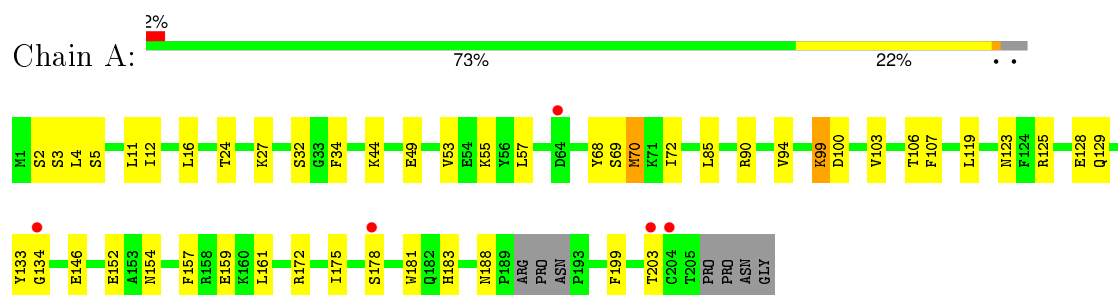
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	O	0	0
			1	1		

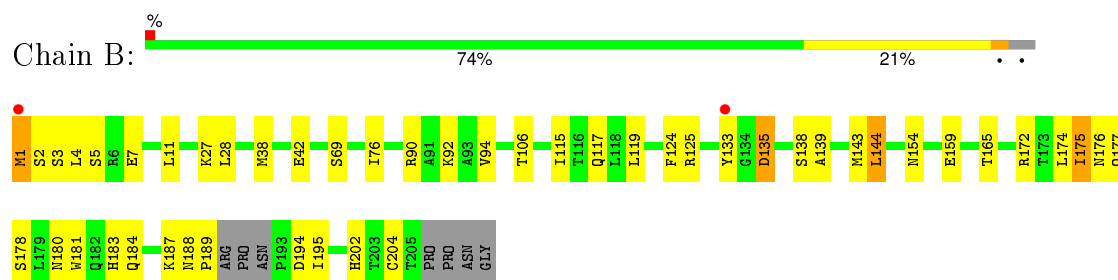
### 3 Residue-property plots [i](#)

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 ( $\text{RSRZ} > 2$ ). 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.

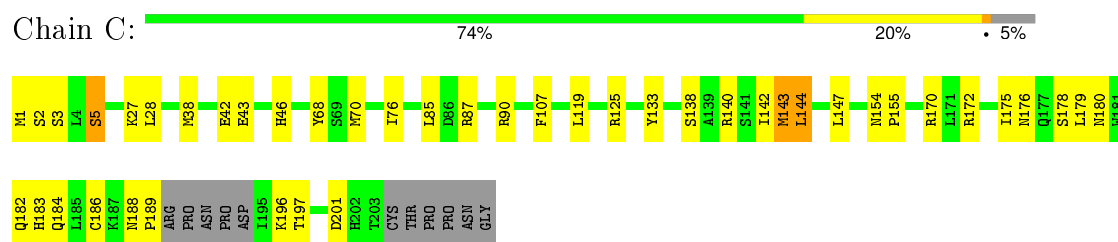
#### • Molecule 1: ASPR2 protein



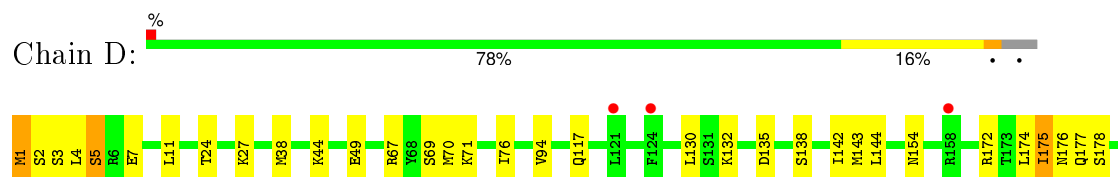
#### • Molecule 1: ASPR2 protein

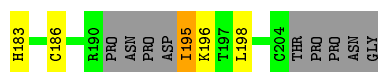


#### • Molecule 1: ASPR2 protein

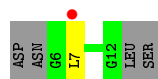


#### • Molecule 1: ASPR2 protein





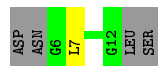
- Molecule 2: AFP homolog 2



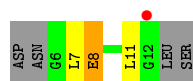
- Molecule 2: AFP homolog 2



- Molecule 2: AFP homolog 2



- Molecule 2: AFP homolog 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	81.74Å 65.04Å 107.73Å 90.00° 105.40° 90.00°	Depositor
Resolution (Å)	48.49 – 3.10 48.50 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.49-3.10) 99.7 (48.50-3.10)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.16 (at 3.12Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.232 , 0.286 0.237 , 0.284	Depositor DCC
$R_{free}$ test set	973 reflections (5.10%)	DCC
Wilson B-factor (Å <sup>2</sup> )	54.0	Xtriage
Anisotropy	0.936	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 49.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 20040 reflections	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	6914	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.84% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.



## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z  > 5$	RMSZ	# $ Z  > 5$
1	A	0.44	0/1722	0.75	3/2311 (0.1%)
1	B	0.43	0/1725	0.66	1/2313 (0.0%)
1	C	0.44	0/1675	0.69	0/2248
1	D	0.38	0/1696	0.60	1/2276 (0.0%)
2	E	0.29	0/46	0.76	0/60
2	F	0.34	0/46	0.75	0/60
2	G	0.38	0/46	0.60	0/60
2	H	0.29	0/46	0.68	0/60
All	All	0.42	0/7002	0.68	5/9388 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	ASN	N-CA-C	-5.60	95.88	111.00
1	A	70	MET	CA-CB-CG	5.48	122.61	113.30
1	B	4	LEU	CA-CB-CG	5.47	127.89	115.30
1	D	4	LEU	CA-CB-CG	5.47	127.89	115.30
1	A	134	GLY	N-CA-C	5.43	126.67	113.10

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	1689	0	1688	37	0
1	B	1693	0	1709	31	0
1	C	1645	0	1641	23	0
1	D	1665	0	1664	24	0
2	E	47	0	49	2	0
2	F	47	0	49	3	0
2	G	47	0	49	1	0
2	H	47	0	49	3	0
3	A	14	0	0	0	0
3	B	7	0	0	0	0
3	C	6	0	0	0	0
3	D	6	0	0	0	0
3	G	1	0	0	0	0
All	All	6914	0	6898	96	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 (96) 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:69:SER:HB3	1:A:154:ASN:HD21	1.53	0.72
1:A:128:GLU:HG3	1:A:129:GLN:HG3	1.71	0.71
1:C:27:LYS:HG3	1:D:27:LYS:HG3	1.75	0.69
1:A:159:GLU:OE1	1:A:159:GLU:N	2.19	0.68
1:A:5:SER:HB3	1:B:172:ARG:HH22	1.59	0.68
1:A:128:GLU:HG3	1:A:129:GLN:H	1.67	0.59
1:A:69:SER:HA	1:A:72:ILE:HD12	1.85	0.59
1:B:1:MET:HG3	1:B:5:SER:HB2	1.84	0.58
1:A:199:PHE:CZ	1:B:7:GLU:HG2	2.39	0.58
1:C:179:LEU:O	1:C:183:HIS:N	2.34	0.58
1:D:1:MET:HG3	1:D:5:SER:HB2	1.86	0.57
1:D:76:ILE:HA	1:D:143:MET:CE	2.34	0.57
1:A:4:LEU:HD23	1:B:172:ARG:HH12	1.69	0.57
1:B:115:ILE:HG12	2:F:11:LEU:HD21	1.86	0.56
1:C:5:SER:HB3	1:D:172:ARG:HH22	1.70	0.56
1:B:188:ASN:HB3	1:B:204:CYS:SG	2.45	0.55
1:A:172:ARG:NH2	1:B:5:SER:HB3	2.21	0.55
1:B:135:ASP:HB3	1:B:138:SER:H	1.70	0.55
1:A:69:SER:HB3	1:A:154:ASN:ND2	2.21	0.54
1:C:5:SER:CB	1:D:172:ARG:HH22	2.21	0.54
1:D:44:LYS:HD2	1:D:49:GLU:HG3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:SER:HB2	1:B:177:GLN:HB2	1.89	0.53
1:C:1:MET:HG3	1:C:5:SER:HB2	1.90	0.53
1:D:76:ILE:HA	1:D:143:MET:HE1	1.90	0.53
1:A:128:GLU:HG3	1:A:129:GLN:N	2.24	0.53
1:B:188:ASN:N	1:B:189:PRO:HD3	2.23	0.53
1:C:182:GLN:NE2	1:D:177:GLN:OE1	2.43	0.52
1:A:27:LYS:HG3	1:B:27:LYS:HG3	1.90	0.52
1:C:46:HIS:O	1:C:140:ARG:NH1	2.42	0.52
1:C:179:LEU:O	1:C:183:HIS:HB2	2.10	0.52
1:A:24:THR:HG21	1:B:28:LEU:HD13	1.92	0.51
1:B:117:GLN:HE21	1:D:117:GLN:HE21	1.57	0.51
1:D:138:SER:O	1:D:142:ILE:HG12	2.11	0.51
1:D:130:LEU:HD11	2:H:11:LEU:HD22	1.92	0.51
1:B:11:LEU:HD21	1:B:174:LEU:HD12	1.93	0.50
1:A:99:LYS:HD2	1:A:100:ASP:OD1	2.12	0.50
