



# Full wwPDB X-ray Structure Validation Report ⓘ

May 20, 2026 – 10:15 am BST

PDB ID : 9TQ6 / pdb\_00009tq6  
Title : Antibody-antigen complex  
Authors : Alzari, P.M.; Vulliez-Le Normand, B.; Martinez, M.  
Deposited on : 2025-12-19  
Resolution : 2.40 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>.

---

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

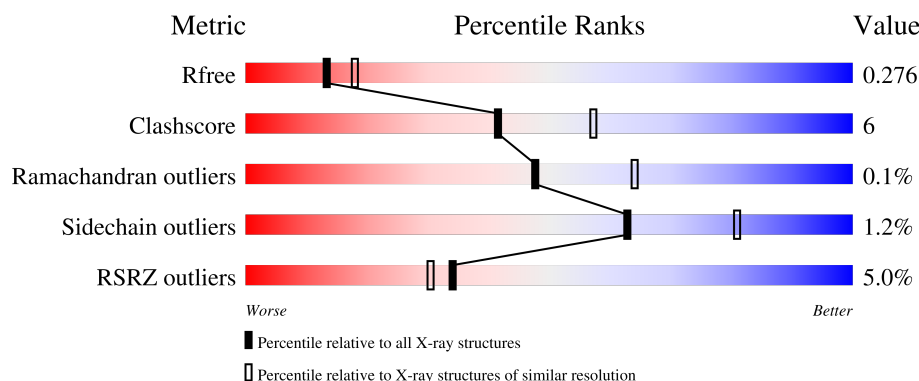
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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))
$R_{free}$	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	239	<div> <div>8%</div> <div> <div></div> <div>74%</div> <div>13%</div> <div>13%</div> </div> </div>
1	H	239	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>11%</div> <div>13%</div> </div> </div>
2	B	239	<div> <div>8%</div> <div> <div></div> <div>76%</div> <div>15%</div> <div>9%</div> </div> </div>
2	L	239	<div> <div></div> <div> <div></div> <div>74%</div> <div>18%</div> <div>8%</div> </div> </div>
3	P	11	<div> <div></div> <div> <div></div> <div>55%</div> <div>9%</div> <div>9%</div> <div>27%</div> </div> </div>

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
3	Q	11	 A horizontal bar chart showing the quality of chain Q. The bar is divided into four segments: a green segment (55%), a yellow segment (9%), an orange segment (9%), and a grey segment (27%).

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6765 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 Antibody Fab fragment heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1585	1012	254	314	5			
1	H	208	Total	C	N	O	S	0	0	0
			1585	1012	254	314	5			

- Molecule 2 is a protein called Antibody Fab fragment light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	217	Total	C	N	O	S	0	0	0
			1690	1056	283	343	8			
2	L	219	Total	C	N	O	S	0	0	0
			1707	1065	287	347	8			

- Molecule 3 is a protein called Beta-nerve growth factor.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	8	Total	C	N	O	0	0	0
			76	50	11	15			
3	Q	8	Total	C	N	O	0	0	0
			76	50	11	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	47	ACE	-	acetylation	UNP P01138
Q	47	ACE	-	acetylation	UNP P01138

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	O	0	0
			2	2		

*Continued on next page...*

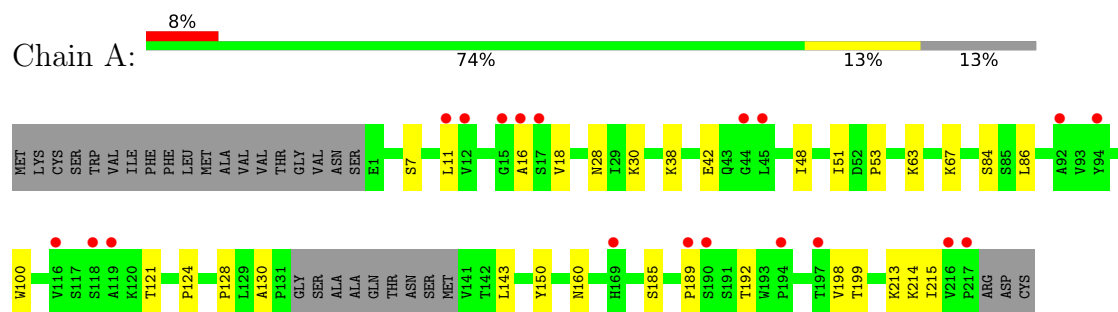
*Continued from previous page...*

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	O 1	0	0
4	H	22	Total 22	O 22	0	0
4	L	17	Total 17	O 17	0	0
4	P	3	Total 3	O 3	0	0
4	Q	1	Total 1	O 1	0	0

