



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2024 – 03:51 PM EDT

PDB ID : 1XPG  
Title : Crystal Structure of T. maritima Cobalamin-Independent Methionine Synthase complexed with Zn<sup>2+</sup> and Methyltetrahydrofolate  
Authors : Pejchal, R.; Ludwig, M.L.  
Deposited on : 2004-10-08  
Resolution : 2.59 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](mailto: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>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
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)	:	2.37.1

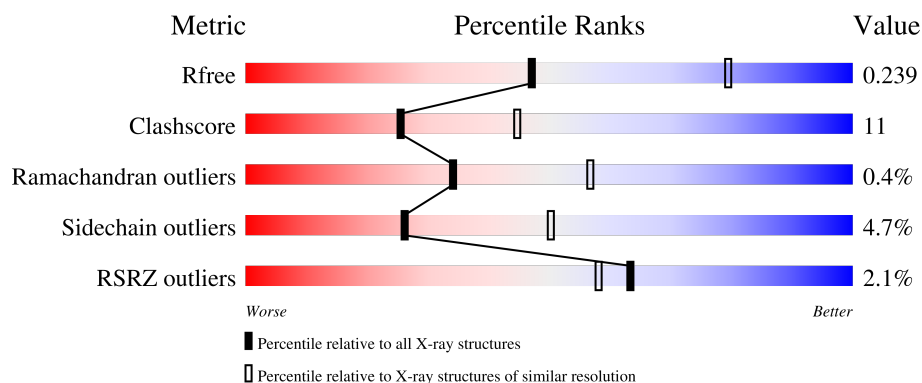
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

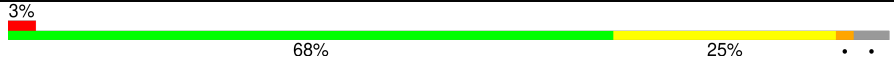

The reported resolution of this entry is 2.59 Å.

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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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  $\geq 3$ , 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	765	
1	B	765	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11993 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 5-methyltetrahydropteroyltriglutamate--homocysteine methyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	734	Total	C	N	O	S	Se	0	0	0
			5805	3773	948	1060	7	17			
1	B	734	Total	C	N	O	S	Se	0	0	0
			5827	3779	951	1073	7	17			

There are 102 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-31	MET	-	EXPRESSION TAG	UNP Q9X112
A	-30	HIS	-	EXPRESSION TAG	UNP Q9X112
A	-29	HIS	-	EXPRESSION TAG	UNP Q9X112
A	-28	HIS	-	EXPRESSION TAG	UNP Q9X112
A	-27	HIS	-	EXPRESSION TAG	UNP Q9X112
A	-26	HIS	-	EXPRESSION TAG	UNP Q9X112
A	-25	HIS	-	EXPRESSION TAG	UNP Q9X112
A	-24	GLY	-	EXPRESSION TAG	UNP Q9X112
A	-23	LYS	-	EXPRESSION TAG	UNP Q9X112
A	-22	PRO	-	EXPRESSION TAG	UNP Q9X112
A	-21	ILE	-	EXPRESSION TAG	UNP Q9X112
A	-20	PRO	-	EXPRESSION TAG	UNP Q9X112
A	-19	ASN	-	EXPRESSION TAG	UNP Q9X112
A	-18	PRO	-	EXPRESSION TAG	UNP Q9X112
A	-17	LEU	-	EXPRESSION TAG	UNP Q9X112
A	-16	LEU	-	EXPRESSION TAG	UNP Q9X112
A	-15	GLY	-	EXPRESSION TAG	UNP Q9X112
A	-14	LEU	-	EXPRESSION TAG	UNP Q9X112
A	-13	ASP	-	EXPRESSION TAG	UNP Q9X112
A	-12	SER	-	EXPRESSION TAG	UNP Q9X112
A	-11	THR	-	EXPRESSION TAG	UNP Q9X112
A	-10	GLU	-	EXPRESSION TAG	UNP Q9X112
A	-9	ASN	-	EXPRESSION TAG	UNP Q9X112
A	-8	LEU	-	EXPRESSION TAG	UNP Q9X112

