



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 06:55 PM GMT

PDB ID : 1D5B
Title : UNLIGANDED MATURE OXY-COPE CATALYTIC ANTIBODY
Authors : Mundorff, E.C.; Hanson, M.A.; Schultz, P.G.; Stevens, R.C.
Deposited on : 1999-10-06
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at 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.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

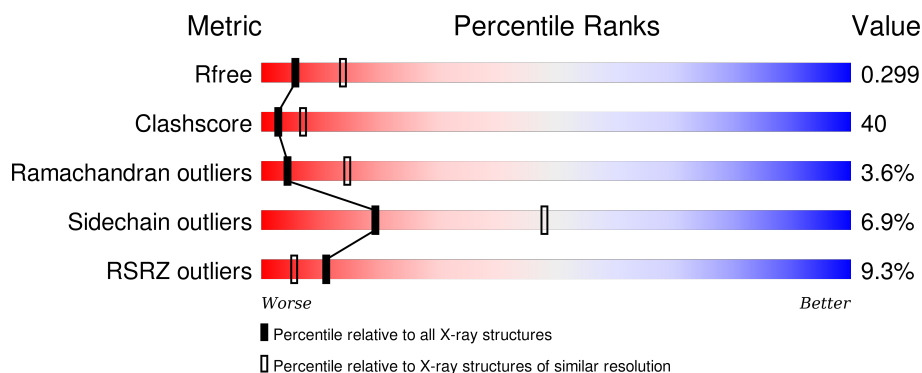
1 Overall quality at a glance

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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	211	<div> <div>7%</div> <div>36%</div> <div>56%</div> <div>7%</div> </div>
1	L	211	<div> <div>7%</div> <div>40%</div> <div>52%</div> <div>8%</div> </div>
2	B	221	<div> <div>12%</div> <div>51%</div> <div>43%</div> <div>5%</div> </div>
2	H	221	<div> <div>11%</div> <div>48%</div> <div>45%</div> <div>6%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6717 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 chimeric OXY-COPE catalytic ANTIBODY AZ-28 (light chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1656	1040	273	336	7			
1	L	211	Total	C	N	O	S	0	0	0
			1656	1040	273	336	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	108	ARG	-	NOT IN SWS ENTRY	UNP P01834
L	108	ARG	-	NOT IN SWS ENTRY	UNP P01834

- Molecule 2 is a protein called chimeric OXY-COPE catalytic ANTIBODY AZ-28 (HEAVY chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	221	Total	C	N	O	S	0	0	0
			1665	1056	273	330	6			
2	H	221	Total	C	N	O	S	0	0	0
			1665	1056	273	330	6			

- Molecule 3 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cd	0	0
			1	1		

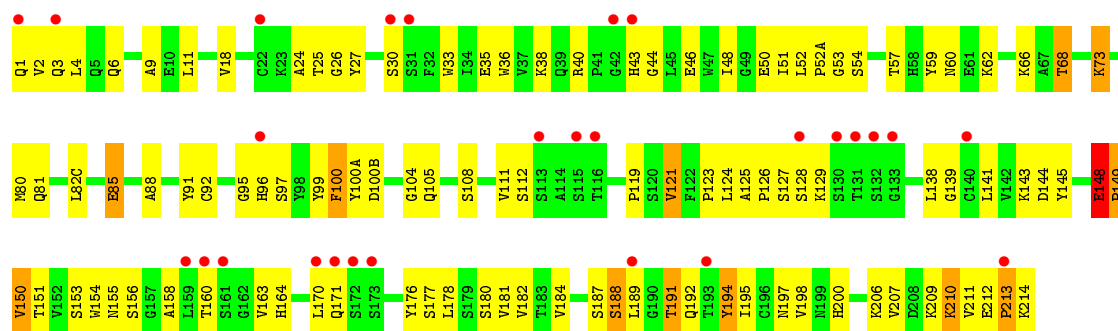
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	26	Total	O	0	0
			26	26		

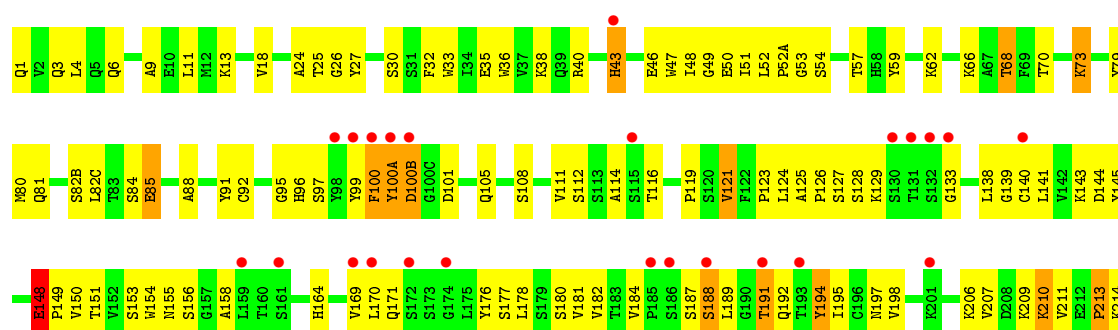
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	13	Total 13	O 13	0	0
4	H	17	Total 17	O 17	0	0
4	L	18	Total 18	O 18	0	0