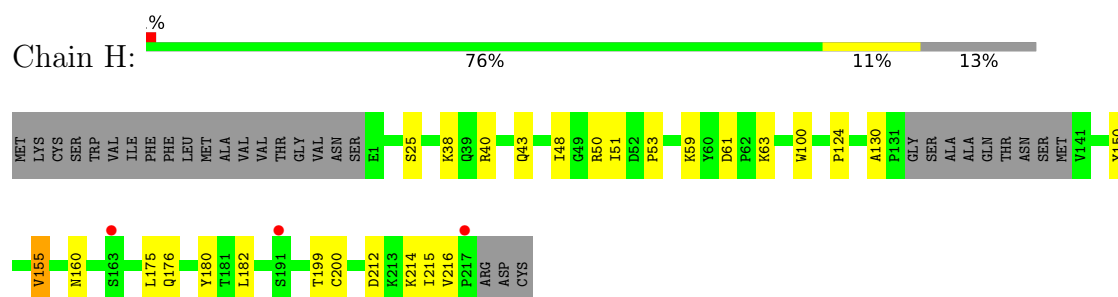
### 3 Residue-property plots [i](#)

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.

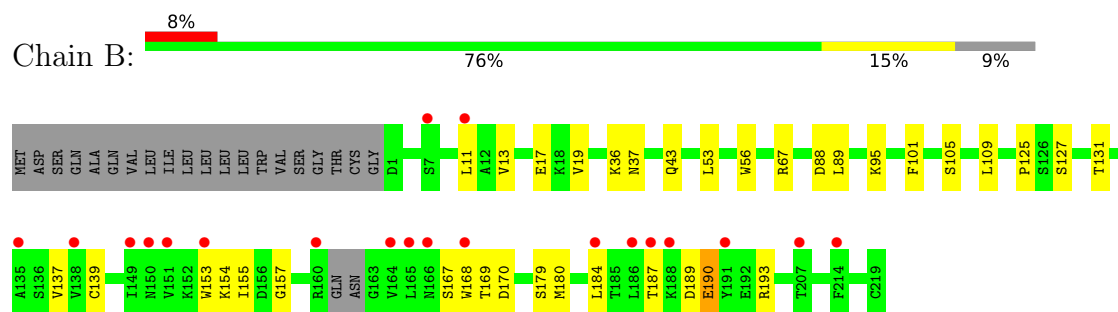
- Molecule 1: Antibody Fab fragment heavy chain



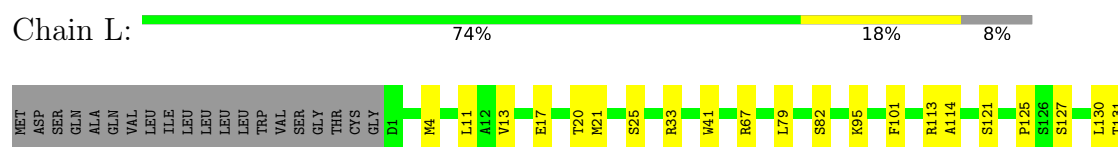
- Molecule 1: Antibody Fab fragment heavy chain



- Molecule 2: Antibody Fab fragment light chain

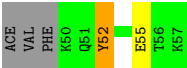


- Molecule 2: Antibody Fab fragment light chain





● Molecule 3: Beta-nerve growth factor



● Molecule 3: Beta-nerve growth factor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.56Å 91.28Å 142.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.64 – 2.40 45.64 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.64-2.40) 99.7 (45.64-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.86 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, $R_{free}$	0.235 , 0.276 0.235 , 0.276	Depositor DCC
$R_{free}$ test set	1876 reflections (3.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	57.1	Xtriage
Anisotropy	0.669	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 49.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6765	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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: NIY

