



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 14, 2026 – 10:08 AM JST

PDB ID : 9ULM / pdb_00009ulm
Title : Mogamulizumab in complex with CCR4 N-terminus peptide (N2-C29)
Authors : Heo, Y.-S.
Deposited on : 2025-04-20
Resolution : 2.01 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

<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
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

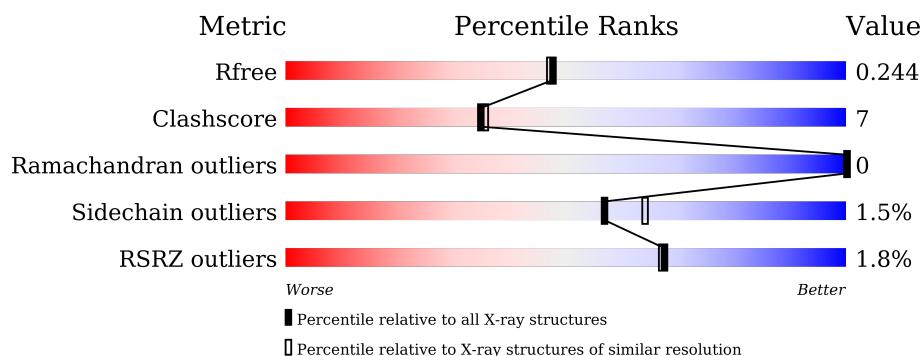
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (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 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	A	231	<div> <div>83%</div> <div>10%</div> <div>7%</div> </div>
1	C	231	<div> <div>81%</div> <div>11%</div> <div>8%</div> </div>
2	B	219	<div> <div>88%</div> <div>11%</div> <div>.</div> </div>
2	D	219	<div> <div>83%</div> <div>16%</div> <div>.</div> </div>
3	I	130	<div> <div>14%</div> <div>55%</div> <div>22%</div> <div>.</div> <div>21%</div> </div>
3	K	130	<div> <div>%</div> <div>79%</div> <div>13%</div> <div>8%</div> </div>

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Mol	Chain	Length	Quality of chain
4	X	28	<div><div><div>4%</div><div>32%</div><div>7%</div><div>61%</div></div></div>
4	Y	28	<div><div><div>25%</div><div>14%</div><div>61%</div></div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9355 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 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1607	1017	270	314	6			
1	C	213	Total	C	N	O	S	0	0	0
			1603	1015	269	313	6			

- Molecule 2 is a protein called light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	216	Total	C	N	O	S	0	0	0
			1677	1058	282	332	5			
2	D	217	Total	C	N	O	S	0	0	0
			1681	1060	283	333	5			

- Molecule 3 is a protein called anti-kappa VHH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	K	120	Total	C	N	O	S	0	0	0
			915	567	160	184	4			
3	I	103	Total	C	N	O	S	0	0	0
			800	499	137	160	4			

- Molecule 4 is a protein called C-C chemokine receptor type 4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	X	11	Total	C	N	O	0	0	0
			99	66	12	21			
4	Y	11	Total	C	N	O	0	0	0
			99	66	12	21			

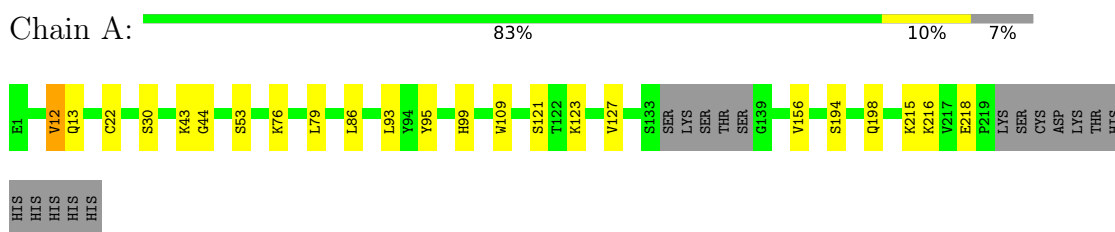
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	177	Total 177	O 177	0	0
5	B	192	Total 192	O 192	0	0
5	C	161	Total 161	O 161	0	0
5	D	178	Total 178	O 178	0	0
5	K	105	Total 105	O 105	0	0
5	I	31	Total 31	O 31	0	0
5	X	13	Total 13	O 13	0	0
5	Y	17	Total 17	O 17	0	0

