



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 9, 2021 – 08:02 PM EST

PDB ID : 7N0Y  
Title : Rigidity of loop 1 contributes to equipotency of globular and ribbon isomers of alpha-conotoxin AusIA  
Authors : Ho, T.N.T.; Abraham, N.; Lewis, R.J.  
Deposited on : 2021-05-26  
Resolution : 2.58 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.

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

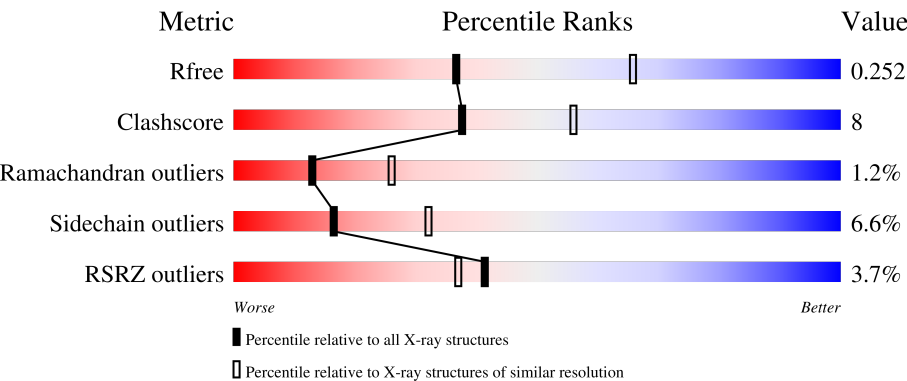
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:  
*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.58 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))
$R_{free}$	130704	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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	205	<div><div>4%</div><div>82%</div><div>15%</div><div>.</div></div>
1	B	205	<div><div>%</div><div>83%</div><div>13%</div><div>.</div></div>
1	C	205	<div><div>4%</div><div>80%</div><div>18%</div><div>.</div></div>
1	D	205	<div><div>2%</div><div>77%</div><div>17%</div><div>..</div></div>
1	E	205	<div><div>6%</div><div>79%</div><div>16%</div><div>...</div></div>

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Mol	Chain	Length	Quality of chain
2	G	16	 <p>12% 25% 50% 6% 6% 12%</p>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8245 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 Acetylcholine-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	205	Total	C	N	O	S	0	0	0
			1637	1023	280	329	5			
1	B	205	Total	C	N	O	S	0	0	0
			1637	1023	280	329	5			
1	C	204	Total	C	N	O	S	0	0	0
			1629	1019	278	327	5			
1	D	200	Total	C	N	O	S	0	0	0
			1602	1005	274	318	5			
1	E	202	Total	C	N	O	S	0	0	0
			1613	1010	276	322	5			

- Molecule 2 is a protein called Globular alpha-conotoxin AusIA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	14	Total	C	N	O	S	0	0	0
			106	60	26	16	4			

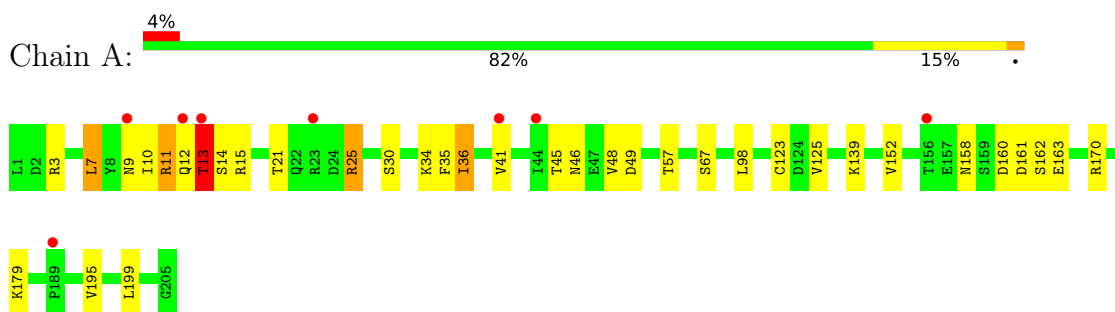
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	O	0	0
			2	2		
3	B	5	Total	O	0	0
			5	5		
3	C	4	Total	O	0	0
			4	4		
3	D	4	Total	O	0	0
			4	4		
3	E	6	Total	O	0	0
			6	6		

