



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:53 AM UTC

PDB ID : 1U71 / pdb_00001u71
Title : Understanding the Role of Leu22 Variants in Methotrexate Resistance: Comparison of Wild-type and Leu22Arg Variant Mouse and Human Dihydrofolate Reductase Ternary Crystal Complexes with Methotrexate and NADPH
Authors : Cody, V.; Luft, J.R.; Pangborn, W.
Deposited on : 2004-08-02
Resolution : 2.20 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

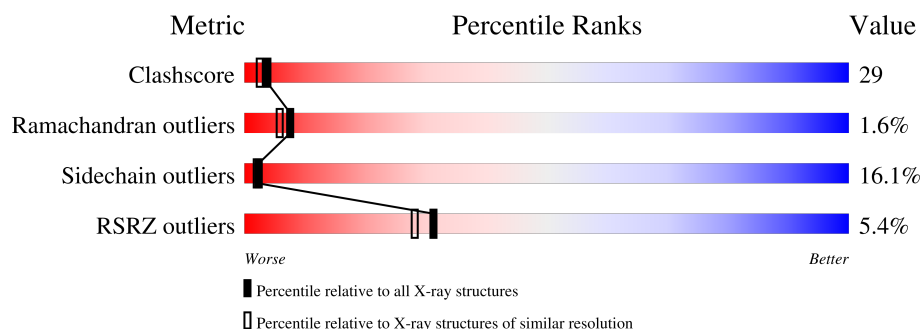
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	186	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	SO4	A	189	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 1607 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 Dihydrofolate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	186	Total	C	N	O	S	0	0	0
			1505	963	256	279	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	ARG	LEU	engineered mutation	UNP P00374

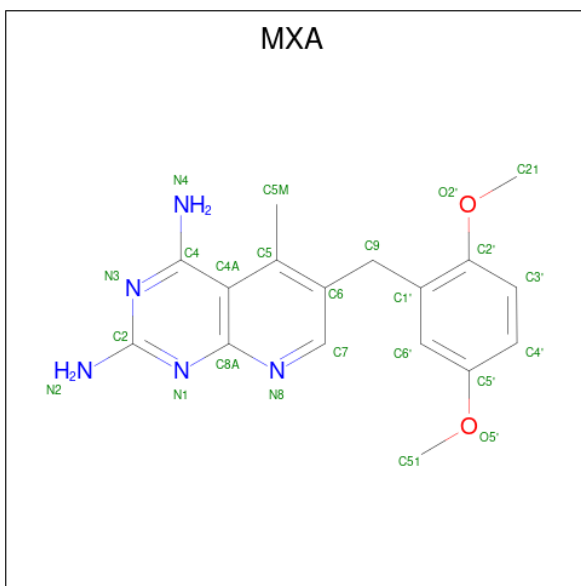
- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 6-(2,5-DIMETHOXY-BENZYL)-5-METHYL-PYRIDO[2,3-D]PYRIMIDINE-

2,4-DIAMINE (CCD ID: MXA) (formula: $C_{17}H_{19}N_5O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			24	17	5	2		

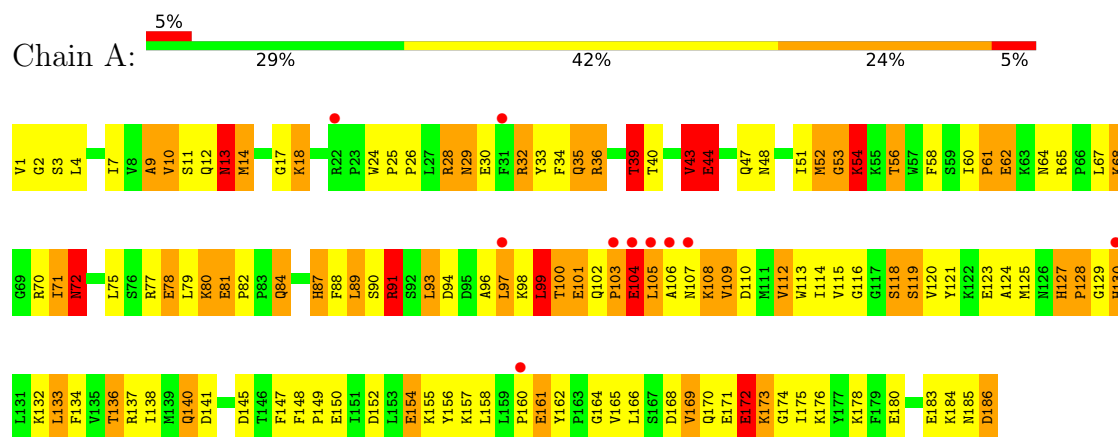
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	68	Total	O	0	0
			68	68		