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	TYR	-	EXPRESSION TAG	UNP Q9X112
A	-6	PHE	-	EXPRESSION TAG	UNP Q9X112
A	-5	GLN	-	EXPRESSION TAG	UNP Q9X112
A	-4	GLY	-	EXPRESSION TAG	UNP Q9X112
A	-3	ILE	-	EXPRESSION TAG	UNP Q9X112
A	-2	ASP	-	EXPRESSION TAG	UNP Q9X112
A	-1	PRO	-	EXPRESSION TAG	UNP Q9X112
A	0	PHE	-	EXPRESSION TAG	UNP Q9X112
A	1	THR	-	EXPRESSION TAG	UNP Q9X112
A	38	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	43	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	45	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	73	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	93	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	102	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	170	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	344	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	421	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	468	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	510	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	529	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	537	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	619	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	682	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	721	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	728	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
A	733	ALA	-	EXPRESSION TAG	UNP Q9X112
B	-31	MET	-	EXPRESSION TAG	UNP Q9X112
B	-30	HIS	-	EXPRESSION TAG	UNP Q9X112
B	-29	HIS	-	EXPRESSION TAG	UNP Q9X112
B	-28	HIS	-	EXPRESSION TAG	UNP Q9X112
B	-27	HIS	-	EXPRESSION TAG	UNP Q9X112
B	-26	HIS	-	EXPRESSION TAG	UNP Q9X112
B	-25	HIS	-	EXPRESSION TAG	UNP Q9X112
B	-24	GLY	-	EXPRESSION TAG	UNP Q9X112
B	-23	LYS	-	EXPRESSION TAG	UNP Q9X112
B	-22	PRO	-	EXPRESSION TAG	UNP Q9X112
B	-21	ILE	-	EXPRESSION TAG	UNP Q9X112
B	-20	PRO	-	EXPRESSION TAG	UNP Q9X112
B	-19	ASN	-	EXPRESSION TAG	UNP Q9X112
B	-18	PRO	-	EXPRESSION TAG	UNP Q9X112
B	-17	LEU	-	EXPRESSION TAG	UNP Q9X112

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	LEU	-	EXPRESSION TAG	UNP Q9X112
B	-15	GLY	-	EXPRESSION TAG	UNP Q9X112
B	-14	LEU	-	EXPRESSION TAG	UNP Q9X112
B	-13	ASP	-	EXPRESSION TAG	UNP Q9X112
B	-12	SER	-	EXPRESSION TAG	UNP Q9X112
B	-11	THR	-	EXPRESSION TAG	UNP Q9X112
B	-10	GLU	-	EXPRESSION TAG	UNP Q9X112
B	-9	ASN	-	EXPRESSION TAG	UNP Q9X112
B	-8	LEU	-	EXPRESSION TAG	UNP Q9X112
B	-7	TYR	-	EXPRESSION TAG	UNP Q9X112
B	-6	PHE	-	EXPRESSION TAG	UNP Q9X112
B	-5	GLN	-	EXPRESSION TAG	UNP Q9X112
B	-4	GLY	-	EXPRESSION TAG	UNP Q9X112
B	-3	ILE	-	EXPRESSION TAG	UNP Q9X112
B	-2	ASP	-	EXPRESSION TAG	UNP Q9X112
B	-1	PRO	-	EXPRESSION TAG	UNP Q9X112
B	0	PHE	-	EXPRESSION TAG	UNP Q9X112
B	1	THR	-	EXPRESSION TAG	UNP Q9X112
B	38	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	43	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	45	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	73	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	93	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	102	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	170	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	344	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	421	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	468	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	510	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	529	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	537	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	619	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	682	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	721	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	728	MSE	MET	MODIFIED RESIDUE	UNP Q9X112
B	733	ALA	-	EXPRESSION TAG	UNP Q9X112

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).