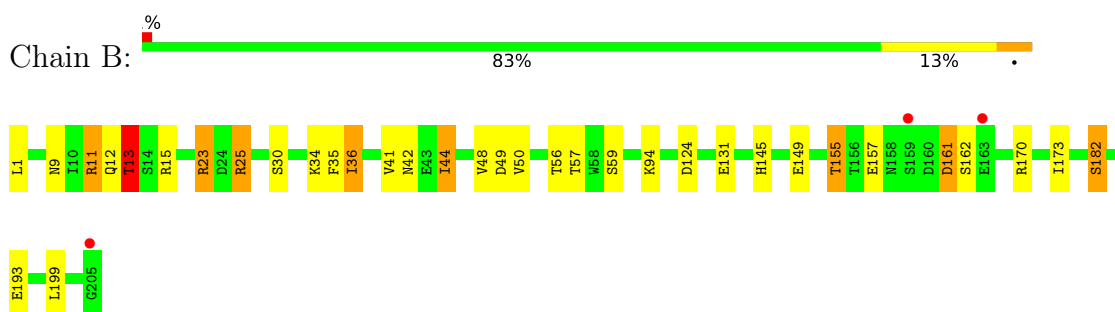
### 3 Residue-property plots [i](#)

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.

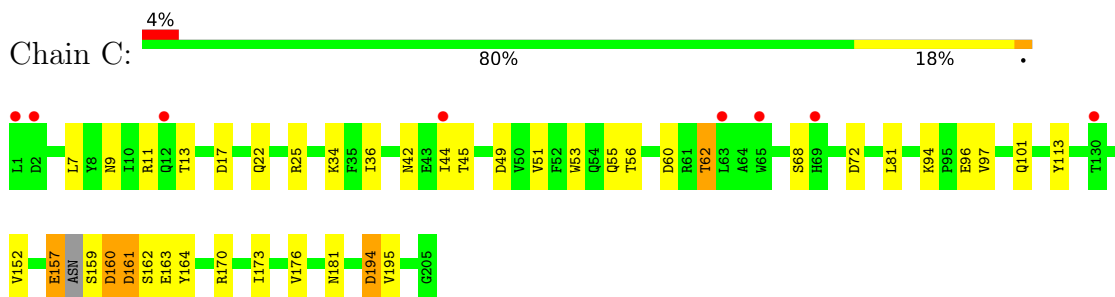
- Molecule 1: Acetylcholine-binding protein



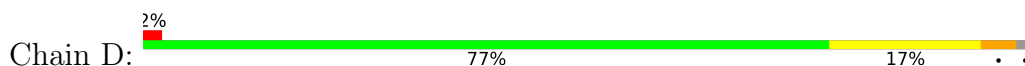
- Molecule 1: Acetylcholine-binding protein

