



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 17, 2017 – 07:01 PM EDT

PDB ID : 5C7E  
Title : Crystal structure of the rice Topless related protein 2 (TPR2) N-terminal domain (1-209) in complex with Arabidopsis IAA10 peptide  
Authors : Ke, J.; Ma, H.; Gu, X.; Brunzelle, J.S.; Xu, H.E.; Melcher, K.  
Deposited on : unknown  
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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029824
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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-20029824

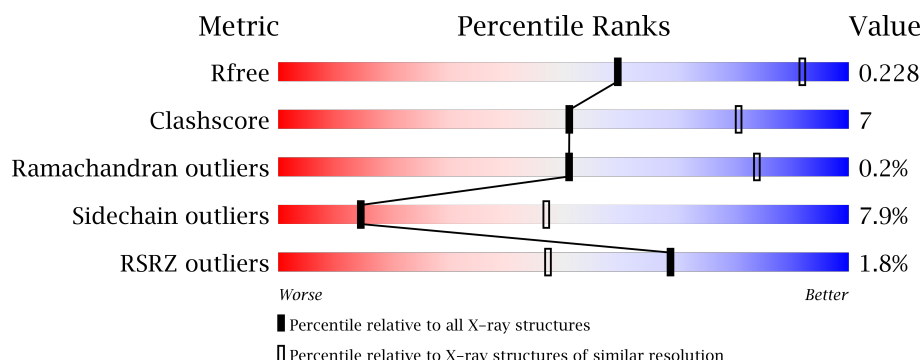
# 1 Overall quality at a glance

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.





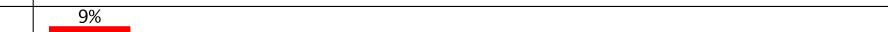
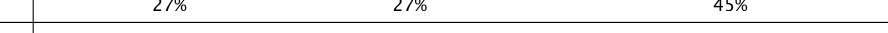



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	100719	1001 (3.12-3.08)
Clashscore	112137	1099 (3.12-3.08)
Ramachandran outliers	110173	1057 (3.12-3.08)
Sidechain outliers	110143	1057 (3.12-3.08)
RSRZ outliers	101464	1006 (3.12-3.08)

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>0.2%</div> <div> <div></div> <div>76%</div> <div>20%</div> <div>..</div> </div> </div>
1	B	209	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>..</div> </div> </div>
1	C	209	<div> <div>0.2%</div> <div> <div></div> <div>80%</div> <div>16%</div> <div>..</div> </div> </div>
1	D	209	<div> <div>0.2%</div> <div> <div></div> <div>78%</div> <div>18%</div> <div>..</div> </div> </div>
1	E	209	<div> <div>2%</div> <div> <div></div> <div>72%</div> <div>24%</div> <div>...</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	209	
2	G	11	
2	H	11	
2	I	11	
2	J	11	
2	K	11	
2	L	11	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10658 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	205	Total	C	N	O	S	0	0	0
			1710	1103	287	314	6			
1	B	205	Total	C	N	O	S	0	0	0
			1704	1101	281	316	6			
1	C	206	Total	C	N	O	S	0	0	0
			1725	1113	288	318	6			
1	D	206	Total	C	N	O	S	0	0	0
			1720	1108	287	319	6			
1	E	205	Total	C	N	O	S	0	0	0
			1714	1105	286	317	6			
1	F	205	Total	C	N	O	S	0	0	0
			1718	1108	287	317	6			

- Molecule 2 is a protein called Auxin-responsive protein IAA10.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	10	Total	C	N	O	0	0	0
			65	41	10	14			
2	H	6	Total	C	N	O	0	0	0
			43	29	6	8			
2	I	9	Total	C	N	O	0	0	0
			68	42	9	17			
2	J	9	Total	C	N	O	0	0	0
			57	37	9	11			
2	K	7	Total	C	N	O	0	0	0
			46	31	7	8			
2	L	8	Total	C	N	O	0	0	0
			59	38	8	13			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	Total Zn 1 1	0	0
3	E	1	Total Zn 1 1	0	0
3	B	1	Total Zn 1 1	0	0
3	C	1	Total Zn 1 1	0	0
3	A	1	Total Zn 1 1	0	0
3	F	1	Total Zn 1 1	0	0