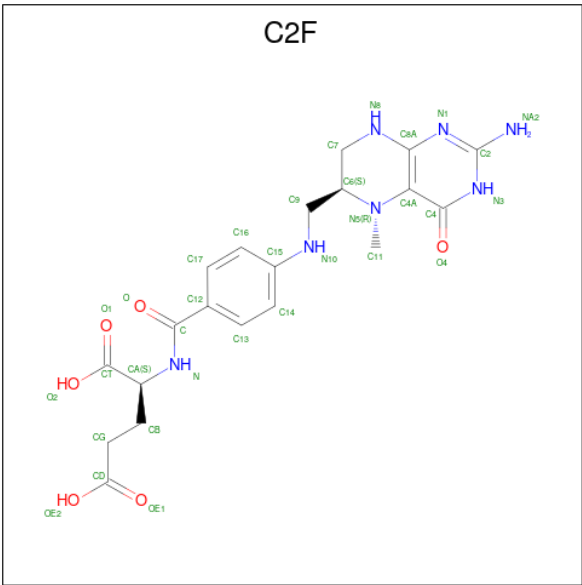


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

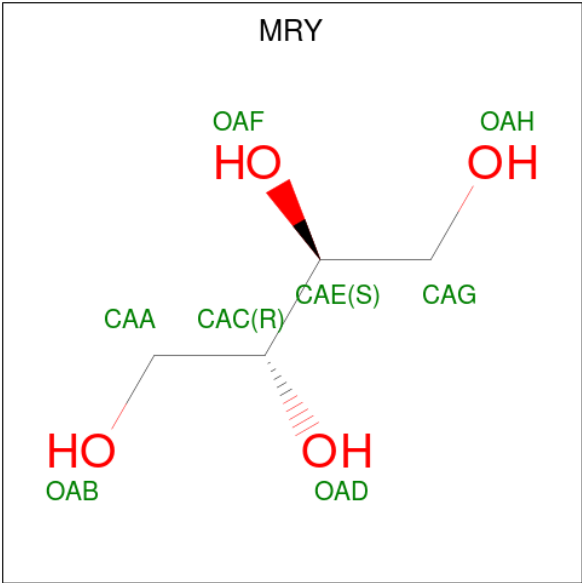
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is 5-METHYL-5,6,7,8-TETRAHYDROFOLIC ACID (three-letter code: C2F) (formula: C<sub>20</sub>H<sub>25</sub>N<sub>7</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			33	20	7	6		
4	B	1	Total	C	N	O	0	0
			33	20	7	6		

- Molecule 5 is MESO-ERYTHRITOL (three-letter code: MRY) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			8	4	4		
5	B	1	Total	C	O	0	0
			8	4	4		

- Molecule 6 is water.