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: Dihydrofolate reductase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	85.14Å 85.14Å 78.04Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	87.1 (50.00-2.20) 64.1 (50.00-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.00 (at 2.18Å)	Xtriage
Refinement program	PROLSQ	Depositor
R, R_{free}	0.211 , 0.236 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	10.6	Xtriage
Anisotropy	0.157	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 130.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.28$, $\langle L^2 \rangle = 0.11$	Xtriage
Estimated twinning fraction	0.238 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	1607	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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: SO4, MXA

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	1.40	9/1540 (0.6%)	3.00	173/2076 (8.3%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	14	MET	C-N	-5.54	1.28	1.33
1	A	129	GLY	N-CA	5.52	1.52	1.45
1	A	150	GLU	C-O	5.38	1.30	1.23
1	A	71	ILE	CA-CB	5.38	1.60	1.54
1	A	17	GLY	C-N	-5.28	1.26	1.33
1	A	145	ASP	N-CA	5.22	1.53	1.46
1	A	119	SER	CB-OG	5.15	1.52	1.42
1	A	36	ARG	NE-CZ	5.06	1.38	1.33
1	A	40	THR	CA-CB	5.00	1.60	1.53

All (173) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	91	ARG	CD-NE-CZ	29.31	165.43	124.40
1	A	140	GLN	OE1-CD-NE2	18.95	141.55	122.60
1	A	12	GLN	OE1-CD-NE2	12.73	135.33	122.60
1	A	174	GLY	CA-C-N	11.91	138.60	123.12
1	A	174	GLY	C-N-CA	11.91	138.60	123.12
1	A	44	GLU	CB-CG-CD	11.73	132.55	112.60
1	A	58	PHE	CA-CB-CG	-11.33	102.47	113.80
1	A	145	ASP	CB-CA-C	10.47	125.10	109.29
1	A	9	ALA	CA-C-O	-10.25	109.85	120.71
1	A	3	SER	CA-C-O	-9.92	110.74	121.56
1	A	3	SER	N-CA-C	9.86	124.40	110.50
1	A	113	TRP	CA-C-O	-9.63	109.96	120.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	PHE	O-C-N	9.63	134.23	123.25
1	A	140	GLN	CB-CG-CD	-9.51	96.43	112.60
1	A	115	VAL	CA-C-O	-9.46	110.68	120.14
1	A	58	PHE	CA-C-O	-9.36	107.86	119.38
1	A	13	ASN	OD1-CG-ND2	-9.18	113.42	122.60
1	A	10	VAL	CA-CB-CG1	9.00	125.71	110.40
1	A	9	ALA	CA-C-N	8.84	135.44	122.71
1	A	9	ALA	C-N-CA	8.84	135.44	122.71
1	A	148	PHE	CA-C-O	-8.78	111.60	120.64
1	A	109	VAL	N-CA-C	8.75	120.87	108.36
1	A	13	ASN	CA-CB-CG	8.71	121.31	112.60
1	A	88	PHE	CA-C-O	-8.59	112.05	121.33
1	A	56	THR	CA-CB-OG1	-8.57	96.75	109.60
1	A	93	LEU	CA-C-N	8.53	131.53	120.44
1	A	93	LEU	C-N-CA	8.53	131.53	120.44
1	A	171	GLU	O-C-N	8.45	133.83	123.44
1	A	158	LEU	CA-C-O	-8.33	110.92	120.58
1	A	29	ASN	CA-CB-CG	8.23	120.83	112.60
1	A	3	SER	O-C-N	8.19	132.52	122.86
1	A	53	GLY	CA-C-N	8.18	131.07	120.44
1	A	53	GLY	C-N-CA	8.18	131.07	120.44
1	A	62	GLU	CA-C-N	8.13	131.98	120.28
1	A	62	GLU	C-N-CA	8.13	131.98	120.28
1	A	138	ILE	O-C-N	8.02	133.55	122.97