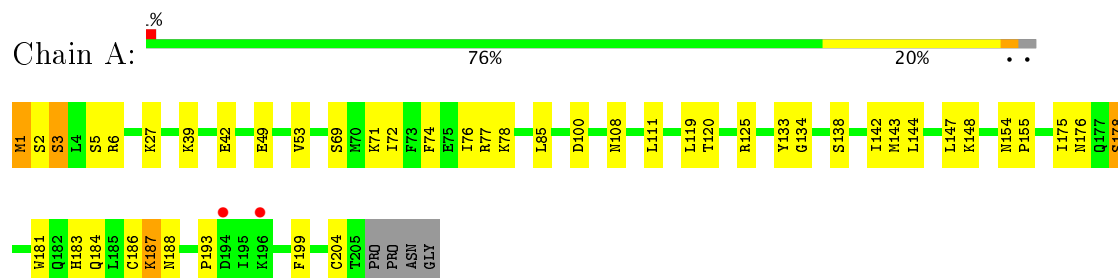
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total O 4 4	0	0
4	B	1	Total O 1 1	0	0
4	C	7	Total O 7 7	0	0
4	D	4	Total O 4 4	0	0
4	E	5	Total O 5 5	0	0
4	F	1	Total O 1 1	0	0
4	L	1	Total O 1 1	0	0

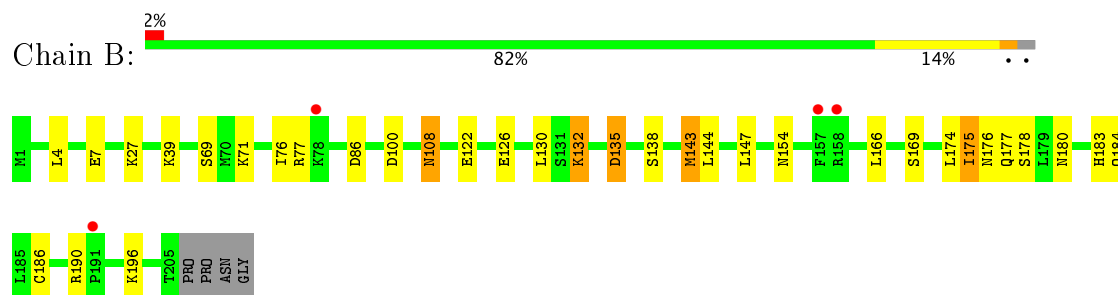
### 3 Residue-property plots [i](#)

