



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 24, 2023 – 02:39 AM EDT

PDB ID : 5T90  
Title : Structural mechanisms for alpha-conotoxin selectivity at the human alpha3beta4 nicotinic acetylcholine receptor  
Authors : Abraham, N.; Healy, M.; Ragnarsson, L.; Brust, A.; Alewood, P.; Lewis, R.  
Deposited on : 2016-09-08  
Resolution : 2.80 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 : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

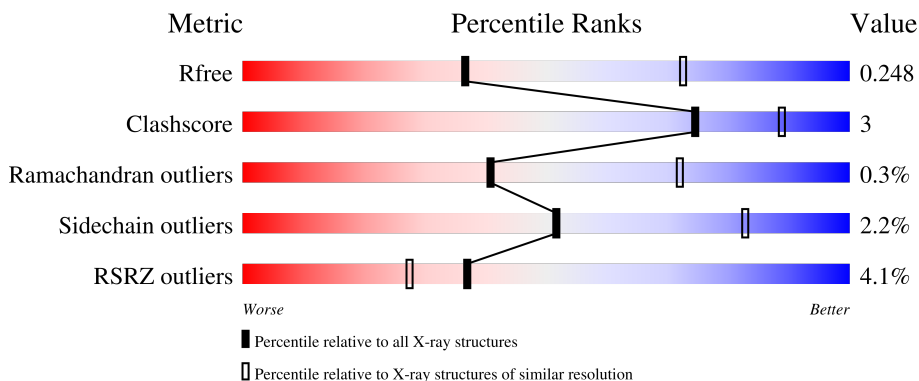
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	210	85% 12% .
1	B	210	3% 88% 6% 6%
1	C	210	2% 85% 11% .
1	D	210	10% 84% 10% 5%
1	E	210	2% 89% 9% .

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Mol	Chain	Length	Quality of chain
2	F	18	 78% 22%
2	G	18	 78% 17% 6%
2	H	18	 39% 83% 17%
2	I	18	 11% 67% 17% 6% 11%
2	J	18	 67% 17% 6% 11%

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8712 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
			Total	C	N	O	S			
1	A	205	Total 1636	C 1023	N 280	O 328	S 5	0	0	0
1	B	197	Total 1576	C 989	N 271	O 311	S 5	0	0	0
1	C	203	Total 1624	C 1017	N 278	O 324	S 5	0	0	0
1	D	200	Total 1604	C 1004	N 275	O 320	S 5	0	0	0
1	E	204	Total 1629	C 1019	N 279	O 326	S 5	0	0	0

- Molecule 2 is a protein called LsIA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	F	18	Total 118	C 66	N 25	O 23	S 4	0	0	1
2	G	18	Total 118	C 66	N 25	O 23	S 4	0	0	1
2	H	18	Total 118	C 66	N 25	O 23	S 4	0	0	1
2	I	16	Total 102	C 56	N 22	O 20	S 4	0	0	1
2	J	16	Total 102	C 56	N 22	O 20	S 4	0	0	1

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	25	Total 25	O 25	0	0
3	B	12	Total 12	O 12	0	0

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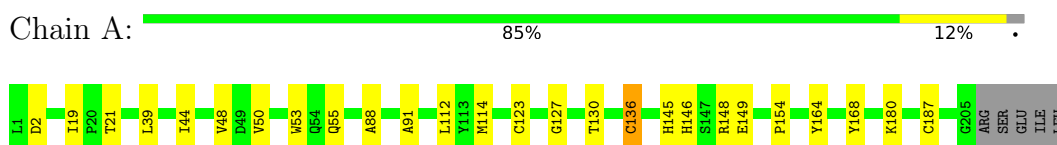
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
3	C	15	Total 15	O 15	0	0
3	D	14	Total 14	O 14	0	0
3	E	15	Total 15	O 15	0	0
3	F	1	Total 1	O 1	0	0
3	G	1	Total 1	O 1	0	0
3	H	1	Total 1	O 1	0	0
3	J	1	Total 1	O 1	0	0

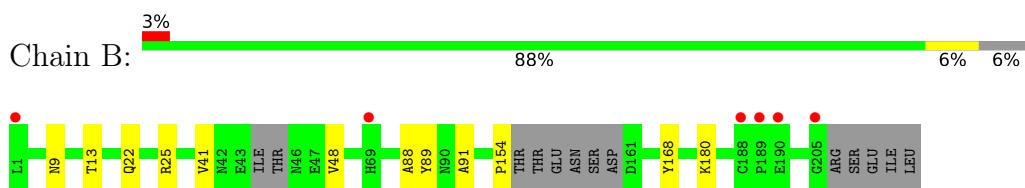
### 3 Residue-property plots [i](#)

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.