1	A	147	PHE	N-CA-C	7.70	122.03	109.72
1	A	72	ASN	CA-CB-CG	7.64	120.24	112.60
1	A	87	HIS	CA-C-O	-7.63	110.79	119.41
1	A	12	GLN	N-CA-CB	7.61	123.26	110.32
1	A	58	PHE	O-C-N	7.58	132.59	122.43
1	A	172	GLU	CG-CD-OE1	7.54	135.74	118.40
1	A	67	LEU	N-CA-C	-7.50	98.57	109.59
1	A	150	GLU	CB-CG-CD	7.47	125.30	112.60
1	A	128	PRO	CB-CA-C	7.47	121.31	110.63
1	A	3	SER	CA-CB-OG	-7.46	96.17	111.10
1	A	81	GLU	O-C-N	7.44	126.47	121.71
1	A	137	ARG	NE-CZ-NH1	-7.44	114.06	121.50
1	A	30	GLU	N-CA-C	-7.41	103.29	111.36
1	A	68	LYS	O-C-N	7.39	131.31	122.96
1	A	145	ASP	N-CA-C	-7.27	103.60	114.64
1	A	14	MET	CA-C-N	7.20	129.20	120.92
1	A	14	MET	C-N-CA	7.20	129.20	120.92
1	A	149	PRO	CB-CA-C	7.17	122.04	111.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	9	ALA	N-CA-C	-7.13	98.11	109.59
1	A	173	LYS	CA-C-O	-7.12	112.18	121.08
1	A	118	SER	O-C-N	7.11	130.54	122.22
1	A	121	TYR	O-C-N	7.04	131.38	122.23
1	A	133	LEU	CA-C-O	-7.03	112.72	120.38
1	A	12	GLN	CA-CB-CG	-6.99	100.12	114.10
1	A	64	ASN	CA-CB-CG	6.99	119.59	112.60
1	A	154	GLU	N-CA-C	-6.96	104.26	112.89
1	A	129	GLY	N-CA-C	-6.96	103.63	111.63
1	A	138	ILE	CB-CA-C	6.95	120.50	110.98
1	A	165	VAL	CA-C-O	-6.93	112.86	120.48
1	A	78	GLU	N-CA-CB	6.89	120.67	110.33
1	A	127	HIS	N-CA-C	-6.88	99.56	109.48
1	A	75	LEU	CA-C-O	-6.87	112.76	120.32
1	A	116	GLY	CA-C-N	6.87	134.06	121.70
1	A	116	GLY	C-N-CA	6.87	134.06	121.70
1	A	62	GLU	CB-CG-CD	6.82	124.20	112.60
1	A	96	ALA	O-C-N	6.78	129.31	122.12
1	A	123	GLU	CA-CB-CG	6.73	127.56	114.10
1	A	51	ILE	O-C-N	-6.71	116.13	123.18
1	A	138	ILE	N-CA-C	-6.65	98.29	107.99
1	A	170	GLN	O-C-N	6.63	132.18	122.97
1	A	11	SER	CA-C-O	-6.60	114.35	121.94
1	A	160	PRO	CA-C-O	6.58	128.58	119.18
1	A	56	THR	N-CA-C	-6.48	103.72	111.69
1	A	178	LYS	CA-C-N	6.47	132.30	122.65
1	A	178	LYS	C-N-CA	6.47	132.30	122.65
1	A	118	SER	CA-CB-OG	-6.44	98.23	111.10
1	A	186	ASP	CA-CB-CG	-6.43	106.17	112.60
1	A	53	GLY	N-CA-C	-6.39	102.73	112.51
1	A	90	SER	CA-CB-OG	-6.39	98.33	111.10
1	A	10	VAL	CB-CA-C	-6.33	99.75	110.71
1	A	132	LYS	O-C-N	6.31	130.97	123.27
1	A	4	LEU	O-C-N	-6.29	115.85	123.27
1	A	120	VAL	O-C-N	6.23	128.84	121.80
1	A	33	TYR	N-CA-CB	-6.19	100.69	110.28
1	A	175	ILE	CA-C-N	6.17	131.84	122.65
1	A	175	ILE	C-N-CA	6.17	131.84	122.65
1	A	87	HIS	ND1-CE1-NE2	-6.14	102.26	108.40
1	A	119	SER	O-C-N	6.13	128.71	122.09
1	A	40	THR	N-CA-C	6.13	118.60	109.59
1	A	168	ASP	N-CA-CB	-6.12	101.57	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	170	GLN	N-CA-CB	6.10	119.78	110.70
1	A	12	GLN	CA-C-O	-6.08	112.14	119.49
1	A	30	GLU	CA-CB-CG	6.07	126.25	114.10
1	A	67	LEU	CB-CA-C	6.05	119.33	110.26
1	A	157	LYS	CA-CB-CG	-6.02	102.06	114.10
1	A	99	LEU	CA-C-O	6.02	126.94	120.20
1	A	61	PRO	CA-C-N	6.00	132.99	121.54
1	A	61	PRO	C-N-CA	6.00	132.99	121.54