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

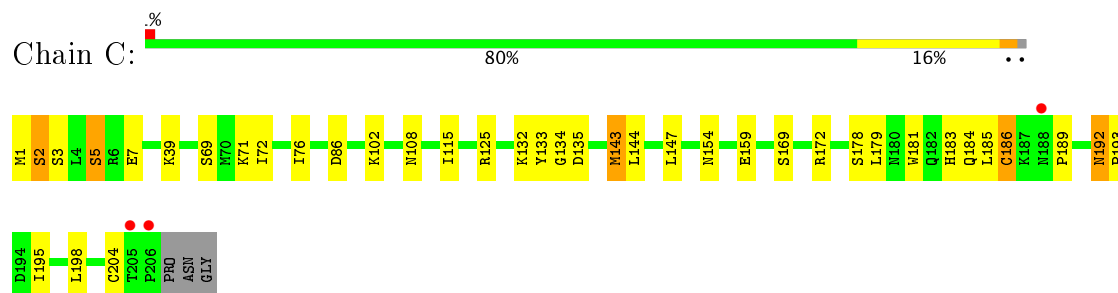
#### • Molecule 1: ASPR2 protein



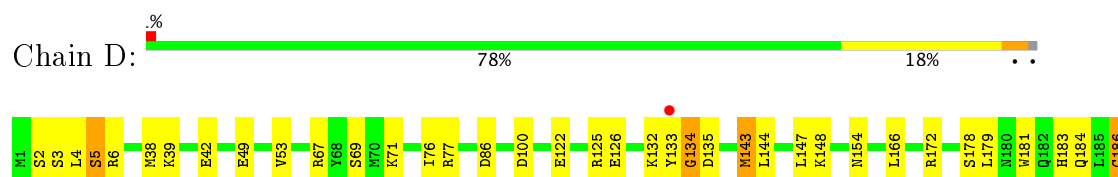
#### • Molecule 1: ASPR2 protein



#### • Molecule 1: ASPR2 protein

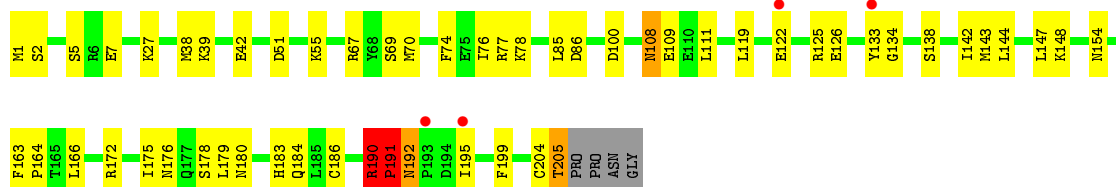
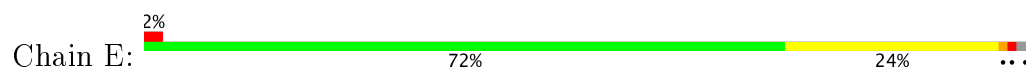


#### • Molecule 1: ASPR2 protein

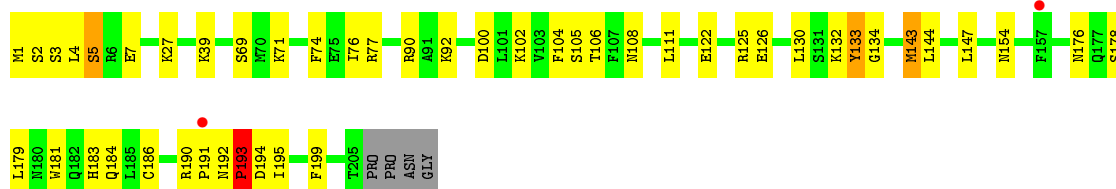
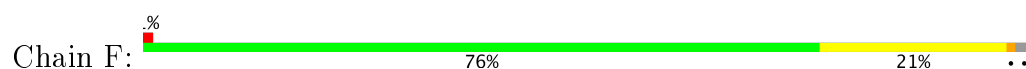




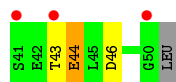
- Molecule 1: ASPR2 protein



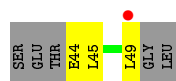
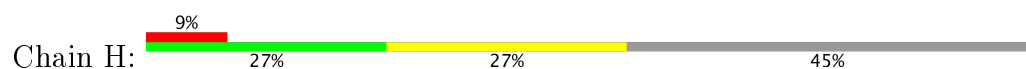
- Molecule 1: ASPR2 protein



- Molecule 2: Auxin-responsive protein IAA10



- Molecule 2: Auxin-responsive protein IAA10



- Molecule 2: Auxin-responsive protein IAA10



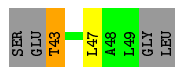
- Molecule 2: Auxin-responsive protein IAA10



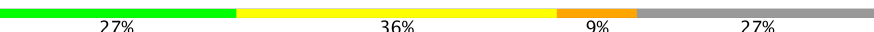


- Molecule 2: Auxin-responsive protein IAA10

Chain K: 



- Molecule 2: Auxin-responsive protein IAA10

Chain L: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	162.67Å 162.67Å 157.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.99 – 3.10 44.99 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.99-3.10) 99.9 (44.99-3.10)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.28 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.196 , 0.229 0.196 , 0.228	Depositor DCC
$R_{free}$ test set	2236 reflections (5.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	98.5	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 63.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10658	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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

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

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	0.40	0/1745	0.65	0/2344
1	B	0.36	0/1739	0.64	1/2338 (0.0%)
1	C	0.44	0/1761	0.69	1/2367 (0.0%)
1	D	0.39	0/1755	0.68	3/2359 (0.1%)
1	E	0.47	1/1749 (0.1%)	0.72	3/2351 (0.1%)
1	F	0.39	0/1753	0.63	0/2355
2	G	0.38	0/64	0.68	0/86
2	H	0.36	0/42	0.79	0/56
2	I	0.65	0/67	1.21	0/90
2	J	0.51	0/56	1.06	0/76
2	K	0.32	0/45	0.79	0/61
2	L	0.36	0/58	0.91	0/78
All	All	0.41	1/10834 (0.0%)	0.68	8/14561 (0.1%)