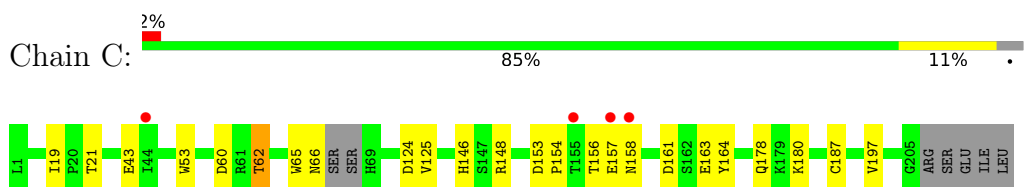
- Molecule 1: Acetylcholine-binding protein



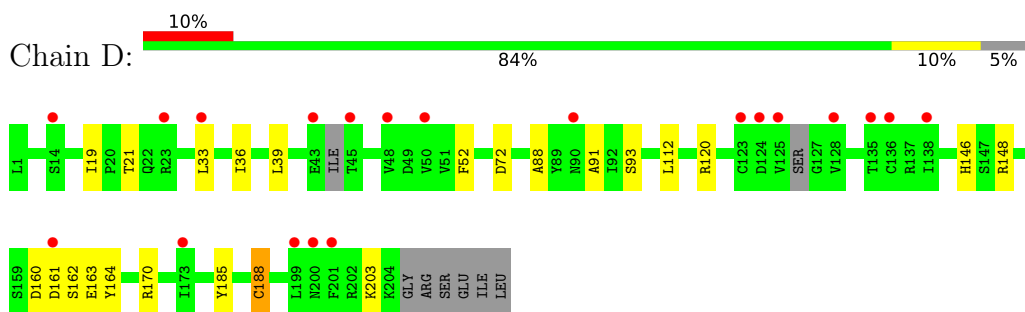
- Molecule 1: Acetylcholine-binding protein



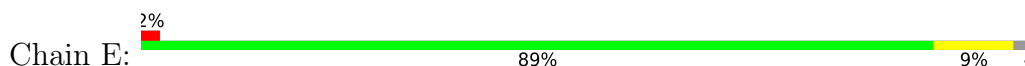
- Molecule 1: Acetylcholine-binding protein

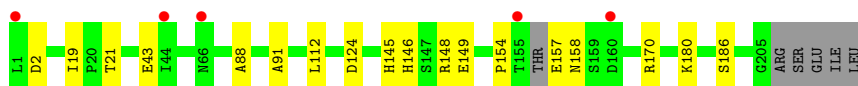


- Molecule 1: Acetylcholine-binding protein

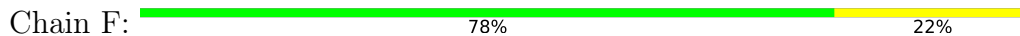


- Molecule 1: Acetylcholine-binding protein

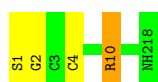




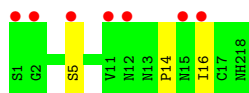
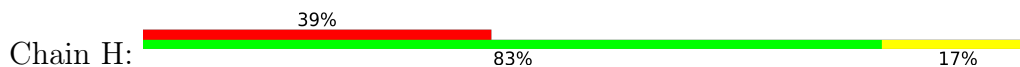
● Molecule 2: LsIA



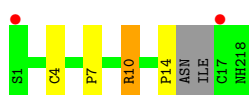
● Molecule 2: LsIA



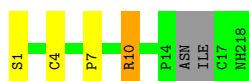
● Molecule 2: LsIA



● Molecule 2: LsIA



● Molecule 2: LsIA



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	115.94Å 124.50Å 154.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.44 – 2.80 48.44 – 2.80	Depositor EDS
% Data completeness (in resolution range)	96.7 (48.44-2.80) 96.7 (48.44-2.80)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.38 (at 2.81Å)	Xtrriage
Refinement program	BUSTER 2.10.2	Depositor
R, $R_{free}$	0.218 , 0.245 0.224 , 0.248	Depositor DCC
$R_{free}$ test set	1354 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	51.7	Xtrriage
Anisotropy	1.001	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 53.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8712	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	70.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NH2