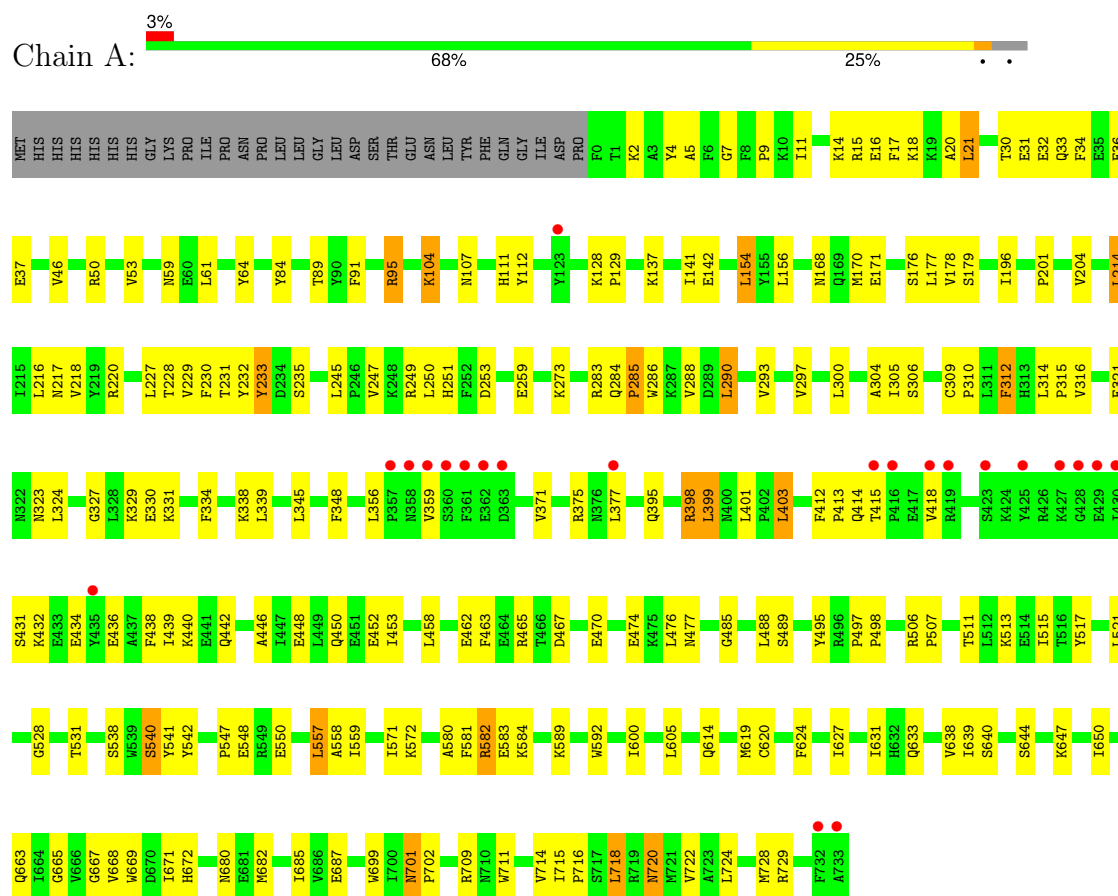
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	133	Total 133	O 133	0	0
6	B	129	Total 129	O 129	0	0