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	C	0	1
1	D	0	3
1	F	0	1
2	K	0	1
All	All	0	6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	191	PRO	N-CD	5.87	1.56	1.47

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	190	ARG	N-CA-C	7.60	131.53	111.00
1	E	192	ASN	N-CA-C	-6.06	94.64	111.00
1	D	190	ARG	N-CA-C	-5.71	95.59	111.00
1	E	190	ARG	N-CA-CB	-5.34	100.98	110.60
1	D	135	ASP	N-CA-C	5.22	125.10	111.00
1	B	190	ARG	N-CA-C	-5.20	96.96	111.00
1	D	186	CYS	CA-CB-SG	5.19	123.34	114.00
1	C	134	GLY	N-CA-C	-5.15	100.23	113.10

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	192	ASN	Peptide
1	D	132	LYS	Peptide
1	D	134	GLY	Peptide
1	D	188	ASN	Peptide
1	F	193	PRO	Peptide
2	K	43	THR	Peptide

## 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	1710	0	1707	27	0
1	B	1704	0	1696	18	0
1	C	1725	0	1730	25	0
1	D	1720	0	1727	24	0
1	E	1714	0	1712	33	1
1	F	1718	0	1723	26	1
2	G	65	0	62	3	0
2	H	43	0	45	3	0
2	I	68	0	65	3	0
2	J	57	0	55	1	0
2	K	46	0	48	0	0
2	L	59	0	58	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	4	0	0	0	0
4	B	1	0	0	0	0
4	C	7	0	0	1	0
4	D	4	0	0	0	0
4	E	5	0	0	0	0
4	F	1	0	0	0	0
4	L	1	0	0	0	0
All	All	10658	0	10628	139	1