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	0.31	0/1631	0.56	0/2239
1	H	0.45	0/1631	0.64	0/2239
2	B	0.30	0/1726	0.51	0/2336
2	L	0.43	0/1744	0.62	0/2362
3	P	0.46	0/61	0.74	0/78
3	Q	0.46	0/61	0.49	0/78
All	All	0.38	0/6854	0.59	0/9332

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

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	1585	0	1537	20	0
1	H	1585	0	1537	17	0
2	B	1690	0	1634	23	0
2	L	1707	0	1649	23	0
3	P	76	0	59	2	0
3	Q	76	0	60	2	0
4	A	2	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	1	0	0	0	0
4	H	22	0	0	1	0
4	L	17	0	0	1	0
4	P	3	0	0	0	0
4	Q	1	0	0	0	0
All	All	6765	0	6476	84	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 6.

All (84) 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:38:LYS:HB2	1:A:48:ILE:HD11	1.65	0.79
2:B:13:VAL:HG11	2:B:19:VAL:HG21	1.74	0.69
1:H:50:ARG:NH2	1:H:59:LYS:HE2	2.09	0.66
2:B:168:TRP:CD2	2:B:180:MET:HG3	2.33	0.64
1:H:155:VAL:HG12	1:H:182:LEU:HD21	1.82	0.61
3:Q:52:NIY:HB3	3:Q:55:GLU:HG3	1.82	0.61
2:L:13:VAL:HG22	2:L:17:GLU:HB2	1.83	0.60
2:L:125:PRO:HD3	2:L:137:VAL:HG22	1.83	0.60
1:H:40:ARG:HB2	1:H:43:GLN:HG3	1.85	0.58
3:P:52:NIY:HB3	3:P:55:GLU:HG3	1.85	0.58
1:H:130:ALA:HB2	1:H:215:ILE:HG22	1.85	0.58
1:A:143:LEU:CD1	1:A:198:VAL:HG21	2.35	0.57
1:H:100:TRP:HE1	2:L:95:LYS:NZ	2.03	0.56
2:L:33:ARG:HB2	4:L:312:HOH:O	2.06	0.55
1:A:130:ALA:HB2	1:A:215:ILE:HG22	1.88	0.54
2:L:150:ASN:HB2	2:L:202:THR:HB	1.89	0.54
1:H:100:TRP:HE1	2:L:95:LYS:HZ3	1.54	0.53
2:B:155:ILE:HD11	2:B:184:LEU:HD21	1.91	0.52
1:H:124:PRO:HB3	1:H:150:TYR:HB3	1.91	0.52
2:B:168:TRP:CE2	2:B:180:MET:HG3	2.45	0.52
3:P:52:NIY:HD2	3:P:55:GLU:OE2	2.10	0.52
1:H:175:LEU:HD13	1:H:180:TYR:CZ	2.46	0.51
2:B:169:THR:HG22	2:B:179:SER:H	1.76	0.51
2:L:113:ARG:HD3	2:L:114:ALA:O	2.11	0.51
3:Q:52:NIY:HD2	3:Q:55:GLU:OE1	2.11	0.51
1:H:199:THR:HA	1:H:214:LYS:HA	1.93	0.50
2:L:137:VAL:HG12	2:L:153:TRP:CH2	2.47	0.50
1:A:124:PRO:HB3	1:A:150:TYR:HB3	1.94	0.50