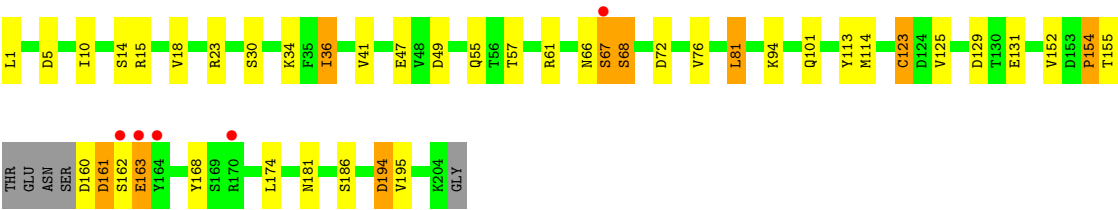


- Molecule 1: Acetylcholine-binding protein

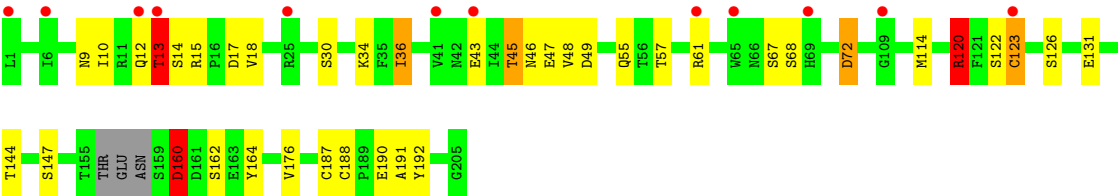
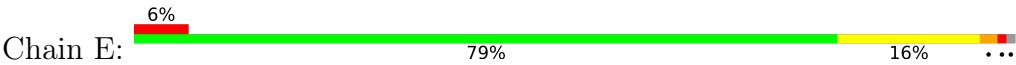


- Molecule 1: Acetylcholine-binding protein





● Molecule 1: Acetylcholine-binding protein



● Molecule 2: Globular alpha-conotoxin AusIA



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.30Å 73.30Å 347.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.93 – 2.58 46.88 – 2.58	Depositor EDS
% Data completeness (in resolution range)	99.8 (46.93-2.58) 99.9 (46.88-2.58)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.33 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.204 , 0.253 0.212 , 0.252	Depositor DCC
$R_{free}$ test set	2004 reflections (5.65%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	70.5	Xtriage
Anisotropy	0.204	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 50.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8245	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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	0.72	0/1673	0.91	0/2282
1	B	0.74	1/1673 (0.1%)	0.92	1/2282 (0.0%)
1	C	0.77	1/1664 (0.1%)	0.92	1/2268 (0.0%)
1	D	0.75	0/1637	0.97	1/2233 (0.0%)
1	E	0.76	0/1648	0.96	3/2246 (0.1%)
2	G	0.83	0/109	1.18	0/147
All	All	0.75	2/8404 (0.0%)	0.94	6/11458 (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	D	0	2
2	G	0	2
All	All	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	182	SER	CB-OG	-5.25	1.35	1.42
1	C	96	GLU	CD-OE1	5.01	1.31	1.25

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	194	ASP	CB-CG-OD1	8.34	125.80	118.30
1	C	194	ASP	CB-CG-OD1	7.82	125.34	118.30
1	E	72	ASP	CB-CG-OD1	-5.98	112.92	118.30
1	E	160	ASP	N-CA-C	5.76	126.56	111.00
1	B	1	LEU	CA-CB-CG	5.73	128.49	115.30
1	E	120	ARG	NE-CZ-NH1	5.17	122.88	120.30



There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	66	ASN	Peptide
1	D	67	SER	Peptide
2	G	2	CYS	Peptide
2	G	4	ALA	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	1637	0	1583	24	0
1	B	1637	0	1581	22	0
1	C	1629	0	1574	28	0
1	D	1602	0	1555	28	0
1	E	1613	0	1561	23	0
2	G	106	0	91	12	0
3	A	2	0	0	0	0
3	B	5	0	0	0	0
3	C	4	0	0	0	0
3	D	4	0	0	0	0
3	E	6	0	0	0	0
All	All	8245	0	7945	124	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 8.