1	A	7	ILE	CB-CA-C	5.97	119.05	110.33
1	A	162	TYR	CA-C-O	-5.96	113.14	119.22
1	A	91	ARG	CG-CD-NE	5.94	125.08	112.00
1	A	124	ALA	CA-C-N	5.93	132.02	120.99
1	A	124	ALA	C-N-CA	5.93	132.02	120.99
1	A	137	ARG	CA-C-O	-5.91	113.82	120.43
1	A	28	ARG	CB-CA-C	5.89	120.70	110.68
1	A	18	LYS	N-CA-CB	5.88	121.12	110.95
1	A	180	GLU	CG-CD-OE1	5.86	131.87	118.40
1	A	52	MET	CA-C-O	-5.80	115.07	121.33
1	A	101	GLU	N-CA-CB	5.79	118.73	110.16
1	A	129	GLY	CA-C-O	-5.79	117.41	122.16
1	A	28	ARG	N-CA-CB	-5.78	101.39	110.06
1	A	134	PHE	O-C-N	-5.77	115.86	123.13
1	A	54	LYS	CB-CA-C	-5.75	101.85	110.88
1	A	81	GLU	CA-CB-CG	5.72	125.54	114.10
1	A	147	PHE	CA-CB-CG	-5.71	108.09	113.80
1	A	128	PRO	CA-C-N	-5.70	113.45	122.92
1	A	128	PRO	C-N-CA	-5.70	113.45	122.92
1	A	115	VAL	N-CA-CB	-5.68	105.74	112.39
1	A	140	GLN	CG-CD-OE1	-5.68	109.44	120.80
1	A	2	GLY	N-CA-C	-5.68	99.73	113.18
1	A	164	GLY	CA-C-O	5.63	125.53	119.06
1	A	129	GLY	O-C-N	5.60	128.38	123.06
1	A	51	ILE	CB-CG1-CD1	-5.55	102.15	113.80
1	A	103	PRO	N-CA-C	-5.55	106.83	113.98
1	A	34	PHE	N-CA-CB	5.54	118.10	110.07
1	A	36	ARG	CD-NE-CZ	-5.52	116.68	124.40
1	A	170	GLN	CA-C-N	-5.51	112.30	121.89
1	A	170	GLN	C-N-CA	-5.51	112.30	121.89
1	A	97	LEU	CB-CA-C	5.50	119.48	110.95
1	A	161	GLU	CG-CD-OE2	-5.50	105.75	118.40
1	A	96	ALA	CA-C-O	-5.48	114.74	120.55
1	A	130	HIS	CA-CB-CG	5.48	119.28	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	162	TYR	CA-CB-CG	-5.46	104.07	113.90
1	A	81	GLU	CG-CD-OE1	5.42	130.88	118.40
1	A	34	PHE	CA-C-N	5.40	127.47	120.44
1	A	34	PHE	C-N-CA	5.40	127.47	120.44
1	A	17	GLY	O-C-N	5.39	128.57	123.56
1	A	78	GLU	CA-CB-CG	5.38	124.85	114.10
1	A	32	ARG	CA-C-N	5.36	128.46	120.31
1	A	32	ARG	C-N-CA	5.36	128.46	120.31
1	A	168	ASP	CB-CA-C	5.34	119.87	110.36
1	A	110	ASP	CA-C-O	-5.33	112.89	120.51
1	A	118	SER	N-CA-C	5.33	117.84	111.71
1	A	51	ILE	N-CA-CB	-5.32	104.01	111.25
1	A	65	ARG	NE-CZ-NH2	-5.32	114.41	119.20
1	A	80	LYS	CG-CD-CE	5.31	123.51	111.30
1	A	94	ASP	O-C-N	5.30	127.53	122.07
1	A	105	LEU	N-CA-CB	-5.30	102.64	110.53
1	A	54	LYS	CA-CB-CG	-5.29	103.52	114.10
1	A	169	VAL	CB-CA-C	5.29	119.30	111.26
1	A	84	GLN	CB-CG-CD	5.28	121.58	112.60
1	A	172	GLU	CA-CB-CG	5.26	124.63	114.10
1	A	39	THR	CA-CB-OG1	-5.26	101.71	109.60
1	A	136	THR	N-CA-CB	5.24	119.16	110.47
1	A	112	VAL	CB-CA-C	5.23	118.37	110.63
1	A	48	ASN	CA-C-O	5.21	127.87	121.72
1	A	39	THR	OG1-CB-CG2	5.20	119.70	109.30
1	A	113	TRP	O-C-N	5.18	129.47	123.30
1	A	150	GLU	CA-CB-CG	5.17	124.44	114.10
1	A	43	VAL	CB-CA-C	-5.16	102.82	111.29
1	A	78	GLU	CG-CD-OE2	-5.16	106.54	118.40
1	A	96	ALA	N-CA-CB	5.15	117.69	110.12
1	A	141	ASP	O-C-N	5.14	129.08	123.22
1	A	140	GLN	CG-CD-NE2	-5.13	108.71	116.40
1	A	137	ARG	CD-NE-CZ	-5.05	117.33	124.40
1	A	65	ARG	N-CA-C	5.02	117.43	109.40
1	A	3	SER	N-CA-CB	-5.01	102.95	110.17