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	0.37	0/1672	0.62	0/2282
1	B	0.36	0/1610	0.59	0/2193
1	C	0.38	0/1659	0.62	0/2263
1	D	0.37	0/1637	0.60	0/2229
1	E	0.37	0/1664	0.61	0/2269
2	F	0.41	0/118	0.56	0/160
2	G	0.38	0/118	0.57	0/160
2	H	0.38	0/118	0.52	0/160
2	I	0.36	0/101	0.52	0/135
2	J	0.38	0/101	0.54	0/135
All	All	0.37	0/8798	0.60	0/11986

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	1636	0	1581	13	0
1	B	1576	0	1526	7	0
1	C	1624	0	1570	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1604	0	1545	10	0
1	E	1629	0	1573	10	0
2	F	118	0	107	3	0
2	G	118	0	107	2	0
2	H	118	0	107	2	0
2	I	102	0	89	3	0
2	J	102	0	89	2	0
3	A	25	0	0	0	0
3	B	12	0	0	0	0
3	C	15	0	0	0	0
3	D	14	0	0	0	0
3	E	15	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
3	J	1	0	0	0	0
All	All	8712	0	8294	54	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 3.

All (54) 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:157:GLU:H	1:C:158:ASN:HA	1.31	0.92
1:C:157:GLU:N	1:C:158:ASN:HA	1.99	0.76
1:D:185:TYR:HB2	1:D:188:CYS:SG	2.35	0.67
1:A:127:GLY:O	1:A:130:THR:HG22	1.95	0.67
1:E:112:LEU:HD21	2:H:14:PRO:HB3	1.77	0.66
1:A:44:ILE:O	1:E:170:ARG:HD2	1.99	0.63
1:A:187:CYS:HB2	2:H:16:ILE:HG12	1.82	0.59
1:E:157:GLU:HB2	1:E:158:ASN:HA	1.83	0.59
1:C:178:GLN:HG2	1:C:197:VAL:HG22	1.86	0.57
1:A:88:ALA:HB3	1:A:91:ALA:HB2	1.90	0.53
1:B:22:GLN:O	1:B:25:ARG:HG2	2.10	0.52
1:B:41:VAL:HG22	1:B:48:VAL:HG12	1.91	0.52
1:A:112:LEU:HD21	2:F:14:PRO:HB3	1.93	0.51
1:A:48:VAL:HG11	1:A:136:CYS:SG	2.50	0.51
1:C:187:CYS:HB2	2:F:16:ILE:HG12	1.92	0.51
1:B:168:TYR:CD1	1:E:124:ASP:HB2	2.46	0.51
2:I:4:CYS:HA	2:I:10:ARG:HG3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:4:CYS:HA	2:J:10:ARG:HG3	1.94	0.49
1:C:43:GLU:HA	1:C:125:VAL:HG23	1.95	0.48
1:D:112:LEU:HD21	2:I:14:PRO:HB3	1.95	0.48
1:D:170:ARG:HH21	1:D:203:LYS:HD2	1.80	0.47
2:G:4:CYS:HA	2:G:10:ARG:HG3	1.97	0.46
1:E:19:ILE:HG12	1:E:21:THR:HG23	1.98	0.46
1:E:186:SER:OG	2:G:1:SER:HB2	2.15	0.46
1:A:19:ILE:HG12	1:A:21:THR:HG23	1.98	0.46
1:A:168:TYR:CD1	1:C:124:ASP:HB3	2.51	0.46
1:C:154:PRO:HD3	1:C:180:LYS:HB2	1.99	0.45
1:C:19:ILE:HG12	1:C:21:THR:HG23	1.99	0.45
1:A:53:TRP:CH2	2:F:7:PRO:HG3	2.53	0.43
1:D:19:ILE:HG12	1:D:21:THR:HG23	2.00	0.43
1:D:33:LEU:HD22	1:D:52:PHE:CD1	2.54	0.43
1:D:88:ALA:HB3	1:D:91:ALA:HB2	2.01	0.43
1:D:93:SER:HB2	1:D:120:ARG:HB2	2.01	0.42