All (124) 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:36:ILE:HD11	1:C:53:TRP:CD1	1.91	1.06
1:C:60:ASP:OD1	1:C:62:THR:HG23	1.59	1.03
1:A:41:VAL:CG1	1:A:125:VAL:HG21	1.91	1.00
2:G:2:CYS:N	2:G:13:HIS:HB2	1.81	0.94
1:D:55:GLN:HB2	1:D:114:MET:HE3	1.52	0.90
1:D:15:ARG:HB2	1:D:18:VAL:HG12	1.53	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:15:ARG:HB3	1:E:18:VAL:HG13	1.53	0.88
1:A:45:THR:O	1:A:45:THR:OG1	1.86	0.87
1:B:42:ASN:OD1	1:B:44:ILE:HG12	1.77	0.84
1:D:154:PRO:O	1:D:155:THR:HG23	1.81	0.80
1:D:76:VAL:HG21	1:D:81:LEU:HD11	1.64	0.77
1:A:41:VAL:HG13	1:A:125:VAL:HG21	1.63	0.77
1:B:11:ARG:NH2	1:E:17:ASP:OD2	2.18	0.76
1:A:9:ASN:O	1:A:12:GLN:O	2.04	0.76
1:D:15:ARG:CB	1:D:18:VAL:HG12	2.15	0.76
1:E:9:ASN:O	1:E:12:GLN:O	2.04	0.76
1:B:9:ASN:O	1:B:12:GLN:O	2.04	0.75
1:C:159:SER:N	1:C:176:VAL:O	2.21	0.74
1:E:15:ARG:CB	1:E:18:VAL:HG13	2.17	0.74
1:A:152:VAL:CG1	1:A:195:VAL:HG23	2.18	0.74
1:A:11:ARG:NH2	1:C:17:ASP:OD2	2.22	0.73
1:E:187:CYS:SG	2:G:2:CYS:HA	2.29	0.72
1:A:125:VAL:O	1:A:125:VAL:HG22	1.90	0.71
1:C:36:ILE:HD11	1:C:53:TRP:NE1	2.04	0.71
1:D:152:VAL:CG1	1:D:195:VAL:HG23	2.21	0.70
1:C:9:ASN:O	1:C:13:THR:HG22	1.91	0.70
1:B:35:PHE:CE1	1:B:199:LEU:HD22	2.29	0.68
1:A:41:VAL:CG1	1:A:125:VAL:CG2	2.71	0.67
1:A:152:VAL:HG12	1:A:195:VAL:HG23	1.75	0.67
1:D:15:ARG:HB2	1:D:18:VAL:CG1	2.26	0.66
1:C:7:LEU:C	1:C:7:LEU:HD13	2.16	0.66
1:B:11:ARG:CZ	1:B:11:ARG:HB2	2.26	0.65
1:C:157:GLU:OE2	1:C:160:ASP:O	2.15	0.65
1:D:152:VAL:HG12	1:D:195:VAL:HG23	1.78	0.65
1:E:15:ARG:HB3	1:E:18:VAL:CG1	2.28	0.62
1:C:36:ILE:HD12	1:C:51:VAL:HG12	1.81	0.61
1:C:97:VAL:O	1:D:94:LYS:HD3	2.03	0.58
1:E:160:ASP:HA	1:E:176:VAL:HB	1.85	0.58
1:D:41:VAL:O	1:D:47:GLU:O	2.22	0.58
1:C:152:VAL:CG1	1:C:195:VAL:HG23	2.35	0.57
1:C:181:ASN:HB3	1:C:194:ASP:OD1	2.05	0.57
1:C:152:VAL:HG12	1:C:195:VAL:HG23	1.87	0.56
1:D:162:SER:C	1:D:163:GLU:HG3	2.27	0.54