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	1505	0	1513	88	0
2	A	10	0	0	3	0
3	A	24	0	19	1	0
4	A	68	0	0	6	0
All	All	1607	0	1532	88	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 29.

All (88) 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:105:LEU:HA	1:A:108:LYS:HG3	1.34	1.05
1:A:35:GLN:HE21	1:A:70:ARG:HH12	1.19	0.88
1:A:103:PRO:HD2	4:A:207:HOH:O	1.75	0.86
1:A:47:GLN:HG3	4:A:235:HOH:O	1.78	0.83
1:A:35:GLN:NE2	1:A:70:ARG:HH12	1.77	0.82
1:A:28:ARG:O	1:A:32:ARG:HG3	1.82	0.80
1:A:35:GLN:HE22	1:A:70:ARG:HH22	1.29	0.79
1:A:93:LEU:HD11	1:A:114:ILE:HD11	1.67	0.76
1:A:105:LEU:HA	1:A:108:LYS:CG	2.18	0.72
1:A:79:LEU:O	1:A:91:ARG:NH2	2.24	0.71
1:A:26:PRO:O	1:A:173:LYS:HE2	1.91	0.71
1:A:97:LEU:O	1:A:100:THR:HB	1.92	0.70
1:A:184:LYS:HE3	1:A:186:ASP:OD1	1.92	0.70
1:A:35:GLN:NE2	1:A:70:ARG:HH22	1.90	0.69
1:A:89:LEU:HG	1:A:91:ARG:HH11	1.58	0.68
1:A:35:GLN:O	1:A:39:THR:OG1	2.12	0.67
1:A:26:PRO:O	1:A:173:LYS:CE	2.42	0.67
1:A:72:ASN:H	1:A:87:HIS:HD2	1.40	0.67
1:A:154:GLU:OE2	4:A:222:HOH:O	2.12	0.66
1:A:99:LEU:HA	1:A:102:GLN:HG2	1.78	0.66
1:A:35:GLN:HE21	1:A:70:ARG:NH1	1.93	0.65
1:A:184:LYS:HG2	1:A:185:ASN:N	2.10	0.65
1:A:105:LEU:CA	1:A:108:LYS:HG3	2.21	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:GLN:C	1:A:104:GLU:H	2.04	0.64
1:A:104:GLU:O	1:A:108:LYS:NZ	2.31	0.62
1:A:72:ASN:H	1:A:87:HIS:CD2	2.16	0.61
1:A:102:GLN:O	1:A:106:ALA:HB2	2.00	0.61
1:A:61:PRO:HD2	3:A:187:MXA:H213	1.85	0.59
1:A:93:LEU:HD11	1:A:114:ILE:CD1	2.33	0.59
1:A:1:VAL:HG13	1:A:100:THR:HG23	1.85	0.58
1:A:71:ILE:HD13	1:A:105:LEU:HD22	1.86	0.58
1:A:161:GLU:O	4:A:228:HOH:O	2.17	0.58
1:A:102:GLN:C	1:A:104:GLU:N	2.60	0.58
1:A:101:GLU:O	1:A:102:GLN:C	2.47	0.57
1:A:130:HIS:HE1	1:A:183:GLU:OE1	1.88	0.57
1:A:130:HIS:CE1	1:A:183:GLU:OE1	2.58	0.56
1:A:98:LYS:O	1:A:102:GLN:NE2	2.38	0.56
1:A:71:ILE:HD13	1:A:105:LEU:CD2	2.35	0.56
1:A:127:HIS:HB2	4:A:257:HOH:O	2.05	0.55
1:A:35:GLN:NE2	1:A:70:ARG:NH1	2.51	0.55
1:A:81:GLU:HB2	1:A:82:PRO:HD2	1.87	0.55
1:A:1:VAL:CG1	1:A:100:THR:CG2	2.86	0.53
1:A:93:LEU:O	1:A:97:LEU:HG	2.08	0.53
1:A:184:LYS:HE3	1:A:186:ASP:CG	2.34	0.52
1:A:156:TYR:CD2	1:A:184:LYS:HB2	2.45	0.52
1:A:97:LEU:O	1:A:98:LYS:C	2.51	0.52
1:A:105:LEU:O	1:A:108:LYS:N	2.40	0.52
1:A:29:ASN:HB2	1:A:172:GLU:OE1	2.12	0.50
1:A:89:LEU:HD11	1:A:91:ARG:HH12	1.75	0.49
1:A:1:VAL:HG11	1:A:100:THR:HG21	1.93	0.49
1:A:89:LEU:CG	1:A:91:ARG:HH11	2.25	0.49
1:A:89:LEU:CG	1:A:91:ARG:NH1	2.76	0.48
1:A:186:ASP:OD1	1:A:186:ASP:C	2.53	0.48
1:A:89:LEU:HD21	1:A:91:ARG:NH1	2.28	0.48
1:A:105:LEU:HB3	1:A:109:VAL:CG2	2.43	0.48
1:A:54:LYS:HE3	4:A:240:HOH:O	2.13	0.48
1:A:60:ILE:O	1:A:61:PRO:C	2.56	0.47
1:A:89:LEU:HD11	1:A:91:ARG:NH1	2.28	0.47
1:A:100:THR:O	1:A:106:ALA:HA	2.15	0.47
1:A:54:LYS:HG3	1:A:79:LEU:HD11	1.96	0.47
1:A:105:LEU:C	1:A:108:LYS:HG3	2.40	0.47
1:A:184:LYS:CE	1:A:186:ASP:OD1	2.62	0.47
1:A:52:MET:HB2	1:A:56:THR:HB	1.97	0.46
1:A:105:LEU:HB3	1:A:109:VAL:HG23	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:VAL:HG11	1:A:100:THR:CG2	2.45	0.46
1:A:35:GLN:NE2	1:A:70:ARG:NH2	2.61	0.46
1:A:105:LEU:O	1:A:106:ALA:C	2.59	0.45
1:A:184:LYS:HG2	1:A:185:ASN:H	1.81	0.45
1:A:29:ASN:OD1	1:A:32:ARG:NH1	2.50	0.45
1:A:169:VAL:HG11	1:A:176:LYS:HD3	1.99	0.44
1:A:24:TRP:HB2	1:A:25:PRO:HD2	1.99	0.44
1:A:43:VAL:O	1:A:44:GLU:C	2.61	0.44
1:A:1:VAL:HG21	1:A:100:THR:HG21	1.99	0.43
1:A:77:ARG:N	2:A:189:SO4:O1	2.44	0.42
1:A:125:MET:HE3	1:A:125:MET:HB3	1.81	0.42
1:A:9:ALA:HA	1:A:136:THR:HB	2.01	0.42
1:A:36:ARG:HG2	1:A:36:ARG:NH1	2.35	0.42
1:A:156:TYR:CE2	1:A:184:LYS:HB2	2.55	0.42
1:A:1:VAL:CG1	1:A:100:THR:HG21	2.48	0.42
1:A:81:GLU:HB2	1:A:82:PRO:CD	2.49	0.42
1:A:100:THR:O	1:A:106:ALA:CA	2.68	0.42
1:A:77:ARG:HB2	2:A:189:SO4:O1	2.20	0.42
1:A:98:LYS:C	1:A:102:GLN:HE21	2.24	0.42
1:A:152:ASP:OD1	1:A:155:LYS:CE	2.68	0.41
1:A:99:LEU:HD23	1:A:105:LEU:HD12	2.03	0.41
1:A:79:LEU:HB2	1:A:89:LEU:HD11	2.03	0.41
1:A:13:ASN:O	1:A:14:MET:HB2	2.21	0.41
1:A:53:GLY:HA3	2:A:188:SO4:S	2.61	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	184/186 (99%)	171 (93%)	10 (5%)	3 (2%)	7 6

