



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 07:07 am BST

PDB ID : 4H0H
Title : Crystal structure of mimicry-recognizing 2D10 scFv with peptide
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Deposited on : 2012-09-08
Resolution : 2.00 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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 : 2.11
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) : 2.11

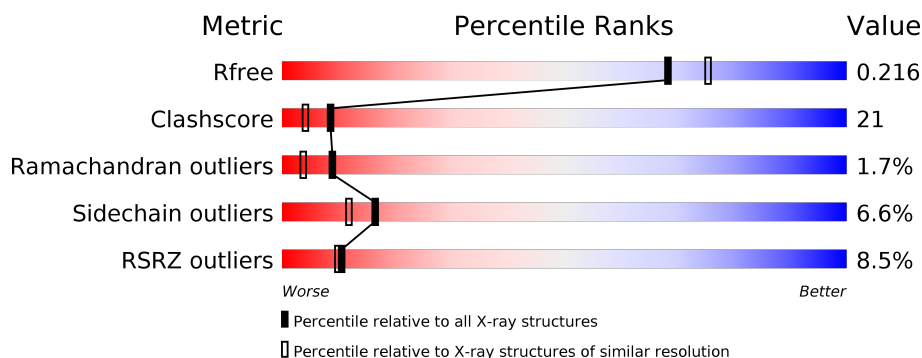
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 ≥ 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	B	251	<div> <div>6%</div> <div> <div></div> <div>76%</div> <div>17%</div> <div>• •</div> </div> </div>
2	D	12	<div> <div>58%</div> <div> <div>8%</div> <div>17%</div> <div>8%</div> <div>25%</div> <div>42%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2118 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 2D10 scFv.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	241	Total	C	N	O	S	0	0	0
			1848	1169	305	366	8			

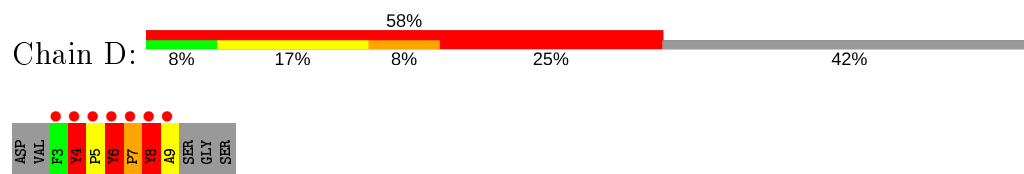
- Molecule 2 is a protein called peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	7	Total	C	N	O	0	0	0
			66	49	7	10			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	203	Total	O	0	0
			203	203		
3	D	1	Total	O	0	0
			1	1		

- Molecule 1: 2D10 scFv



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	80.33 Å 80.33 Å 74.98 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.50 – 2.00 25.50 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (25.50-2.00) 100.0 (25.50-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.72 (at 1.99 Å)	Xtriage
Refinement program	REFMAC 5.5.0102, CNS 1.1	Depositor
R, R_{free}	0.195 , 0.215 0.195 , 0.216	Depositor DCC
R_{free} test set	1897 reflections (9.84%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2118	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	B	1.28	4/1891 (0.2%)	0.91	1/2557 (0.0%)
2	D	2.50	3/71 (4.2%)	1.74	2/98 (2.0%)
All	All	1.35	7/1962 (0.4%)	0.95	3/2655 (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
2	D	0	2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	8	TYR	C-O	12.46	1.47	1.23
2	D	8	TYR	C-N	7.09	1.50	1.34
2	D	4	TYR	CB-CG	-7.04	1.41	1.51
1	B	153	SER	CB-OG	5.69	1.49	1.42
1	B	225	GLU	CG-CD	5.61	1.60	1.51
1	B	193	TYR	CG-CD1	5.44	1.46	1.39
1	B	103	TYR	CB-CG	5.17	1.59	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	108	ASP	CB-CG-OD1	5.61	123.35	118.30
2	D	8	TYR	CA-C-N	-5.51	105.07	117.20
2	D	4	TYR	CB-CA-C	-5.08	100.24	110.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	4	TYR	Peptide
2	D	8	TYR	Mainchain

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	B	1848	0	1800	70	0
2	D	66	0	54	32	0
3	B	203	0	0	4	0
3	D	1	0	0	0	0
All	All	2118	0	1854	78	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 21.

