



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:23 PM UTC

PDB ID : 4LAS / pdb_00004las
Title : Crystal structure of a therapeutic single chain antibody in complex with 4-hydroxymethamphetamine
Authors : Celical, R.; Gokulan, K.; Peterson, E.C.; Varughese, K.I.
Deposited on : 2013-06-20
Resolution : 2.33 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 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 : 2022.3.0, CSD as543be (2022)
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

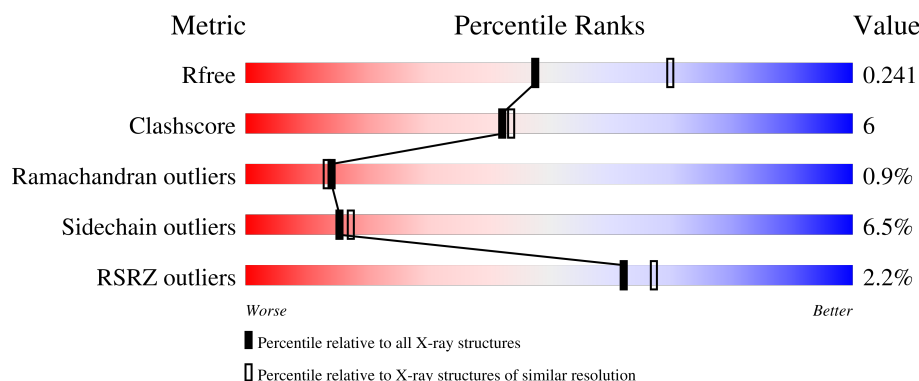
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

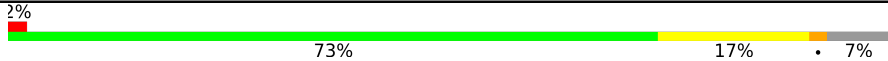
The reported resolution of this entry is 2.33 Å.

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	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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 ≥ 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	H	249	

2 Entry composition [i](#)

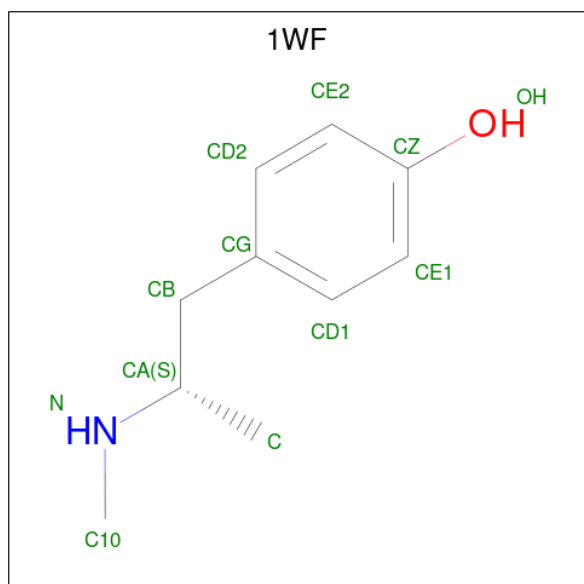
There are 3 unique types of molecules in this entry. The entry contains 1858 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Single chain antibody fragment scFv6H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	232	Total	C	N	O	S	0	0	0
			1775	1117	286	364	8			

- Molecule 2 is 4-[(2S)-2-(methylamino)propyl]phenol (CCD ID: 1WF) (formula: C₁₀H₁₅NO).

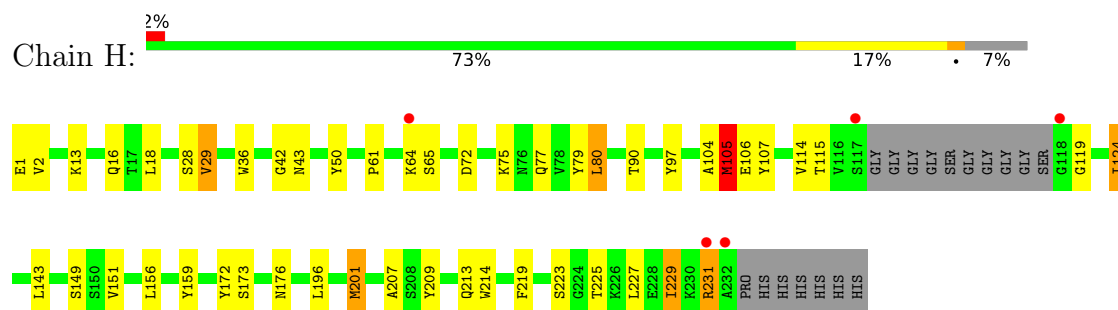


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	H	1	Total	C	N	O	0	0
			12	10	1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	71	Total	O	0	0
			71	71		