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	GLU
1	A	43	VAL
1	A	104	GLU

5.3.2 Protein sidechains [i](#)

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	168/168 (100%)	141 (84%)	27 (16%)	2 2

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	A	13	ASN
1	A	18	LYS
1	A	35	GLN
1	A	39	THR
1	A	54	LYS
1	A	62	GLU
1	A	68	LYS
1	A	72	ASN
1	A	78	GLU
1	A	80	LYS
1	A	84	GLN
1	A	89	LEU
1	A	91	ARG
1	A	99	LEU
1	A	100	THR
1	A	104	GLU
1	A	107	ASN
1	A	108	LYS
1	A	112	VAL
1	A	118	SER
1	A	119	SER

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Mol	Chain	Res	Type
1	A	128	PRO
1	A	133	LEU
1	A	140	GLN
1	A	166	LEU
1	A	172	GLU

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

Mol	Chain	Res	Type
1	A	5	ASN
1	A	12	GLN
1	A	35	GLN
1	A	47	GLN
1	A	87	HIS
1	A	185	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](#)

3 ligands 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
2	SO4	A	188	-	4,4,4	0.61	0	6,6,6	0.65	0
3	MXA	A	187	-	25,26,26	1.16	2 (8%)	36,37,37	1.65	9 (25%)
2	SO4	A	189	-	4,4,4	0.51	0	6,6,6	0.71	0