All (78) 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)
2:D:5:PRO:O	2:D:6:TYR:HD2	1.24	1.15
1:B:100:LYS:HZ2	2:D:7:PRO:HB2	1.17	1.09
1:B:100:LYS:NZ	2:D:7:PRO:HB2	1.72	1.04
2:D:5:PRO:O	2:D:6:TYR:CD2	2.14	1.00
1:B:53:ASN:CG	2:D:5:PRO:HD3	1.84	0.97
1:B:53:ASN:ND2	2:D:5:PRO:HD3	1.82	0.94
1:B:53:ASN:OD1	2:D:5:PRO:HD3	1.68	0.91
1:B:40:GLN:HE22	1:B:182:GLN:HE22	1.21	0.87
1:B:170:HIS:HD2	1:B:172:ASN:H	1.16	0.87
1:B:163:ARG:NH2	1:B:214:ASP:HB2	1.92	0.85
1:B:32:ASP:C	2:D:5:PRO:HB3	1.97	0.84
1:B:123:THR:O	1:B:124:THR:HB	1.79	0.81
1:B:163:ARG:CZ	1:B:214:ASP:HB2	2.10	0.80
1:B:100:LYS:HZ2	2:D:7:PRO:CB	1.93	0.80
1:B:163:ARG:NE	1:B:214:ASP:HB2	1.99	0.78
1:B:178:HIS:HD2	1:B:194:LYS:H	1.31	0.76
1:B:170:HIS:HE1	2:D:9:ALA:O	1.70	0.74
1:B:1:MET:HG2	1:B:1:MET:O	1.85	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:ASN:CG	2:D:5:PRO:CD	2.55	0.74
1:B:13:VAL:HG11	1:B:19:VAL:CG1	2.19	0.73
2:D:5:PRO:C	2:D:6:TYR:HD2	1.93	0.72
1:B:178:HIS:CD2	1:B:194:LYS:H	2.08	0.71
1:B:53:ASN:ND2	2:D:5:PRO:CD	2.55	0.70
1:B:53:ASN:OD1	2:D:5:PRO:CD	2.39	0.69
1:B:53:ASN:HD22	1:B:53:ASN:C	1.96	0.69
1:B:163:ARG:HH21	1:B:214:ASP:HB2	1.57	0.69
1:B:100:LYS:CG	2:D:7:PRO:HB3	2.23	0.68
1:B:218:LYS:HD3	3:B:354:HOH:O	1.96	0.65
1:B:13:VAL:HG21	1:B:87:LEU:CD1	2.26	0.65
1:B:163:ARG:HE	1:B:214:ASP:HB2	1.62	0.65
1:B:100:LYS:HG3	2:D:7:PRO:HB3	1.79	0.64
1:B:53:ASN:HD21	2:D:5:PRO:HD3	1.58	0.63
1:B:13:VAL:HG11	1:B:19:VAL:HG12	1.82	0.61
2:D:5:PRO:C	2:D:6:TYR:CD2	2.70	0.61
1:B:170:HIS:CE1	2:D:9:ALA:O	2.54	0.60
1:B:154:LEU:CD2	1:B:250:ILE:HD11	2.31	0.60
1:B:100:LYS:HZ3	2:D:7:PRO:HB2	1.65	0.60
1:B:48:TRP:HE1	1:B:51:ASN:ND2	2.00	0.60
1:B:13:VAL:HG21	1:B:87:LEU:HD13	1.82	0.60
2:D:8:TYR:HD2	2:D:9:ALA:H	1.49	0.60
1:B:170:HIS:CD2	1:B:172:ASN:H	2.09	0.58
1:B:163:ARG:NE	1:B:214:ASP:CB	2.66	0.58
1:B:167:SER:O	1:B:237:HIS:HE1	1.88	0.56
2:D:7:PRO:O	2:D:8:TYR:HD1	1.89	0.56
1:B:100:LYS:HE2	3:B:499:HOH:O	2.05	0.56
1:B:12:LEU:HD21	1:B:124:THR:O	2.06	0.55
1:B:100:LYS:HG2	2:D:7:PRO:HB3	1.88	0.55
1:B:36:LEU:HD12	1:B:51:ASN:HD21	1.73	0.54
1:B:34:ILE:HD12	2:D:5:PRO:HB2	1.90	0.54
1:B:33:TYR:N	2:D:5:PRO:HB3	2.22	0.53
1:B:100:LYS:HG2	2:D:7:PRO:CB	2.39	0.53
1:B:19:VAL:HG13	1:B:87:LEU:HD11	1.90	0.53
1:B:101:ASN:ND2	1:B:104:GLY:H	2.07	0.53
1:B:1:MET:O	1:B:2:GLU:O	2.27	0.52
1:B:181:LEU:HD13	1:B:230:TYR:CZ	2.44	0.52
1:B:1:MET:CG	1:B:1:MET:O	2.57	0.52
1:B:163:ARG:HE	1:B:214:ASP:CB	2.22	0.52
1:B:3:ILE:N	1:B:3:ILE:HD12	2.26	0.51
2:D:7:PRO:O	2:D:8:TYR:CD1	2.63	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:ASN:ND2	3:B:461:HOH:O	2.45	0.49
1:B:178:HIS:HD2	1:B:194:LYS:N	2.05	0.48
1:B:154:LEU:HD23	1:B:250:ILE:HD11	1.96	0.47
1:B:100:LYS:CG	2:D:7:PRO:CB	2.92	0.47
2:D:7:PRO:O	2:D:8:TYR:HB2	2.14	0.47
1:B:163:ARG:HE	1:B:214:ASP:CA	2.29	0.46
1:B:6:GLN:HG3	1:B:6:GLN:O	2.14	0.46
1:B:53:ASN:ND2	1:B:53:ASN:C	2.68	0.46
1:B:238:VAL:HG13	1:B:239:PRO:HA	1.99	0.44
1:B:13:VAL:HG21	1:B:87:LEU:HD12	1.99	0.44
1:B:32:ASP:O	2:D:5:PRO:HB3	2.19	0.43
1:B:154:LEU:HD21	1:B:250:ILE:HD11	2.01	0.43
1:B:247:LYS:CE	3:B:414:HOH:O	2.67	0.42
1:B:53:ASN:HD22	1:B:55:TYR:H	1.67	0.41
1:B:39:LYS:HB2	1:B:49:ILE:HD11	2.02	0.41
1:B:3:ILE:N	1:B:3:ILE:CD1	2.83	0.41
1:B:101:ASN:C	1:B:101:ASN:HD22	2.24	0.41
1:B:34:ILE:CD1	2:D:5:PRO:HD2	2.51	0.41
1:B:102:TYR:CD1	2:D:6:TYR:CG	3.09	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	B	237/251 (94%)	229 (97%)	6 (2%)	2 (1%)	19	13
2	D	5/12 (42%)	1 (20%)	2 (40%)	2 (40%)	0	0
All	All	242/263 (92%)	230 (95%)	8 (3%)	4 (2%)	9	4