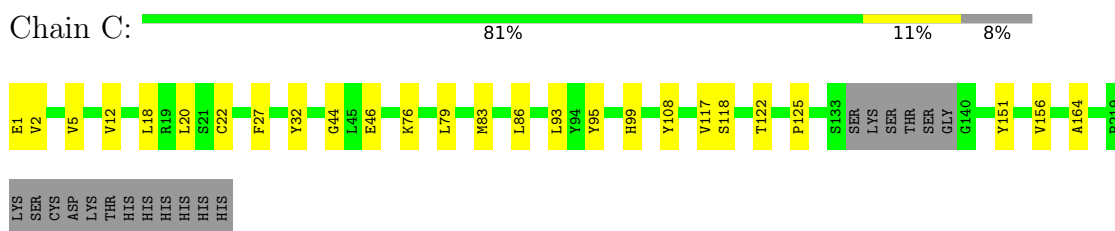
3 Residue-property plots

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.

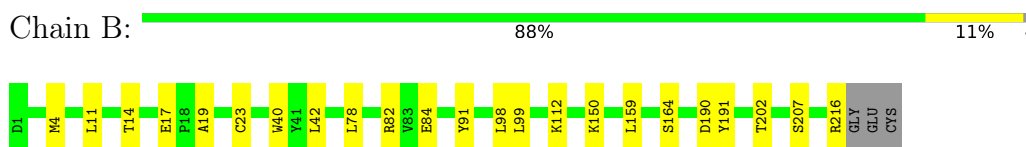
- Molecule 1: heavy chain



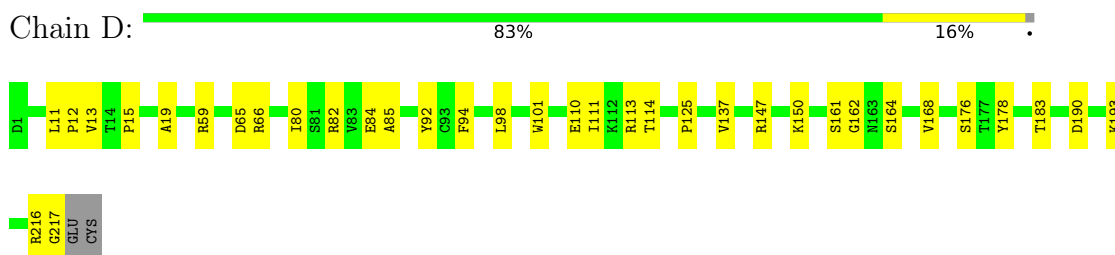
- Molecule 1: heavy chain



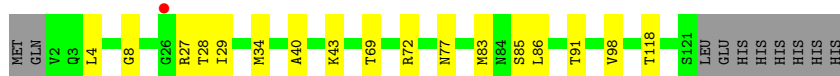
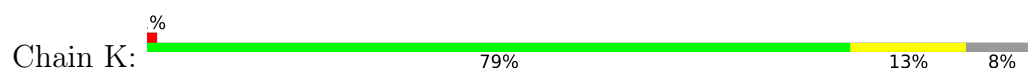
- Molecule 2: light chain



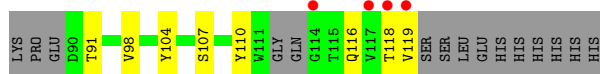
- Molecule 2: light chain



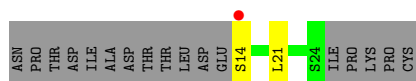
- Molecule 3: anti-kappa VHH



• Molecule 3: anti-kappa VHH



• Molecule 4: C-C chemokine receptor type 4



• Molecule 4: C-C chemokine receptor type 4



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	65.39Å 71.19Å 73.45Å 81.24° 64.41° 71.05°	Depositor
Resolution (Å)	29.55 – 2.01 29.55 – 2.01	Depositor EDS
% Data completeness (in resolution range)	93.7 (29.55-2.01) 93.7 (29.55-2.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.01Å)	Xtriage
Refinement program	PHENIX (1.20_4459: ???)	Depositor
R, R_{free}	0.197 , 0.244 0.196 , 0.244	Depositor DCC
R_{free} test set	3526 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	29.1	Xtriage
Anisotropy	0.469	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,-h+l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9355	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.33	0/1647	0.52	0/2243
1	C	0.29	0/1643	0.46	0/2238
2	B	0.33	0/1716	0.54	0/2332
2	D	0.30	0/1720	0.50	0/2337
3	I	0.31	0/812	0.54	1/1092 (0.1%)
3	K	0.35	0/934	0.50	0/1264
4	X	0.26	0/102	0.65	0/138
4	Y	0.34	0/102	0.48	0/138
All	All	0.32	0/8676	0.51	1/11782 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	34	MET	CA-CB-CG	-5.08	103.95	114.10