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 (139) 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:C:189:PRO:HB3	1:C:204:CYS:SG	2.18	0.83
1:A:71:LYS:HE2	2:G:44:GLU:HG2	1.62	0.81
1:F:190:ARG:HB3	1:F:191:PRO:HD2	1.63	0.79
1:F:102:LYS:O	1:F:105:SER:HB2	1.84	0.77
1:E:190:ARG:HB3	1:E:191:PRO:HD3	1.64	0.77
1:E:1:MET:HB3	1:E:2:SER:HA	1.67	0.74
1:E:199:PHE:CZ	1:F:7:GLU:HG2	2.30	0.67
1:C:183:HIS:O	1:C:186:CYS:HB2	1.95	0.66
1:E:125:ARG:HH21	1:E:134:GLY:HA3	1.59	0.66
1:E:183:HIS:HA	1:E:186:CYS:SG	2.34	0.66
1:B:135:ASP:HB3	1:B:138:SER:H	1.61	0.65
1:C:76:ILE:HA	1:C:143:MET:CE	2.29	0.63
1:C:7:GLU:HG2	1:D:199:PHE:CZ	2.33	0.63
1:C:181:TRP:HB2	1:D:181:TRP:HB2	1.80	0.63
1:F:108:ASN:HB3	1:F:111:LEU:HB3	1.79	0.62
1:C:76:ILE:HA	1:C:143:MET:HE1	1.81	0.62
1:D:3:SER:HB3	1:D:6:ARG:HB3	1.81	0.62
1:D:125:ARG:NH2	1:D:133:TYR:HD2	1.97	0.62
1:E:1:MET:CB	1:E:2:SER:HA	2.29	0.61
1:F:1:MET:SD	1:F:4:LEU:HD23	2.41	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:LEU:O	1:B:7:GLU:HB2	2.02	0.60
1:B:71:LYS:HE2	2:H:44:GLU:O	2.02	0.59
1:E:163:PHE:CG	1:E:164:PRO:HD2	2.38	0.59
1:F:71:LYS:HE2	2:L:44:GLU:O	2.02	0.58
1:A:69:SER:HB3	1:A:154:ASN:HD21	1.68	0.57
1:E:76:ILE:HA	1:E:143:MET:CE	2.34	0.57
1:F:190:ARG:HB3	1:F:191:PRO:CD	2.33	0.57
1:E:179:LEU:HB3	1:E:195:ILE:HD13	1.88	0.56
1:B:183:HIS:HA	1:B:186:CYS:SG	2.46	0.56
1:C:69:SER:HB3	1:C:154:ASN:HD21	1.71	0.56
1:E:191:PRO:CG	1:E:192:ASN:H	2.18	0.56
1:C:115:ILE:HG12	2:I:49:LEU:HD21	1.88	0.56
1:B:122:GLU:HB3	1:B:126:GLU:OE2	2.06	0.55
1:F:69:SER:HB3	1:F:154:ASN:HD21	1.71	0.55
1:C:5:SER:HB3	1:D:172:ARG:HH22	1.71	0.55
1:A:125:ARG:HH21	1:A:134:GLY:CA	2.20	0.55
1:F:76:ILE:HA	1:F:143:MET:CE	2.36	0.55
1:C:179:LEU:HD13	1:C:195:ILE:HD13	1.89	0.55
1:A:3:SER:HG	1:A:6:ARG:H	1.55	0.55
1:E:69:SER:HB3	1:E:154:ASN:HD21	1.73	0.55
1:C:125:ARG:NE	1:C:133:TYR:O	2.37	0.54
1:E:7:GLU:HG2	1:F:199:PHE:CZ	2.41	0.54
1:E:191:PRO:HG2	1:E:192:ASN:H	1.73	0.54
1:A:199:PHE:CZ	1:B:7:GLU:HG2	2.43	0.54
1:D:76:ILE:HA	1:D:143:MET:CE	2.38	0.54
1:D:188:ASN:CG	1:D:189:PRO:HD3	2.27	0.54
1:A:108:ASN:HB3	1:A:111:LEU:HB3	1.90	0.54
1:B:130:LEU:HD21	2:H:49:LEU:HB3	1.89	0.53
1:E:76:ILE:HA	1:E:143:MET:HE1	1.90	0.53
1:D:76:ILE:HA	1:D:143:MET:HE1	1.90	0.52
1:D:42:GLU:OE2	1:D:148:LYS:HE2	2.10	0.52
1:A:42:GLU:OE2	1:A:148:LYS:HE2	2.10	0.52
1:E:180:ASN:HB3	1:F:181:TRP:NE1	2.25	0.52
1:A:71:LYS:HD2	2:G:46:ASP:O	2.10	0.52
1:F:183:HIS:HA	1:F:186:CYS:SG	2.50	0.51
1:A:1:MET:HB2	1:B:196:LYS:O	2.10	0.51
1:E:108:ASN:HB3	1:E:111:LEU:HB3	1.92	0.51
1:A:76:ILE:HA	1:A:143:MET:HE1	1.92	0.51
1:E:133:TYR:CD1	1:E:134:GLY:N	2.79	0.50
1:A:133:TYR:CD1	1:A:134:GLY:N	2.79	0.50
1:F:125:ARG:NE	1:F:133:TYR:O	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:74:PHE:HB2	1:F:104:PHE:CE2	2.46	0.50
1:A:125:ARG:HH21	1:A:134:GLY:HA3	1.75	0.50
1:C:183:HIS:HE1	1:C:189:PRO:CB	2.24	0.50
1:A:120:THR:O	1:C:102:LYS:NZ	2.41	0.49
1:E:125:ARG:HH21	1:E:134:GLY:CA	2.23	0.49
1:A:138:SER:O	1:A:142:ILE:HG12	2.12	0.49
1:A:76:ILE:HA	1:A:143:MET:CE	2.43	0.49