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	MXA	A	187	-	-	1/8/8/8	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	187	MXA	C4-N3	-2.72	1.29	1.33
3	A	187	MXA	C4'-C5'	-2.56	1.33	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	187	MXA	C51-O5'-C5'	-3.38	110.26	117.50
3	A	187	MXA	C1'-C9-C6	3.35	121.65	112.64
3	A	187	MXA	C5M-C5-C6	-2.93	117.48	120.81
3	A	187	MXA	N8-C8A-N1	-2.71	112.81	115.77
3	A	187	MXA	C9-C6-C7	-2.66	117.95	122.20
3	A	187	MXA	O5'-C5'-C4'	2.50	131.69	119.82
3	A	187	MXA	O5'-C5'-C6'	-2.37	108.68	120.00
3	A	187	MXA	C4A-C4-N4	2.23	126.67	122.70
3	A	187	MXA	C9-C6-C5	2.11	123.51	121.19

There are no chirality outliers.

All (1) torsion outliers are listed below:

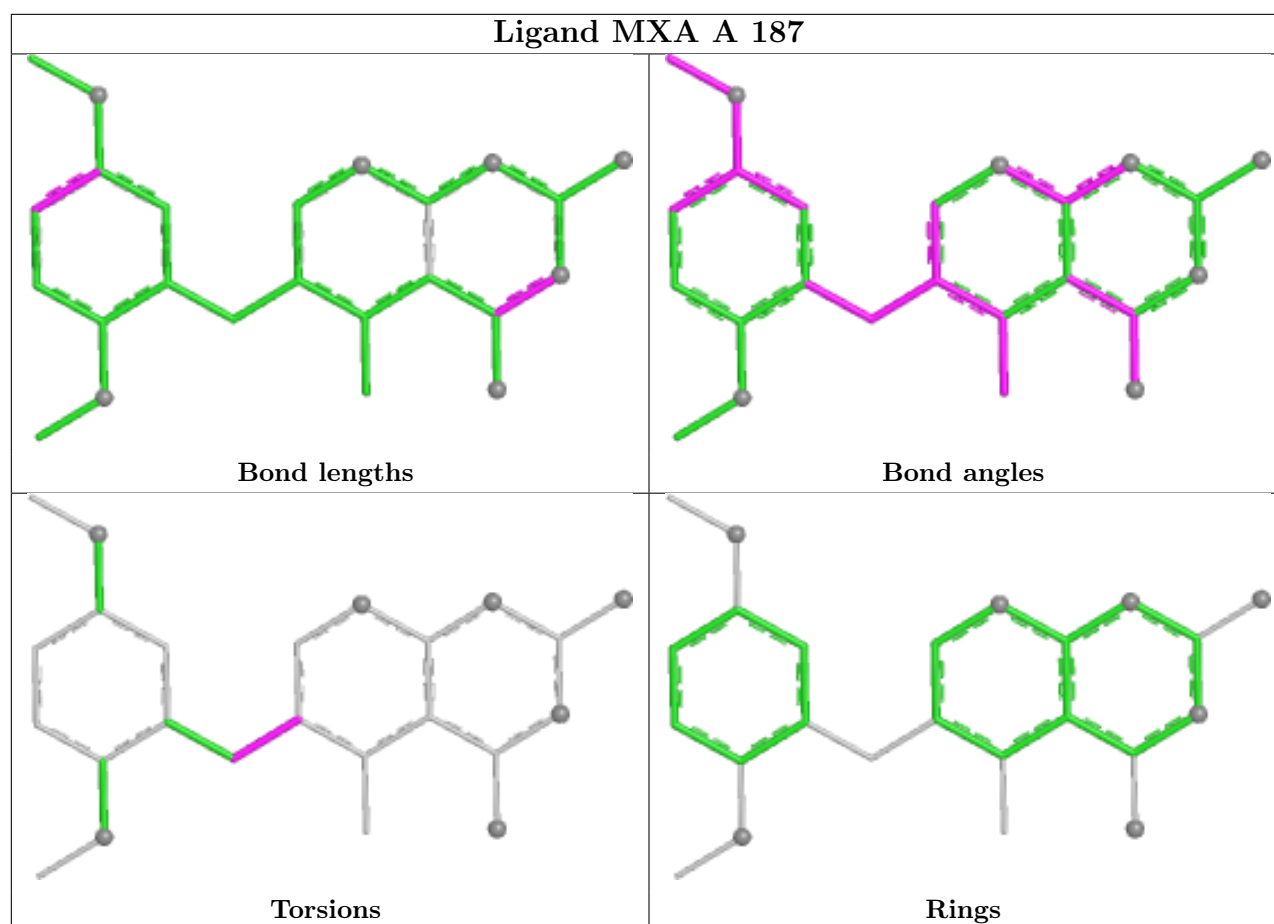
Mol	Chain	Res	Type	Atoms
3	A	187	MXA	C7-C6-C9-C1'