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	2	GLU
1	B	124	THR
2	D	7	PRO
2	D	6	TYR

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	B	207/209 (99%)	196 (95%)	11 (5%)	22	18
2	D	6/10 (60%)	3 (50%)	3 (50%)	0	0
All	All	213/219 (97%)	199 (93%)	14 (7%)	16	12

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	6	GLN
1	B	12	LEU
1	B	53	ASN
1	B	66	LYS
1	B	100	LYS
1	B	101	ASN
1	B	122	LYS
1	B	124	THR
1	B	163	ARG
1	B	168	LEU
1	B	218	LYS
2	D	4	TYR
2	D	6	TYR
2	D	8	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	51	ASN

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Mol	Chain	Res	Type
1	B	53	ASN
1	B	101	ASN
1	B	170	HIS
1	B	178	HIS
1	B	182	GLN
1	B	186	GLN
1	B	237	HIS

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](#)

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, 95th 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 ²)	Q<0.9
1	B	241/251 (96%)	0.04	14 (5%) 23 22	17, 26, 43, 65	0
2	D	7/12 (58%)	10.17	7 (100%) 0 0	89, 92, 94, 96	7 (100%)
All	All	248/263 (94%)	0.33	21 (8%) 10 10	17, 26, 58, 96	7 (2%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	9	ALA	13.3
2	D	8	TYR	13.0
2	D	7	PRO	11.0
2	D	4	TYR	9.1
2	D	6	TYR	9.1
1	B	137	GLY	8.5
2	D	5	PRO	8.5
2	D	3	PHE	7.1
1	B	138	GLY	6.4
1	B	136	GLY	6.4
1	B	1	MET	5.2
1	B	42	HIS	3.5
1	B	125	GLY	3.4
1	B	121	ALA	3.3
1	B	55	TYR	2.9
1	B	123	THR	2.9
1	B	124	THR	2.8
1	B	139	SER	2.6
1	B	148	PHE	2.2
1	B	120	SER	2.1
1	B	76	SER	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.