1:B:76:ILE:HA	1:B:143:MET:CE	2.42	0.49
1:A:123:ASN:OD1	1:A:125:ARG:HG3	2.12	0.49
1:B:187:LYS:C	1:B:189:PRO:HD3	2.32	0.49
1:A:157:PHE:O	1:A:161:LEU:HG	2.12	0.49
1:D:183:HIS:HA	1:D:186:CYS:SG	2.52	0.48
1:C:76:ILE:HA	1:C:143:MET:CE	2.42	0.48
1:D:195:ILE:CG2	1:D:196:LYS:N	2.77	0.48
1:B:195:ILE:HG22	1:B:202:HIS:CB	2.44	0.48
1:A:4:LEU:HD23	1:B:172:ARG:NH1	2.29	0.48
1:A:49:GLU:O	1:A:53:VAL:HG23	2.14	0.47
1:B:183:HIS:CE1	1:B:189:PRO:HB3	2.50	0.47
1:A:181:TRP:HB2	1:B:181:TRP:HB2	1.97	0.47
1:A:152:GLU:HA	1:A:161:LEU:HD11	1.97	0.46
1:A:199:PHE:CE1	1:B:7:GLU:HG2	2.50	0.46
1:B:125:ARG:NE	1:B:133:TYR:O	2.48	0.46
1:A:85:LEU:HD21	1:A:119:LEU:HD23	1.97	0.46
1:C:42:GLU:HG2	1:C:144:LEU:HD11	1.98	0.45
1:A:44:LYS:HD2	1:A:49:GLU:HG3	1.99	0.45
1:C:28:LEU:HD13	1:D:24:THR:HG21	1.98	0.45
1:A:94:VAL:HG11	1:C:90:ARG:HD2	1.97	0.45
1:C:180:ASN:O	1:C:184:GLN:CB	2.64	0.45
1:C:197:THR:HG22	1:D:7:GLU:OE2	2.17	0.45
1:B:90:ARG:HD2	1:D:94:VAL:HG11	1.98	0.45
2:F:11:LEU:HD23	2:F:11:LEU:HA	1.71	0.45
1:B:42:GLU:HG2	1:B:144:LEU:HD11	1.99	0.45
1:B:69:SER:HB3	1:B:154:ASN:HD21	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:LEU:HA	1:A:57:LEU:HD23	1.82	0.44
1:C:125:ARG:NH2	1:C:133:TYR:HD1	2.16	0.44
1:C:172:ARG:HH22	1:D:5:SER:HB3	1.81	0.44
1:D:175:ILE:HA	1:D:175:ILE:HD12	1.84	0.44
2:F:8:GLU:CD	2:F:8:GLU:H	2.21	0.43
1:A:133:TYR:CD1	1:A:133:TYR:C	2.91	0.43
1:D:174:LEU:HA	1:D:174:LEU:HD23	1.87	0.43
1:D:69:SER:HB3	1:D:154:ASN:HD21	1.83	0.43
1:A:32:SER:HB2	1:A:34:PHE:CD2	2.54	0.43
1:B:175:ILE:HA	1:B:175:ILE:HD12	1.82	0.43
1:A:68:TYR:CE1	2:E:7:LEU:HD21	2.54	0.42
1:C:182:GLN:NE2	1:C:201:ASP:OD1	2.52	0.42
1:D:71:LYS:HD3	2:H:7:LEU:HB3	2.01	0.42
1:A:55:LYS:HE2	1:A:55:LYS:HB3	1.68	0.42
1:C:68:TYR:CE1	2:G:7:LEU:HD11	2.54	0.42
1:D:11:LEU:HD21	1:D:174:LEU:HD12	2.02	0.42
1:A:68:TYR:CD1	2:E:7:LEU:HD21	2.55	0.42
1:A:183:HIS:O	1:A:183:HIS:ND1	2.53	0.42
1:A:175:ILE:HD13	1:A:175:ILE:HA	1.93	0.42
1:B:133:TYR:CE2	1:B:139:ALA:HB1	2.54	0.41
1:A:12:ILE:O	1:A:16:LEU:HG	2.20	0.41
1:A:125:ARG:NE	1:A:133:TYR:O	2.51	0.41
1:A:106:THR:OG1	1:A:107:PHE:N	2.52	0.41
1:C:170:ARG:NE	1:D:198:LEU:O	2.49	0.41
1:C:85:LEU:HD21	1:C:119:LEU:HD23	2.03	0.41
1:C:154:ASN:HA	1:C:155:PRO:HD2	1.94	0.41
1:C:138:SER:O	1:C:142:ILE:HG12	2.21	0.41
1:A:181:TRP:NE1	1:B:180:ASN:HB3	2.36	0.41
1:B:124:PHE:HZ	1:B:133:TYR:CE1	2.39	0.41
2:H:8:GLU:HG2	2:H:8:GLU:H	1.47	0.41
1:D:195:ILE:HG22	1:D:196:LYS:N	2.36	0.40
1:C:188:ASN:HA	1:C:189:PRO:HA	1.82	0.40
1:B:94:VAL:HG22	1:B:119:LEU:HD11	2.03	0.40
1:B:92:LYS:HA	1:B:92:LYS:HD2	1.70	0.40

There are no symmetry-related clashes.