1:E:85:LEU:HD21	1:E:119:LEU:HD23	1.94	0.49
1:C:135:ASP:OD1	4:C:501:HOH:O	2.20	0.49
1:A:27:LYS:HG3	1:B:27:LYS:HG3	1.96	0.48
1:E:204:CYS:O	1:E:205:THR:HG22	2.13	0.48
2:I:42:GLU:O	2:I:42:GLU:HG2	2.14	0.48
1:D:183:HIS:CE1	1:D:193:PRO:HB3	2.49	0.48
1:A:49:GLU:O	1:A:53:VAL:HG23	2.13	0.48
1:C:1:MET:SD	1:C:3:SER:HB3	2.54	0.48
1:B:76:ILE:HA	1:B:143:MET:CE	2.43	0.47
1:E:190:ARG:CB	1:E:191:PRO:HD3	2.38	0.47
1:A:186:CYS:HB2	1:A:204:CYS:H	1.79	0.47
1:D:179:LEU:HB3	1:D:195:ILE:HD13	1.96	0.47
1:E:27:LYS:HG3	1:F:27:LYS:HG3	1.96	0.47
2:G:44:GLU:HG2	2:G:44:GLU:O	2.14	0.47
1:D:69:SER:HB3	1:D:154:ASN:HD21	1.80	0.47
1:C:2:SER:HA	1:C:5:SER:OG	2.14	0.47
1:C:192:ASN:CB	1:C:193:PRO:HA	2.45	0.46
1:C:125:ARG:NH2	1:C:133:TYR:HD1	2.13	0.46
1:E:179:LEU:HD22	1:E:195:ILE:HG21	1.96	0.46
1:F:71:LYS:HD3	2:L:45:LEU:HA	1.98	0.46
1:C:172:ARG:HH22	1:D:5:SER:CB	2.27	0.46
1:B:69:SER:HB3	1:B:154:ASN:HD21	1.81	0.46
1:B:132:LYS:HG2	1:B:132:LYS:H	1.48	0.45
1:F:179:LEU:HD13	1:F:195:ILE:HD13	1.98	0.45
1:C:198:LEU:HD11	1:D:4:LEU:CD1	2.46	0.45
1:F:76:ILE:HA	1:F:143:MET:HE2	1.97	0.45
1:C:71:LYS:HE2	2:I:44:GLU:O	2.16	0.45
1:A:178:SER:HB3	1:B:177:GLN:HB2	1.98	0.45
1:D:189:PRO:HB2	1:D:190:ARG:O	2.17	0.45
1:D:122:GLU:HB3	1:D:126:GLU:OE2	2.17	0.45
1:D:77:ARG:HD3	1:D:100:ASP:HB3	1.98	0.45
1:A:74:PHE:CE2	1:A:78:LYS:HE3	2.52	0.44
1:C:183:HIS:HE1	1:C:189:PRO:HB2	1.82	0.44
1:F:77:ARG:HD3	1:F:100:ASP:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:HIS:CE1	1:A:193:PRO:HG3	2.53	0.44
1:A:77:ARG:HD3	1:A:100:ASP:HB3	1.99	0.44
1:A:85:LEU:HD21	1:A:119:LEU:HD23	2.00	0.43
1:E:1:MET:HE2	1:E:7:GLU:OE1	2.18	0.43
1:F:122:GLU:HB3	1:F:126:GLU:OE2	2.18	0.43
1:A:181:TRP:NE1	1:B:180:ASN:HB3	2.33	0.43
1:A:187:LYS:HA	1:A:188:ASN:HA	1.57	0.43
2:H:45:LEU:HD12	2:H:45:LEU:O	2.18	0.43
1:E:77:ARG:HD3	1:E:100:ASP:HB3	1.99	0.43
1:E:172:ARG:HH22	1:F:5:SER:HB3	1.83	0.43
1:B:77:ARG:HD3	1:B:100:ASP:HB3	2.00	0.43
1:E:51:ASP:O	1:E:55:LYS:HB2	2.19	0.43
1:C:7:GLU:HG2	1:D:199:PHE:CE1	2.53	0.43
1:D:187:LYS:HG2	1:D:187:LYS:H	1.70	0.43
1:E:190:ARG:CB	1:E:191:PRO:CD	2.97	0.43
1:F:1:MET:C	1:F:3:SER:H	2.22	0.43
1:D:183:HIS:HA	1:D:186:CYS:SG	2.60	0.42
1:F:92:LYS:HA	1:F:92:LYS:HD2	1.85	0.42
1:D:71:LYS:HE3	2:J:46:ASP:O	2.19	0.42
1:A:154:ASN:HA	1:A:155:PRO:HD2	1.92	0.42
1:D:49:GLU:O	1:D:53:VAL:HG23	2.21	0.41
1:E:190:ARG:HB3	1:E:191:PRO:CD	2.42	0.41
1:C:69:SER:HA	1:C:72:ILE:HD12	2.02	0.41
1:A:69:SER:HA	1:A:72:ILE:HD12	2.03	0.41
1:E:122:GLU:HB3	1:E:126:GLU:OE2	2.21	0.41
1:B:174:LEU:HA	1:B:174:LEU:HD23	1.85	0.41
1:C:183:HIS:CE1	1:C:189:PRO:CB	3.03	0.41
1:D:125:ARG:HH21	1:D:134:GLY:HA3	1.85	0.41
1:E:74:PHE:CE2	1:E:78:LYS:HE3	2.56	0.41
1:F:125:ARG:NH2	1:F:134:GLY:O	2.54	0.41
1:B:175:ILE:HA	1:B:175:ILE:HD12	1.80	0.41
1:F:130:LEU:HD21	2:L:49:LEU:HB3	2.03	0.41
1:E:138:SER:O	1:E:142:ILE:HG12	2.20	0.41
1:F:192:ASN:HA	1:F:193:PRO:HA	1.79	0.40
1:B:108:ASN:N	1:B:108:ASN:OD1	2.55	0.40
1:E:42:GLU:OE2	1:E:148:LYS:HE2	2.20	0.40
1:D:183:HIS:CE1	1:D:189:PRO:HB3	2.57	0.40