- Molecule 1: Single chain antibody fragment scFv6H4



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	34.28Å 65.30Å 48.57Å 90.00° 98.36° 90.00°	Depositor
Resolution (Å)	33.94 – 2.33 33.94 – 2.33	Depositor EDS
% Data completeness (in resolution range)	93.7 (33.94-2.33) 93.7 (33.94-2.33)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.47 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.7.0029, CNS	Depositor
R, R_{free}	0.160 , 0.238 0.170 , 0.241	Depositor DCC
R_{free} test set	405 reflections (4.43%)	wwPDB-VP
Wilson B-factor (Å ²)	26.6	Xtriage
Anisotropy	0.440	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	1858	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.07% 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

Bond lengths and bond angles in the following residue types are not validated in this section: 1WF

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	H	1.15	8/1821 (0.4%)	1.01	3/2473 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	149	SER	C-N	10.70	1.47	1.33
1	H	124	ILE	C-O	-6.59	1.16	1.24
1	H	106	GLU	C-O	-6.16	1.16	1.23
1	H	207	ALA	C-O	-6.00	1.18	1.23
1	H	105	MET	C-O	-5.97	1.16	1.23
1	H	149	SER	C-O	-5.43	1.17	1.23
1	H	227	LEU	C-O	-5.18	1.17	1.24
1	H	107	TYR	C-O	-5.07	1.18	1.24

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	149	SER	O-C-N	8.85	133.28	123.10
1	H	149	SER	CA-C-N	-5.41	112.58	120.75
1	H	149	SER	C-N-CA	-5.41	112.58	120.75

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	H	1775	0	1681	22	0
2	H	12	0	14	0	0
3	H	71	0	0	0	1
All	All	1858	0	1695	22	1

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 (22) 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:H:29:VAL:HG11	1:H:97:TYR:OH	1.74	0.87
1:H:13:LYS:O	1:H:16:GLN:HB2	1.94	0.68
1:H:196:LEU:C	1:H:196:LEU:HD23	2.32	0.54
1:H:104:ALA:HB2	1:H:172:TYR:CZ	2.43	0.54
1:H:42:GLY:O	1:H:43:ASN:HB2	2.05	0.54
1:H:90:THR:HG23	1:H:115:THR:HA	1.91	0.53
1:H:77:GLN:NE2	1:H:79:TYR:OH	2.44	0.50
1:H:151:VAL:HG11	1:H:213:GLN:HG3	1.94	0.49
1:H:18:LEU:C	1:H:18:LEU:HD23	2.37	0.49
1:H:72:ASP:CG	1:H:75:LYS:HG2	2.39	0.47
1:H:201:MET:HE2	1:H:229:ILE:HG23	1.96	0.46
1:H:36:TRP:CE2	1:H:80:LEU:HB2	2.51	0.45
1:H:18:LEU:HD12	1:H:114:VAL:HG11	1.97	0.45
1:H:173:SER:HB2	1:H:176:ASN:HD22	1.83	0.44
1:H:156:LEU:C	1:H:156:LEU:HD13	2.43	0.43
1:H:229:ILE:HD13	1:H:229:ILE:HG21	1.83	0.42
1:H:214:TRP:HA	1:H:219:PHE:CD1	2.54	0.42
1:H:124:ILE:HD13	1:H:151:VAL:HG12	2.01	0.41
1:H:159:TYR:O	1:H:209:TYR:HA	2.21	0.41
1:H:143:LEU:HD22	1:H:225:THR:HG21	2.02	0.41
1:H:105:MET:HA	1:H:105:MET:HE2	2.02	0.41
1:H:50:TYR:C	1:H:50:TYR:CD1	3.00	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:602:HOH:O	3:H:622:HOH:O[2_545]	2.07	0.13

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	H	228/249 (92%)	219 (96%)	7 (3%)	2 (1%)	14 13

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	119	GLY
1	H	231	ARG

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	H	201/210 (96%)	188 (94%)	13 (6%)	15 17

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	1	GLU
1	H	2	VAL
1	H	28	SER
1	H	29	VAL
1	H	61	PRO
1	H	64	LYS
1	H	65	SER
1	H	80	LEU
1	H	105	MET

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Mol	Chain	Res	Type
1	H	201	MET
1	H	223	SER
1	H	229	ILE
1	H	231	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	39	GLN
1	H	43	ASN
1	H	77	GLN
1	H	160	GLN
1	H	161	GLN
1	H	176	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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
2	1WF	H	500	-	12,12,12	1.71	3 (25%)	14,15,15	1.09	1 (7%)

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
2	1WF	H	500	-	-	2/6/6/6	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	500	1WF	CA-N	-3.70	1.42	1.48
2	H	500	1WF	OH-CZ	-2.40	1.31	1.37
2	H	500	1WF	CD1-CE1	-2.30	1.35	1.38

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	500	1WF	C10-N-CA	3.31	118.34	114.39

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	H	500	1WF	CB-CA-N-C10
2	H	500	1WF	C-CA-N-C10

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 [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, 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	H	232/249 (93%)	-0.20	5 (2%) 62 67	15, 26, 48, 99	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	232	ALA	6.4
1	H	118	GLY	3.9
1	H	117	SER	2.8
1	H	231	ARG	2.7
1	H	64	LYS	2.6

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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	1WF	H	500	12/12	0.94	0.07	19,20,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.