## 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	198/209 (95%)	196 (99%)	2 (1%)	0	100	100
1	B	198/209 (95%)	196 (99%)	2 (1%)	0	100	100
1	C	194/209 (93%)	192 (99%)	2 (1%)	0	100	100
1	D	196/209 (94%)	193 (98%)	3 (2%)	0	100	100
2	E	5/11 (46%)	5 (100%)	0	0	100	100
2	F	5/11 (46%)	5 (100%)	0	0	100	100
2	G	5/11 (46%)	5 (100%)	0	0	100	100
2	H	5/11 (46%)	5 (100%)	0	0	100	100
All	All	806/880 (92%)	797 (99%)	9 (1%)	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	186/197 (94%)	177 (95%)	9 (5%)	31	69
1	B	189/197 (96%)	175 (93%)	14 (7%)	17	51
1	C	180/197 (91%)	164 (91%)	16 (9%)	12	42
1	D	183/197 (93%)	169 (92%)	14 (8%)	16	50
2	E	5/9 (56%)	5 (100%)	0	100	100
2	F	5/9 (56%)	4 (80%)	1 (20%)	1	7

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	G	5/9 (56%)	5 (100%)	0	100	100
2	H	5/9 (56%)	4 (80%)	1 (20%)	1	7
All	All	758/824 (92%)	703 (93%)	55 (7%)	17	52

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	3	SER
1	A	11	LEU
1	A	70	MET
1	A	90	ARG
1	A	99	LYS
1	A	103	VAL
1	A	146	GLU
1	A	203	THR
1	B	1	MET
1	B	2	SER
1	B	3	SER
1	B	38	MET
1	B	106	THR
1	B	135	ASP
1	B	144	LEU
1	B	159	GLU
1	B	165	THR
1	B	175	ILE
1	B	176	ASN
1	B	178	SER
1	B	184	GLN
1	B	194	ASP
1	C	2	SER
1	C	3	SER
1	C	5	SER
1	C	38	MET
1	C	43	GLU
1	C	70	MET
1	C	87	ARG
1	C	107	PHE
1	C	143	MET
1	C	144	LEU
1	C	147	LEU

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Mol	Chain	Res	Type
1	C	175	ILE
1	C	176	ASN
1	C	178	SER
1	C	186	CYS
1	C	196	LYS
1	D	1	MET
1	D	2	SER
1	D	3	SER
1	D	5	SER
1	D	38	MET
1	D	67	ARG
1	D	70	MET
1	D	132	LYS
1	D	135	ASP
1	D	144	LEU
1	D	175	ILE
1	D	176	ASN
1	D	178	SER
1	D	195	ILE
2	F	10	SER
2	H	8	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	30	GLN
1	B	117	GLN
1	C	117	GLN
1	C	182	GLN
1	D	177	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 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, 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	202/209 (96%)	0.01	5 (2%) 61 37	26, 41, 99, 139	0
1	B	202/209 (96%)	-0.04	2 (0%) 84 69	36, 53, 98, 151	0
1	C	198/209 (94%)	-0.09	0 100 100	25, 46, 105, 142	0
1	D	200/209 (95%)	0.11	3 (1%) 76 58	43, 69, 108, 128	0
2	E	7/11 (63%)	0.31	1 (14%) 4 2	46, 54, 65, 67	0
2	F	7/11 (63%)	0.04	0 100 100	52, 58, 69, 78	0
2	G	7/11 (63%)	0.39	0 100 100	42, 50, 56, 65	0
2	H	7/11 (63%)	0.91	1 (14%) 4 2	93, 102, 108, 111	0
All	All	830/880 (94%)	0.01	12 (1%) 78 60	25, 53, 104, 151	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	204	CYS	5.0
1	A	203	THR	4.4
2	H	12	GLY	3.3
1	B	1	MET	2.8
1	A	178	SER	2.7
1	D	158	ARG	2.6
2	E	7	LEU	2.6
1	D	121	LEU	2.4
1	B	133	TYR	2.3
1	D	124	PHE	2.3
1	A	64	ASP	2.3
1	A	134	GLY	2.2

## 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.