1:D:181:ASN:HB3	1:D:194:ASP:OD1	2.07	0.54
1:B:41:VAL:HG22	1:B:48:VAL:HG22	1.89	0.54
1:C:36:ILE:CD1	1:C:51:VAL:HG12	2.38	0.54
1:B:145:HIS:HD2	1:B:149:GLU:OE1	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:44:ILE:HG13	1:C:45:THR:HG23	1.91	0.53
1:C:101:GLN:OE1	1:C:113:TYR:OH	2.26	0.53
2:G:3:CYS:O	2:G:10:ARG:HG2	2.08	0.53
1:B:162:SER:HB3	1:B:173:ILE:CG1	2.38	0.53
1:E:34:LYS:HB3	1:E:164:TYR:HD2	1.74	0.52
2:G:14:PRO:O	2:G:15:CYS:HB3	2.10	0.52
1:B:155:THR:HG23	1:B:157:GLU:H	1.75	0.51
1:A:21:THR:OG1	1:A:25:ARG:O	2.27	0.51
1:D:161:ASP:OD1	1:D:161:ASP:C	2.49	0.51
2:G:2:CYS:N	2:G:13:HIS:CB	2.67	0.51
1:D:154:PRO:O	1:D:155:THR:CG2	2.57	0.50
2:G:3:CYS:HA	2:G:15:CYS:HB2	1.94	0.50
1:D:161:ASP:OD1	1:D:163:GLU:N	2.45	0.50
1:A:3:ARG:O	1:A:7:LEU:HD23	2.12	0.49
1:D:101:GLN:OE1	1:D:113:TYR:OH	2.30	0.49
1:C:34:LYS:HG2	1:C:164:TYR:CZ	2.48	0.49
1:B:35:PHE:CD1	1:B:199:LEU:HD22	2.47	0.48
1:B:161:ASP:N	1:B:161:ASP:OD1	2.47	0.48
1:E:144:THR:HG22	2:G:11:HIS:CD2	2.48	0.48
1:C:34:LYS:HG2	1:C:164:TYR:CE1	2.48	0.48
1:B:173:ILE:O	1:B:173:ILE:HG13	2.14	0.48
1:A:41:VAL:HG12	1:A:125:VAL:HG21	1.89	0.47
1:C:34:LYS:NZ	1:C:55:GLN:OE1	2.48	0.47
1:C:157:GLU:C	1:C:160:ASP:HB2	2.35	0.47
1:C:163:GLU:OE1	1:D:186:SER:HB2	2.15	0.47
1:D:55:GLN:CB	1:D:114:MET:HE3	2.37	0.47
1:E:15:ARG:CB	1:E:18:VAL:CG1	2.91	0.47
1:E:10:ILE:O	1:E:14:SER:HB3	2.16	0.46
1:E:48:VAL:HG12	1:E:123:CYS:SG	2.55	0.46
1:A:125:VAL:O	1:A:125:VAL:CG2	2.60	0.46
1:B:35:PHE:CZ	1:B:199:LEU:HD22	2.49	0.46
1:B:94:LYS:HE3	1:B:94:LYS:HB2	1.78	0.46
1:C:42:ASN:OD1	1:C:44:ILE:HG13	2.15	0.45
1:C:160:ASP:O	1:C:161:ASP:CB	2.64	0.45
1:A:12:GLN:C	1:A:13:THR:HG23	2.37	0.45
1:D:1:LEU:HD22	1:D:5:ASP:HB3	1.98	0.45
1:E:34:LYS:O	1:E:36:ILE:HD13	2.16	0.45
1:B:30:SER:OG	1:B:57:THR:OG1	2.33	0.45
1:A:41:VAL:HG22	1:A:48:VAL:HG23	1.98	0.45
1:D:34:LYS:O	1:D:36:ILE:HD13	2.16	0.45