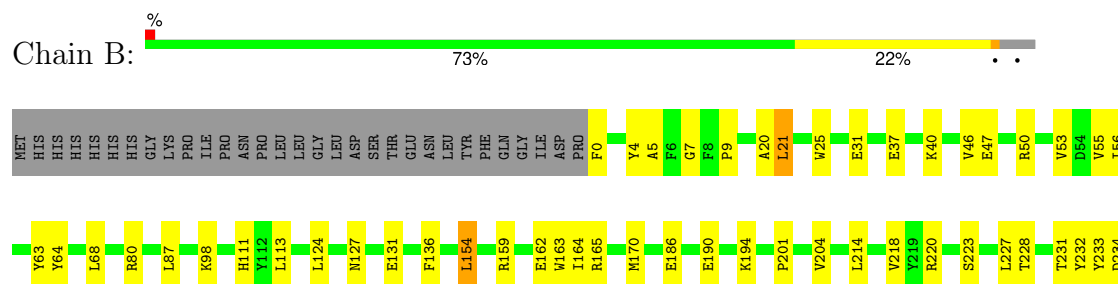
### 3 Residue-property plots

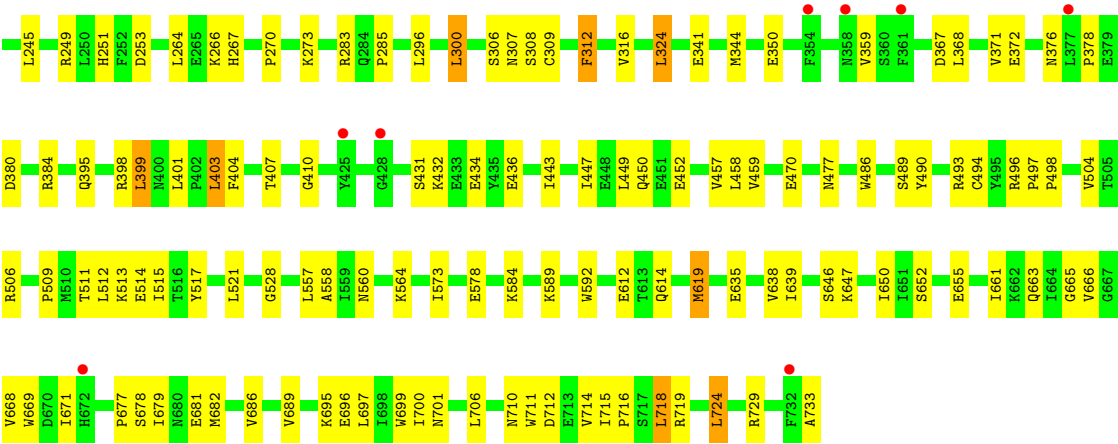
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 5-methyltetrahydropteroyltriglutamate--homocysteine methyltransferase



- Molecule 1: 5-methyltetrahydropteroyltriglutamate--homocysteine methyltransferase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	163.95Å 159.30Å 64.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 2.59 24.85 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.1 (19.99-2.59) 99.2 (24.85-2.59)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.15 (at 2.60Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.196 , 0.248 0.186 , 0.239	Depositor DCC
$R_{free}$ test set	2667 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.9	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 39.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11993	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MRY, C2F, SO4

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	0.34	0/5935	0.57	1/8032 (0.0%)
1	B	0.34	0/5957	0.58	1/8062 (0.0%)
All	All	0.34	0/11892	0.57	2/16094 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	528	GLY	N-CA-C	-5.87	98.42	113.10
1	A	528	GLY	N-CA-C	-5.44	99.51	113.10

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	5805	0	5603	132	0
1	B	5827	0	5617	111	0
2	A	10	0	0	0	0
2	B	5	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	33	0	23	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	33	0	23	4	0
5	A	8	0	10	0	0
5	B	8	0	10	0	0
6	A	133	0	0	3	0
6	B	129	0	0	3	0
All	All	11993	0	11286	245	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 11.