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	188	SO4	1	0
3	A	187	MXA	1	0
2	A	189	SO4	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

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	186/186 (100%)	0.60	10 (5%) 31 28	10, 22, 43, 53	1 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	22	ARG	3.8
1	A	103	PRO	3.2
1	A	105	LEU	2.9
1	A	31	PHE	2.7
1	A	104	GLU	2.6
1	A	130	HIS	2.5
1	A	106	ALA	2.3
1	A	107	ASN	2.3
1	A	97	LEU	2.3
1	A	160	PRO	2.2

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](#)

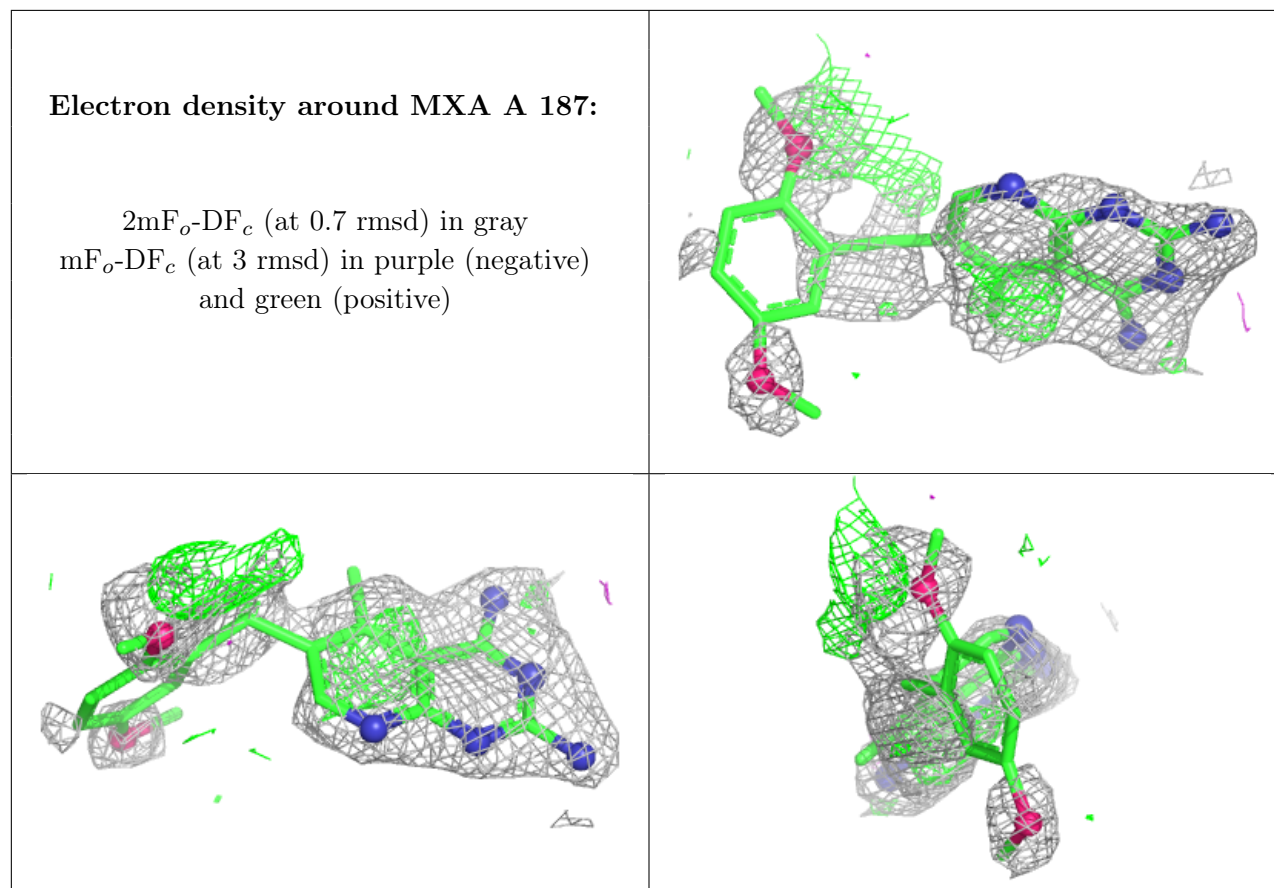
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
3	MXA	A	187	24/24	0.74	0.28	2,5,14,16	24
2	SO4	A	189	5/5	0.85	0.12	45,45,45,45	0
2	SO4	A	188	5/5	0.98	0.06	20,20,20,21	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers ⓘ

There are no such residues in this entry.