2:G:4:ALA:HA	2:G:10:ARG:NH1	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:14:PRO:O	2:G:15:CYS:CB	2.64	0.45
1:A:10:ILE:O	1:A:14:SER:HB3	2.17	0.45
1:A:170:ARG:HD3	1:C:44:ILE:O	2.16	0.44
1:E:12:GLN:C	1:E:13:THR:HG23	2.37	0.44
1:D:129:ASP:OD1	1:D:129:ASP:N	2.45	0.44
1:B:182:SER:HB3	1:B:193:GLU:HG2	1.98	0.44
1:D:67:SER:O	1:D:68:SER:HB3	2.18	0.44
1:D:10:ILE:O	1:D:14:SER:HB3	2.18	0.44
1:C:181:ASN:CB	1:C:194:ASP:OD1	2.66	0.44
1:A:34:LYS:O	1:A:36:ILE:HD13	2.18	0.44
1:B:34:LYS:O	1:B:36:ILE:HD13	2.18	0.44
1:B:157:GLU:HG3	2:G:15:CYS:SG	2.58	0.44
1:A:160:ASP:HB3	1:A:163:GLU:HB2	2.00	0.43
1:B:12:GLN:C	1:B:13:THR:HG23	2.38	0.43
1:B:170:ARG:NH2	1:E:43:GLU:O	2.49	0.43
1:D:174:LEU:N	1:D:174:LEU:CD1	2.81	0.43
1:E:45:THR:O	1:E:47:GLU:N	2.51	0.43
1:E:30:SER:OG	1:E:57:THR:OG1	2.35	0.43
1:A:41:VAL:HG11	1:A:125:VAL:CG2	2.48	0.43
1:B:124:ASP:HB2	1:D:168:TYR:CE1	2.54	0.43
1:E:188:CYS:SG	2:G:13:HIS:CD2	3.13	0.42
1:C:22:GLN:O	1:C:25:ARG:HG3	2.19	0.42
1:E:67:SER:O	1:E:67:SER:OG	2.34	0.42
1:A:35:PHE:CE2	1:A:199:LEU:HD22	2.53	0.42
1:C:162:SER:HB2	1:C:173:ILE:HD12	2.01	0.42
1:B:23:ARG:NH2	1:B:25:ARG:HG3	2.35	0.42
1:E:147:SER:OG	1:E:191:ALA:O	2.32	0.42
1:E:47:GLU:HB3	1:E:120:ARG:NH1	2.35	0.41
1:E:55:GLN:HB2	1:E:114:MET:CE	2.50	0.41
1:A:30:SER:OG	1:A:57:THR:OG1	2.36	0.41
1:D:76:VAL:HG21	1:D:81:LEU:CD1	2.41	0.41
1:D:30:SER:OG	1:D:57:THR:OG1	2.38	0.41
1:E:192:TYR:HB3	2:G:6:ASN:ND2	2.36	0.41
1:A:98:LEU:HD22	1:C:94:LYS:N	2.36	0.40
1:D:123:CYS:SG	1:D:125:VAL:HG12	2.61	0.40
1:A:41:VAL:HG13	1:A:125:VAL:CG2	2.41	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	203/205 (99%)	196 (97%)	4 (2%)	3 (2%)	10	20
1	B	203/205 (99%)	196 (97%)	5 (2%)	2 (1%)	15	31
1	C	200/205 (98%)	195 (98%)	4 (2%)	1 (0%)	29	50
1	D	196/205 (96%)	188 (96%)	6 (3%)	2 (1%)	15	31
1	E	198/205 (97%)	192 (97%)	3 (2%)	3 (2%)	10	20
2	G	12/16 (75%)	9 (75%)	2 (17%)	1 (8%)	1	0
All	All	1012/1041 (97%)	976 (96%)	24 (2%)	12 (1%)	13	26