- Molecule 2: chimeric OXY-COPE catalytic ANTIBODY AZ-28 (HEAVY chain)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.42Å 80.67Å 134.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 19.95 – 2.80	Depositor EDS
% Data completeness (in resolution range)	79.2 (20.00-2.80) 87.5 (19.95-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 2.79Å)	Xtriage
Refinement program	CNS 0.4	Depositor
R, R_{free}	0.232 , 0.285 0.249 , 0.299	Depositor DCC
R_{free} test set	1848 reflections (9.85%)	DCC
Wilson B-factor (Å ²)	50.0	Xtriage
Anisotropy	0.576	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 86.3	EDS
Estimated twinning fraction	0.000 for k,h,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	2 of 18775 reflections (0.011%)	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6717	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 34.33 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 6.9430e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CD

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	0.52	1/1692 (0.1%)	0.72	1/2290 (0.0%)
1	L	0.56	3/1692 (0.2%)	0.74	2/2290 (0.1%)
2	B	0.44	0/1711	0.73	1/2326 (0.0%)
2	H	0.53	0/1711	0.79	2/2326 (0.1%)
All	All	0.52	4/6806 (0.1%)	0.74	6/9232 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	79	GLU	CG-CD	5.67	1.60	1.51
1	L	143	GLU	CD-OE2	-5.35	1.19	1.25
1	A	143	GLU	CB-CG	5.26	1.62	1.52
1	L	79	GLU	CB-CG	5.02	1.61	1.52

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	143	GLU	OE1-CD-OE2	-6.29	115.75	123.30
2	H	148	GLU	N-CA-C	6.00	127.21	111.00
2	B	148	GLU	N-CA-C	5.81	126.69	111.00
1	L	143	GLU	OE1-CD-OE2	-5.57	116.62	123.30
1	L	79	GLU	OE1-CD-OE2	-5.41	116.81	123.30

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	1656	0	1607	153	0
1	L	1656	0	1607	142	0
2	B	1665	0	1609	122	0
2	H	1665	0	1609	129	0
3	A	1	0	0	0	0
4	A	26	0	0	9	0
4	B	13	0	0	1	0
4	H	17	0	0	6	0
4	L	18	0	0	10	0
All	All	6717	0	6432	518	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 40.

The worst 5 of 518 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:80:TYR:CD2	1:A:168:SER:HB3	1.84	1.11
1:A:150:VAL:HG23	1:A:151:ASP:H	1.06	1.10
1:A:80:TYR:HD2	1:A:168:SER:HB3	1.07	1.10
1:L:150:VAL:HG23	1:L:151:ASP:H	1.09	1.08
1:L:185:ASP:HA	1:L:188:LYS:HD3	1.42	1.00

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	209/211 (99%)	175 (84%)	27 (13%)	7 (3%)	5	16
1	L	209/211 (99%)	175 (84%)	26 (12%)	8 (4%)	4	13
2	B	219/221 (99%)	188 (86%)	23 (10%)	8 (4%)	4	14
2	H	219/221 (99%)	187 (85%)	24 (11%)	8 (4%)	4	14
All	All	856/864 (99%)	725 (85%)	100 (12%)	31 (4%)	4	14

5 of 31 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	143	GLU
1	A	150	VAL
2	B	128	SER
1	L	143	GLU
1	L	150	VAL

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	190/190 (100%)	178 (94%)	12 (6%)	22	53
1	L	190/190 (100%)	178 (94%)	12 (6%)	22	53
2	B	185/185 (100%)	172 (93%)	13 (7%)	19	47
2	H	185/185 (100%)	170 (92%)	15 (8%)	15	39
All	All	750/750 (100%)	698 (93%)	52 (7%)	19	48

5 of 52 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	206	LYS
1	L	135	LEU
2	H	191	THR
2	B	210	LYS
1	L	21	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	105	GLN
1	L	69	GLN
2	H	58	HIS
2	B	164	HIS
2	B	204	ASN

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 1 ligands modelled in this entry, 1 is 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 ⓘ

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, 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	A	211/211 (100%)	0.38	15 (7%)	19 10	20, 51, 88, 95	0
1	L	211/211 (100%)	0.22	14 (6%)	22 13	20, 43, 78, 94	0
2	B	221/221 (100%)	0.57	27 (12%)	5 3	17, 52, 87, 95	0
2	H	221/221 (100%)	0.42	24 (10%)	7 3	11, 44, 86, 95	0
All	All	864/864 (100%)	0.40	80 (9%)	11 5	11, 47, 87, 95	0

The worst 5 of 80 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	130	SER	9.4
2	H	130	SER	8.6
2	H	131	THR	7.3
2	H	133	GLY	5.5
2	B	193	THR	5.4

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, 95th 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(\AA^2)	Q<0.9
3	CD	A	212	1/1	0.98	0.14	-1.26	52,52,52,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.