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:PRO:HD3	1:A:213:LYS:HD3	1.94	0.50
2:B:125:PRO:HD3	2:B:137:VAL:HG22	1.93	0.49
1:H:214:LYS:HE3	1:H:216:VAL:HG12	1.95	0.49
2:B:139:CYS:HB2	2:B:153:TRP:CH2	2.48	0.49
2:L:127:SER:O	2:L:131:THR:HG23	2.13	0.48
2:B:37:ASN:O	2:B:56:TRP:HA	2.14	0.48
2:B:11:LEU:HD12	2:B:11:LEU:HA	1.75	0.48
2:B:127:SER:O	2:B:131:THR:HG23	2.13	0.48
2:L:95:LYS:HE3	2:L:101:PHE:HD2	1.79	0.48
1:A:16:ALA:O	1:A:86:LEU:HD12	2.13	0.47
1:A:199:THR:HA	1:A:214:LYS:HA	1.95	0.47
2:B:189:ASP:O	2:B:193:ARG:HG3	2.13	0.47
1:H:160:ASN:ND2	1:H:199:THR:HG22	2.30	0.47
2:L:155:ILE:HD11	2:L:184:LEU:HD21	1.95	0.47
1:H:51:ILE:O	1:H:53:PRO:HD3	2.15	0.47
2:B:169:THR:HG23	2:B:170:ASP:O	2.15	0.46
1:H:61:ASP:OD1	1:H:63:LYS:HE3	2.16	0.46
1:A:143:LEU:HD13	1:A:198:VAL:HG21	1.97	0.46
2:B:67:ARG:HH12	2:B:88:ASP:CG	2.24	0.45
2:B:154:LYS:HD2	2:B:157:GLY:O	2.16	0.45
2:L:67:ARG:HB2	2:L:82:SER:O	2.15	0.45
2:L:151:VAL:HA	2:L:200:GLU:O	2.16	0.45
2:L:214:PHE:C	2:L:214:PHE:CD2	2.94	0.45
1:A:189:PRO:HG2	1:A:192:THR:HG22	1.98	0.45
2:L:191:TYR:CZ	2:L:216:ARG:HG3	2.51	0.45
1:A:198:VAL:HG23	1:A:198:VAL:O	2.17	0.45
2:L:190:GLU:O	2:L:194:HIS:HD2	2.00	0.45
2:B:190:GLU:HA	2:B:193:ARG:HD2	1.97	0.44
1:A:100:TRP:HE1	2:B:95:LYS:NZ	2.15	0.44
1:A:42:GLU:O	1:A:42:GLU:OE1	2.36	0.44
1:A:28:ASN:OD1	1:A:30:LYS:HG2	2.18	0.44
2:B:13:VAL:HG12	2:B:109:LEU:HD21	2.00	0.44
2:L:11:LEU:HD11	2:L:21:MET:HE2	1.98	0.43
2:L:4:MET:SD	2:L:25:SER:HB3	2.58	0.43
2:L:215:ASN:HB3	2:L:217:ASN:OD1	2.17	0.43
1:H:200:CYS:O	1:H:212:ASP:HA	2.18	0.43
1:A:18:VAL:HG12	1:A:86:LEU:HD21	2.01	0.43
2:B:36:LYS:HB3	2:B:56:TRP:CD1	2.54	0.43
1:A:51:ILE:O	1:A:53:PRO:HD3	2.19	0.42
1:H:50:ARG:CZ	1:H:59:LYS:HE2	2.49	0.42
1:A:67:LYS:HD2	1:A:84:SER:O	2.20	0.42

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:LYS:HE3	1:A:63:LYS:HB3	1.71	0.42
2:B:95:LYS:HE3	2:B:101:PHE:HD2	1.85	0.42
2:B:167:SER:O	2:B:180:MET:HG2	2.20	0.42
1:A:160:ASN:OD1	1:A:198:VAL:HA	2.20	0.41
2:L:41:TRP:CE2	2:L:79:LEU:HB2	2.56	0.41
2:L:168:TRP:CE2	2:L:180:MET:HG3	2.55	0.41
2:B:89:LEU:HD12	2:B:109:LEU:O	2.19	0.41
1:H:38:LYS:HB2	1:H:48:ILE:HD11	2.02	0.41
2:B:13:VAL:CG2	2:B:17:GLU:HB2	2.51	0.41
2:B:43:GLN:HB2	2:B:53:LEU:HD11	2.03	0.41
2:L:154:LYS:HG2	2:L:159:GLU:HA	2.03	0.41
1:A:11:LEU:HD22	1:A:121:THR:HG23	2.03	0.40
1:H:176:GLN:HB3	4:H:304:HOH:O	2.21	0.40
2:L:130:LEU:O	2:L:188:LYS:HD2	2.20	0.40
1:A:198:VAL:O	1:A:215:ILE:HG12	2.22	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	204/239 (85%)	198 (97%)	6 (3%)	0	100	100
1	H	204/239 (85%)	197 (97%)	7 (3%)	0	100	100
2	B	213/239 (89%)	204 (96%)	9 (4%)	0	100	100
2	L	217/239 (91%)	206 (95%)	10 (5%)	1 (0%)	24	37
3	P	5/11 (46%)	5 (100%)	0	0	100	100
3	Q	5/11 (46%)	5 (100%)	0	0	100	100
All	All	848/978 (87%)	815 (96%)	32 (4%)	1 (0%)	48	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	162	ASN