All (1) 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:E:67:ARG:NH1	1:F:193:PRO:O[4_555]	2.05	0.15

## 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	203/209 (97%)	199 (98%)	4 (2%)	0	100	100
1	B	203/209 (97%)	202 (100%)	1 (0%)	0	100	100
1	C	204/209 (98%)	202 (99%)	2 (1%)	0	100	100
1	D	204/209 (98%)	202 (99%)	2 (1%)	0	100	100
1	E	203/209 (97%)	198 (98%)	3 (2%)	2 (1%)	18	57
1	F	203/209 (97%)	199 (98%)	3 (2%)	1 (0%)	32	71
2	G	8/11 (73%)	8 (100%)	0	0	100	100
2	H	4/11 (36%)	4 (100%)	0	0	100	100
2	I	7/11 (64%)	7 (100%)	0	0	100	100
2	J	7/11 (64%)	6 (86%)	1 (14%)	0	100	100
2	K	5/11 (46%)	4 (80%)	1 (20%)	0	100	100
2	L	6/11 (54%)	6 (100%)	0	0	100	100
All	All	1257/1320 (95%)	1237 (98%)	17 (1%)	3 (0%)	51	84

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	191	PRO
1	E	190	ARG
1	F	193	PRO



### 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	188/197 (95%)	176 (94%)	12 (6%)	20	56
1	B	188/197 (95%)	174 (93%)	14 (7%)	16	49
1	C	192/197 (98%)	177 (92%)	15 (8%)	15	48
1	D	192/197 (98%)	179 (93%)	13 (7%)	18	53
1	E	190/197 (96%)	175 (92%)	15 (8%)	14	47
1	F	191/197 (97%)	177 (93%)	14 (7%)	16	50
2	G	6/9 (67%)	4 (67%)	2 (33%)	0	0
2	H	4/9 (44%)	4 (100%)	0	100	100
2	I	8/9 (89%)	7 (88%)	1 (12%)	5	22
2	J	5/9 (56%)	3 (60%)	2 (40%)	0	0
2	K	4/9 (44%)	2 (50%)	2 (50%)	0	0
2	L	6/9 (67%)	3 (50%)	3 (50%)	0	0
All	All	1174/1236 (95%)	1081 (92%)	93 (8%)	14	47

All (93) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	SER
1	A	3	SER
1	A	5	SER
1	A	39	LYS
1	A	144	LEU
1	A	147	LEU
1	A	175	ILE
1	A	176	ASN
1	A	178	SER
1	A	184	GLN
1	A	187	LYS
1	B	39	LYS
1	B	86	ASP