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	1607	0	1562	15	0
1	C	1603	0	1559	18	0
2	B	1677	0	1642	15	0
2	D	1681	0	1645	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	I	800	0	735	38	0
3	K	915	0	854	11	0
4	X	99	0	84	2	0
4	Y	99	0	84	4	0
5	A	177	0	0	2	4
5	B	192	0	0	4	2
5	C	161	0	0	6	0
5	D	178	0	0	9	4
5	I	31	0	0	6	0
5	K	105	0	0	3	2
5	X	13	0	0	0	0
5	Y	17	0	0	2	0
All	All	9355	0	8165	124	6

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 7.

The worst 5 of 124 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)
3:I:116:GLN:NE2	5:I:202:HOH:O	1.94	1.00
1:A:43:LYS:HD3	1:A:44:GLY:H	1.31	0.92
1:A:123:LYS:NZ	5:A:301:HOH:O	2.06	0.88
4:Y:14:SER:N	5:Y:101:HOH:O	2.10	0.84
3:I:19:ARG:HH21	3:I:82:GLN:HB3	1.44	0.82

The worst 5 of 6 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)
5:A:440:HOH:O	5:D:356:HOH:O[1_456]	1.96	0.24
5:A:449:HOH:O	5:D:315:HOH:O[1_456]	1.98	0.22
5:A:477:HOH:O	5:B:491:HOH:O[1_654]	1.98	0.22
5:D:451:HOH:O	5:K:270:HOH:O[1_565]	2.12	0.08
5:B:335:HOH:O	5:K:238:HOH:O[1_556]	2.13	0.07

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	210/231 (91%)	209 (100%)	1 (0%)	0	100	100
1	C	209/231 (90%)	208 (100%)	1 (0%)	0	100	100
2	B	214/219 (98%)	212 (99%)	2 (1%)	0	100	100
2	D	215/219 (98%)	214 (100%)	1 (0%)	0	100	100
3	I	91/130 (70%)	89 (98%)	2 (2%)	0	100	100
3	K	118/130 (91%)	117 (99%)	1 (1%)	0	100	100
4	X	9/28 (32%)	9 (100%)	0	0	100	100
4	Y	9/28 (32%)	9 (100%)	0	0	100	100
All	All	1075/1216 (88%)	1067 (99%)	8 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	179/196 (91%)	175 (98%)	4 (2%)	45	50
1	C	179/196 (91%)	177 (99%)	2 (1%)	65	73
2	B	193/195 (99%)	191 (99%)	2 (1%)	68	75
2	D	193/195 (99%)	191 (99%)	2 (1%)	68	75
3	I	83/106 (78%)	80 (96%)	3 (4%)	31	31
3	K	95/106 (90%)	95 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	X	11/27 (41%)	11 (100%)	0	100	100
4	Y	11/27 (41%)	10 (91%)	1 (9%)	9	6
All	All	944/1048 (90%)	930 (98%)	14 (2%)	57	64

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	156	VAL
2	D	13	VAL
4	Y	19	TYR
3	I	65	GLN
3	I	81	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	58	ASN
3	K	3	GLN
3	I	84	ASN
3	K	77	ASN
2	B	50	GLN

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](#)

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, 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	214/231 (92%)	-0.30	0 100 100	22, 30, 45, 62	0
1	C	213/231 (92%)	-0.10	0 100 100	24, 36, 50, 58	0
2	B	216/219 (98%)	-0.30	0 100 100	21, 30, 49, 58	0
2	D	217/219 (99%)	-0.12	0 100 100	24, 36, 49, 56	0
3	I	103/130 (79%)	1.13	18 (17%) 4 3	48, 69, 90, 103	0
3	K	120/130 (92%)	-0.18	1 (0%) 82 82	20, 32, 53, 80	0
4	X	11/28 (39%)	0.80	1 (9%) 15 13	40, 43, 51, 52	0
4	Y	11/28 (39%)	-0.08	0 100 100	31, 34, 41, 48	0
All	All	1105/1216 (90%)	-0.07	20 (1%) 67 67	20, 34, 67, 103	0

The worst 5 of 20 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	I	119	VAL	4.6
3	I	8	GLY	3.0
3	I	79	VAL	3.0
3	I	34	MET	2.6
3	I	86	LEU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.