1:E:145:HIS:HB3	1:E:149:GLU:HB2	2.01	0.42
1:D:161:ASP:HB3	1:D:162:SER:HA	2.01	0.42
1:A:55:GLN:HA	1:A:114:MET:HG3	2.02	0.42
1:E:154:PRO:HD3	1:E:180:LYS:HB2	2.02	0.42
1:A:154:PRO:HD3	1:A:180:LYS:HB2	2.02	0.42
1:B:154:PRO:HD3	1:B:180:LYS:HB2	2.02	0.41
1:C:65:TRP:O	1:C:66:ASN:HB2	2.20	0.41
1:C:146:HIS:CE1	1:C:148:ARG:HB2	2.55	0.41
1:E:146:HIS:CE1	1:E:148:ARG:HB2	2.55	0.41
1:D:146:HIS:CE1	1:D:148:ARG:HB2	2.55	0.41
1:B:88:ALA:HB3	1:B:91:ALA:HB2	2.03	0.41
1:A:146:HIS:CE1	1:A:148:ARG:HB2	2.55	0.41
1:B:9:ASN:O	1:B:13:THR:HG22	2.21	0.41
1:B:89:TYR:CE1	2:I:7:PRO:HG2	2.56	0.41
1:C:53:TRP:CH2	2:J:7:PRO:HG3	2.56	0.41
1:C:156:THR:HB	1:C:158:ASN:HA	2.03	0.41
1:C:60:ASP:OD1	1:C:62:THR:HG22	2.21	0.40
1:C:157:GLU:N	1:C:158:ASN:CA	2.79	0.40
1:D:36:ILE:HD13	1:D:164:TYR:HB3	2.04	0.40
1:A:145:HIS:HB3	1:A:149:GLU:HB2	2.03	0.40
1:E:88:ALA:HB3	1:E:91:ALA:HB2	2.03	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	203/210 (97%)	198 (98%)	5 (2%)	0	100	100
1	B	191/210 (91%)	189 (99%)	2 (1%)	0	100	100
1	C	199/210 (95%)	188 (94%)	10 (5%)	1 (0%)	29	61
1	D	192/210 (91%)	182 (95%)	9 (5%)	1 (0%)	29	61
1	E	200/210 (95%)	195 (98%)	5 (2%)	0	100	100
2	F	16/18 (89%)	15 (94%)	1 (6%)	0	100	100
2	G	16/18 (89%)	15 (94%)	0	1 (6%)	1	3
2	H	16/18 (89%)	16 (100%)	0	0	100	100
2	I	12/18 (67%)	12 (100%)	0	0	100	100
2	J	12/18 (67%)	12 (100%)	0	0	100	100
All	All	1057/1140 (93%)	1022 (97%)	32 (3%)	3 (0%)	41	72

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	161	ASP
1	D	160	ASP
2	G	2	GLY

### 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	191/196 (97%)	185 (97%)	6 (3%)	40	74
1	B	183/196 (93%)	183 (100%)	0	100	100
1	C	189/196 (96%)	185 (98%)	4 (2%)	53	84
1	D	187/196 (95%)	183 (98%)	4 (2%)	53	84
1	E	190/196 (97%)	188 (99%)	2 (1%)	73	92
2	F	15/15 (100%)	14 (93%)	1 (7%)	16	43
2	G	15/15 (100%)	14 (93%)	1 (7%)	16	43
2	H	15/15 (100%)	14 (93%)	1 (7%)	16	43
2	I	13/15 (87%)	12 (92%)	1 (8%)	13	35
2	J	13/15 (87%)	11 (85%)	2 (15%)	2	8
All	All	1011/1055 (96%)	989 (98%)	22 (2%)	52	83

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ASP
1	A	39	LEU
1	A	50	VAL
1	A	123	CYS
1	A	136	CYS
1	A	164	TYR
1	C	62	THR
1	C	153	ASP
1	C	163	GLU
1	C	164	TYR
1	D	39	LEU
1	D	72	ASP
1	D	163	GLU
1	D	188	CYS
1	E	2	ASP
1	E	43	GLU
2	F	10	ARG
2	G	10	ARG
2	H	5	SER
2	I	10	ARG
2	J	1	SER
2	J	10	ARG