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	B	161	ASP
1	C	161	ASP
1	D	154	PRO
1	E	46	ASN
1	E	160	ASP
1	A	161	ASP
2	G	5	ARG
1	D	68	SER
1	A	13	THR
1	E	13	THR
1	B	13	THR

### 5.3.2 Protein sidechains [i](#)

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	191/191 (100%)	178 (93%)	13 (7%)	16	30
1	B	191/191 (100%)	178 (93%)	13 (7%)	16	30
1	C	190/191 (100%)	180 (95%)	10 (5%)	22	43
1	D	187/191 (98%)	176 (94%)	11 (6%)	19	37
1	E	188/191 (98%)	173 (92%)	15 (8%)	12	22
2	G	12/14 (86%)	11 (92%)	1 (8%)	11	21
All	All	959/969 (99%)	896 (93%)	63 (7%)	16	32

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	11	ARG
1	A	13	THR
1	A	15	ARG
1	A	25	ARG
1	A	36	ILE
1	A	49	ASP
1	A	67	SER
1	A	123	CYS
1	A	139	LYS
1	A	158	ASN
1	A	162	SER
1	A	179	LYS
1	B	11	ARG
1	B	13	THR
1	B	15	ARG
1	B	23	ARG
1	B	25	ARG
1	B	36	ILE
1	B	44	ILE
1	B	49	ASP
1	B	50	VAL
1	B	56	THR
1	B	59	SER
1	B	131	GLU
1	B	155	THR
1	C	11	ARG
1	C	49	ASP

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Mol	Chain	Res	Type
1	C	56	THR
1	C	62	THR
1	C	68	SER
1	C	72	ASP
1	C	81	LEU
1	C	157	GLU
1	C	160	ASP
1	C	170	ARG
1	D	23	ARG
1	D	36	ILE
1	D	49	ASP
1	D	61	ARG
1	D	72	ASP
1	D	81	LEU
1	D	123	CYS
1	D	131	GLU
1	D	160	ASP
1	D	161	ASP
1	D	163	GLU
1	E	13	THR
1	E	36	ILE
1	E	45	THR
1	E	49	ASP
1	E	61	ARG
1	E	68	SER
1	E	72	ASP
1	E	120	ARG
1	E	122	SER
1	E	123	CYS
1	E	126	SER
1	E	131	GLU
1	E	160	ASP
1	E	162	SER
1	E	190	GLU
2	G	2	CYS

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

Mol	Chain	Res	Type
1	A	12	GLN
1	B	145	HIS
1	E	55	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides 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	205/205 (100%)	0.37	8 (3%)	39 34	55, 80, 122, 137	0
1	B	205/205 (100%)	0.15	3 (1%)	73 71	55, 81, 120, 147	0
1	C	204/205 (99%)	0.24	8 (3%)	39 34	51, 77, 106, 133	0
1	D	200/205 (97%)	0.12	5 (2%)	57 53	52, 74, 111, 142	0
1	E	202/205 (98%)	0.36	12 (5%)	22 19	51, 77, 109, 134	0
2	G	14/16 (87%)	0.60	2 (14%)	2 2	70, 80, 109, 114	0
All	All	1030/1041 (98%)	0.25	38 (3%)	41 37	51, 78, 117, 147	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	43	GLU	4.0
1	E	69	HIS	3.9
1	E	65	TRP	3.8
1	E	1	LEU	3.8
1	A	41	VAL	3.7
1	C	63	LEU	3.6
1	C	2	ASP	3.4
2	G	2	CYS	3.4
1	E	123	CYS	3.3
1	A	13	THR	3.3
1	A	156	THR	3.2
1	E	61	ARG	3.0
1	C	1	LEU	2.9
1	E	41	VAL	2.9
1	D	162	SER	2.9
1	C	44	ILE	2.9
1	D	164	TYR	2.6
1	D	67	SER	2.5
1	E	13	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	44	ILE	2.5
1	C	12	GLN	2.5
1	A	23	ARG	2.5
1	B	205	GLY	2.4
1	C	69	HIS	2.4
1	B	159	SER	2.4
1	A	189	PRO	2.3
1	E	25	ARG	2.3
2	G	15	CYS	2.3
1	A	9	ASN	2.2
1	E	12	GLN	2.2
1	D	163	GLU	2.2
1	C	130	THR	2.2
1	E	109	GLY	2.1
1	D	170	ARG	2.1
1	B	163	GLU	2.1
1	C	65	TRP	2.0
1	E	6	ILE	2.0
1	A	12	GLN	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 monosaccharides 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.