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Mol	Chain	Res	Type
1	B	108	ASN
1	B	132	LYS
1	B	135	ASP
1	B	143	MET
1	B	144	LEU
1	B	147	LEU
1	B	166	LEU
1	B	169	SER
1	B	175	ILE
1	B	176	ASN
1	B	178	SER
1	B	184	GLN
1	C	2	SER
1	C	5	SER
1	C	39	LYS
1	C	86	ASP
1	C	108	ASN
1	C	132	LYS
1	C	143	MET
1	C	144	LEU
1	C	147	LEU
1	C	159	GLU
1	C	169	SER
1	C	178	SER
1	C	184	GLN
1	C	185	LEU
1	C	186	CYS
1	D	2	SER
1	D	5	SER
1	D	38	MET
1	D	39	LYS
1	D	67	ARG
1	D	86	ASP
1	D	143	MET
1	D	144	LEU
1	D	147	LEU
1	D	166	LEU
1	D	178	SER
1	D	184	GLN
1	D	187	LYS
1	E	5	SER
1	E	38	MET

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Mol	Chain	Res	Type
1	E	39	LYS
1	E	70	MET
1	E	86	ASP
1	E	108	ASN
1	E	109	GLU
1	E	144	LEU
1	E	147	LEU
1	E	166	LEU
1	E	175	ILE
1	E	176	ASN
1	E	178	SER
1	E	184	GLN
1	E	205	THR
1	F	2	SER
1	F	5	SER
1	F	39	LYS
1	F	90	ARG
1	F	106	THR
1	F	132	LYS
1	F	133	TYR
1	F	143	MET
1	F	144	LEU
1	F	147	LEU
1	F	176	ASN
1	F	178	SER
1	F	184	GLN
1	F	194	ASP
2	G	43	THR
2	G	44	GLU
2	I	44	GLU
2	J	47	LEU
2	J	49	LEU
2	K	43	THR
2	K	47	LEU
2	L	42	GLU
2	L	43	THR
2	L	45	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

### 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	205/209 (98%)	-0.08	2 (0%) 82 67	69, 86, 163, 195	0
1	B	205/209 (98%)	0.11	4 (1%) 65 44	81, 113, 149, 171	0
1	C	206/209 (98%)	-0.02	3 (1%) 74 54	63, 81, 123, 155	0
1	D	206/209 (98%)	0.10	3 (1%) 74 54	67, 97, 137, 180	0
1	E	205/209 (98%)	0.04	4 (1%) 65 44	78, 98, 155, 194	0
1	F	205/209 (98%)	0.07	2 (0%) 82 67	83, 105, 158, 198	0
2	G	10/11 (90%)	1.14	3 (30%) 1 0	86, 101, 157, 162	0
2	H	6/11 (54%)	0.71	1 (16%) 2 1	127, 141, 150, 155	0
2	I	9/11 (81%)	0.27	0 100 100	81, 110, 148, 150	0
2	J	9/11 (81%)	0.52	1 (11%) 6 2	107, 121, 156, 158	0
2	K	7/11 (63%)	-0.28	0 100 100	96, 111, 138, 146	0
2	L	8/11 (72%)	0.60	0 100 100	119, 137, 163, 180	0
All	All	1281/1320 (97%)	0.06	23 (1%) 69 47	63, 98, 154, 198	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	191	PRO	6.1
1	C	206	PRO	3.8
1	D	206	PRO	3.2
1	B	157	PHE	3.0
2	G	43	THR	3.0
2	G	50	GLY	2.9
2	J	41	SER	2.9
1	F	191	PRO	2.8
2	H	49	LEU	2.6
1	B	158	ARG	2.6
1	D	133	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	196	LYS	2.5
1	E	122	GLU	2.4
1	E	195	ILE	2.4
1	E	193	PRO	2.3
2	G	41	SER	2.3
1	A	194	ASP	2.3
1	B	78	LYS	2.3
1	C	188	ASN	2.3
1	F	157	PHE	2.2
1	E	133	TYR	2.2
1	B	191	PRO	2.1
1	C	205	THR	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	ZN	D	400	1/1	0.92	0.09	-1.65	110,110,110,110	1
3	ZN	C	400	1/1	0.98	0.15	-2.21	84,84,84,84	1
3	ZN	E	400	1/1	0.79	0.07	-2.48	149,149,149,149	1
3	ZN	B	400	1/1	0.97	0.07	-3.34	119,119,119,119	1
3	ZN	A	400	1/1	0.90	0.07	-3.48	161,161,161,161	1
3	ZN	F	400	1/1	0.87	0.05	-	162,162,162,162	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.