Sometimes 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 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 [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	205/210 (97%)	-0.07	0 <span style="border: 1px solid blue; padding: 2px;">100</span> <span style="border: 1px solid blue; padding: 2px;">100</span>	33, 53, 83, 103	0
1	B	197/210 (93%)	0.26	6 (3%) <span style="border: 1px solid red; padding: 2px;">50</span> <span style="border: 1px solid red; padding: 2px;">40</span>	45, 74, 117, 145	0
1	C	203/210 (96%)	0.11	4 (1%) <span style="border: 1px solid blue; padding: 2px;">65</span> <span style="border: 1px solid blue; padding: 2px;">56</span>	34, 62, 111, 128	0
1	D	200/210 (95%)	0.59	21 (10%) <span style="border: 1px solid red; padding: 2px;">6</span> <span style="border: 1px solid red; padding: 2px;">3</span>	48, 76, 119, 135	0
1	E	204/210 (97%)	0.16	5 (2%) <span style="border: 1px solid blue; padding: 2px;">57</span> <span style="border: 1px solid blue; padding: 2px;">47</span>	38, 65, 104, 118	0
2	F	17/18 (94%)	-0.37	0 <span style="border: 1px solid blue; padding: 2px;">100</span> <span style="border: 1px solid blue; padding: 2px;">100</span>	43, 46, 58, 71	0
2	G	17/18 (94%)	0.57	0 <span style="border: 1px solid blue; padding: 2px;">100</span> <span style="border: 1px solid blue; padding: 2px;">100</span>	68, 76, 86, 92	0
2	H	17/18 (94%)	1.52	7 (41%) <span style="border: 1px solid red; padding: 2px;">0</span> <span style="border: 1px solid red; padding: 2px;">0</span>	60, 72, 87, 93	0
2	I	15/18 (83%)	0.85	2 (13%) <span style="border: 1px solid red; padding: 2px;">3</span> <span style="border: 1px solid red; padding: 2px;">2</span>	82, 88, 97, 99	0
2	J	15/18 (83%)	0.25	0 <span style="border: 1px solid blue; padding: 2px;">100</span> <span style="border: 1px solid blue; padding: 2px;">100</span>	78, 84, 90, 94	0
All	All	1090/1140 (95%)	0.24	45 (4%) <span style="border: 1px solid red; padding: 2px;">37</span> <span style="border: 1px solid red; padding: 2px;">27</span>	33, 68, 109, 145	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	123	CYS	4.8
1	E	160	ASP	3.9
1	D	136	CYS	3.8
1	D	200	ASN	3.7
1	D	124	ASP	3.7
1	D	128	VAL	3.6
2	H	1	SER	3.6
2	H	15	ASN	3.4
1	D	199	LEU	3.3
1	D	201	PHE	3.3
2	H	11	VAL	3.2
1	B	190	GLU	3.1
1	B	188	CYS	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	14	SER	3.0
1	E	1	LEU	2.8
1	C	158	ASN	2.7
1	D	23	ARG	2.6
1	C	157	GLU	2.5
1	B	1	LEU	2.5
1	D	125	VAL	2.4
1	C	155	THR	2.4
1	D	158	ASN	2.4
1	D	161	ASP	2.4
2	H	5	SER	2.4
2	I	17	CYS	2.4
1	C	44	ILE	2.3
1	B	189	PRO	2.3
1	D	173	ILE	2.3
2	H	2	GLY	2.3
1	D	90	ASN	2.2
1	E	155	THR	2.2
1	B	69	HIS	2.2
1	B	205	GLY	2.2
1	E	44	ILE	2.1
1	D	50	VAL	2.1
2	H	16	ILE	2.1
1	D	48	VAL	2.1
2	I	1	SER	2.1
1	D	138	ILE	2.1
1	E	66	ASN	2.1
1	D	135	THR	2.0
1	D	33	LEU	2.0
2	H	12	ASN	2.0
1	D	43	GLU	2.0
1	D	45	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 monosaccharides in this entry.



## 6.4 Ligands

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.