### 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	181/207 (87%)	179 (99%)	2 (1%)	65	82
1	H	181/207 (87%)	179 (99%)	2 (1%)	65	82
2	B	194/213 (91%)	191 (98%)	3 (2%)	57	77
2	L	196/213 (92%)	194 (99%)	2 (1%)	68	84
3	P	6/9 (67%)	6 (100%)	0	100	100
3	Q	6/9 (67%)	6 (100%)	0	100	100
All	All	764/858 (89%)	755 (99%)	9 (1%)	63	81

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	SER
1	A	185	SER
2	B	105	SER
2	B	187	THR
2	B	190	GLU
1	H	25	SER
1	H	155	VAL
2	L	20	THR
2	L	121	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	39	GLN
1	A	169	HIS
2	B	44	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
2	B	48	GLN
2	B	215	ASN
2	L	27	GLN
2	L	43	GLN
2	L	162	ASN
2	L	194	HIS
2	L	195	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	NIY	Q	52	3	13,15,16	1.10	0	13,20,22	1.91	4 (30%)
3	NIY	P	52	3	13,15,16	1.14	0	13,20,22	1.61	2 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NIY	Q	52	3	-	1/7/10/12	0/1/1/1
3	NIY	P	52	3	-	0/7/10/12	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Q	52	NIY	CB-CA-C	-4.60	102.85	111.47
3	P	52	NIY	CB-CA-C	-3.80	104.35	111.47
3	Q	52	NIY	OH-CZ-CE2	-2.56	112.37	119.33
3	Q	52	NIY	CD2-CG-CD1	2.46	121.98	118.54
3	P	52	NIY	O2-NN-CE1	2.27	122.92	119.03
3	Q	52	NIY	OH-CZ-CE1	2.07	127.42	121.31

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	Q	52	NIY	O-C-CA-CB

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	Q	52	NIY	2	0
3	P	52	NIY	2	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data

### 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	208/239 (87%)	0.86	19 (9%) 15 12	54, 103, 144, 159	0
1	H	208/239 (87%)	0.01	3 (1%) 73 69	42, 61, 100, 119	0
2	B	217/239 (90%)	0.99	20 (9%) 14 12	57, 122, 191, 204	0
2	L	219/239 (91%)	0.06	1 (0%) 87 85	46, 65, 102, 145	0
3	P	7/11 (63%)	-0.21	0 100 100	48, 51, 53, 56	0
3	Q	7/11 (63%)	0.46	0 100 100	67, 71, 82, 84	0
All	All	866/978 (88%)	0.47	43 (4%) 34 30	42, 81, 159, 204	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	217	PRO	4.1
2	B	135	ALA	3.8
2	B	186	LEU	3.5
2	B	11	LEU	3.4
1	A	12	VAL	3.4
1	H	217	PRO	3.3
2	B	149	ILE	3.3
2	B	151	VAL	3.2
2	B	207	THR	2.8
1	A	190	SER	2.7
2	B	160	ARG	2.7
2	B	191	TYR	2.7
1	A	197	THR	2.7
2	B	187	THR	2.6
1	A	94	TYR	2.6
2	B	165	LEU	2.5
2	B	184	LEU	2.5
1	A	189	PRO	2.5
1	A	16	ALA	2.5

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	17	SER	2.5
1	A	15	GLY	2.4
1	A	119	ALA	2.4
1	A	118	SER	2.3
1	A	169	HIS	2.3
1	A	45	LEU	2.3
1	A	194	PRO	2.3
1	H	191	SER	2.2
2	B	7	SER	2.2
2	B	150	ASN	2.2
2	B	168	TRP	2.2
1	A	44	GLY	2.2
2	B	188	LYS	2.2
2	L	134	GLY	2.2
2	B	153	TRP	2.1
2	B	138	VAL	2.1
2	B	164	VAL	2.1
2	B	214	PHE	2.1
2	B	166	ASN	2.1
1	H	163	SER	2.1
1	A	116	VAL	2.0
1	A	216	VAL	2.0
1	A	11	LEU	2.0
1	A	92	ALA	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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. 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	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NIY	Q	52	15/16	0.85	0.15	75,88,96,98	0
3	NIY	P	52	15/16	0.95	0.08	47,48,51,52	0

## 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.