All (245) 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:582:ARG:HG3	1:A:627:ILE:HD13	1.61	0.81
1:A:668:VAL:HG23	1:A:682:MSE:HE2	1.65	0.77
1:A:715:ILE:HB	1:A:716:PRO:HD3	1.69	0.75
1:A:665:GLY:HA2	1:A:699:TRP:HB2	1.71	0.71
1:B:403:LEU:HD11	1:B:695:LYS:HD2	1.73	0.69
1:B:450:GLN:HB3	1:B:458:LEU:HD22	1.75	0.67
1:A:620:CYS:HB2	6:A:939:HOH:O	1.95	0.66
1:B:638:VAL:HG22	1:B:663:GLN:HB2	1.78	0.66
1:A:283:ARG:HG3	1:A:309:CYS:SG	2.36	0.65
1:A:498:PRO:HD2	1:A:540:SER:HB3	1.77	0.65
1:A:467:ASP:HB3	1:A:470:GLU:HB2	1.78	0.65
1:A:232:TYR:O	1:A:233:TYR:HB2	1.98	0.64
1:A:669:TRP:HB2	1:A:682:MSE:HE3	1.79	0.64
1:B:4:TYR:HB2	1:B:306:SER:HB3	1.78	0.64
1:A:4:TYR:HB2	1:A:306:SER:HB3	1.79	0.63
1:A:377:LEU:HD11	1:A:557:LEU:HD21	1.79	0.63
1:B:665:GLY:HA2	1:B:699:TRP:HB2	1.80	0.63
1:B:399:LEU:HB3	1:B:401:LEU:HG	1.80	0.62
1:A:196:ILE:HB	1:A:227:LEU:HD23	1.82	0.62
1:A:228:THR:HG23	1:A:249:ARG:HG3	1.81	0.62
1:B:719:ARG:HG3	1:B:719:ARG:HH11	1.64	0.62
1:A:371:VAL:HG12	1:A:375:ARG:HD2	1.82	0.62
1:A:284:GLN:NE2	1:A:542:TYR:H	1.98	0.61
4:A:1882:C2F:O4	4:A:1882:C2F:H112	2.00	0.61
1:B:56:ILE:N	1:B:56:ILE:HD12	2.15	0.61
1:B:232:TYR:O	1:B:233:TYR:HB2	2.01	0.61
1:B:431:SER:OG	1:B:434:GLU:HB2	2.01	0.60
4:B:1883:C2F:O4	4:B:1883:C2F:H112	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:PRO:O	1:A:204:VAL:HG22	2.02	0.60
1:A:399:LEU:HB3	1:A:401:LEU:HG	1.83	0.60
1:B:432:LYS:O	1:B:436:GLU:HG3	2.02	0.59
1:B:159:ARG:HH21	1:B:159:ARG:HG3	1.67	0.59
1:A:714:VAL:HG12	1:A:718:LEU:HD22	1.85	0.58
1:A:571:ILE:HD12	1:A:571:ILE:N	2.19	0.58
1:B:270:PRO:HB2	1:B:273:LYS:HG2	1.86	0.58
1:B:679:ILE:HD12	1:B:679:ILE:N	2.20	0.57
1:A:672:HIS:C	1:A:709:ARG:HH21	2.08	0.56
1:B:560:ASN:OD1	1:B:564:LYS:HE2	2.04	0.56
1:A:171:GLU:HG3	1:A:214:LEU:HD11	1.86	0.56
1:A:624:PHE:HD2	1:A:631:ILE:HD13	1.71	0.56
1:A:638:VAL:HG22	1:A:663:GLN:HB2	1.88	0.56
1:B:324:LEU:HD13	6:B:753:HOH:O	2.05	0.56
1:A:154:LEU:HB3	1:A:170:MSE:HE3	1.88	0.55
1:B:403:LEU:HD23	1:B:404:PHE:H	1.71	0.55
1:A:462:GLU:HG2	1:A:465:ARG:HD2	1.87	0.55
1:A:230:PHE:HA	1:A:251:HIS:HB3	1.88	0.55
1:A:724:LEU:O	1:A:728:MSE:HG2	2.07	0.55
1:A:32:GLU:O	1:A:36:GLU:HG3	2.08	0.54
1:A:305:ILE:HD12	1:A:345:LEU:HD23	1.89	0.54
1:A:314:LEU:HD11	1:A:338:LYS:HE3	1.89	0.54
1:A:229:VAL:HB	1:A:250:LEU:HD23	1.89	0.54
1:A:107:ASN:ND2	1:A:283:ARG:HH12	2.07	0.53
1:A:403:LEU:O	1:A:729:ARG:HD3	2.09	0.53
1:B:341:GLU:HA	1:B:344:MSE:HE3	1.90	0.53
1:B:372:GLU:HG3	1:B:376:ASN:ND2	2.23	0.53
1:A:20:ALA:HB2	1:A:37:GLU:HG2	1.89	0.53
1:B:378:PRO:C	1:B:380:ASP:H	2.11	0.53
1:B:679:ILE:HD12	1:B:679:ILE:H	1.74	0.52
1:A:415:THR:O	1:A:418:VAL:HG12	2.08	0.52
1:A:216:LEU:HG	1:A:220:ARG:NH1	2.24	0.52
1:B:449:LEU:HD21	1:B:715:ILE:CD1	2.39	0.52
1:B:506:ARG:HB2	1:B:558:ALA:HB1	1.91	0.52
1:B:714:VAL:HG12	1:B:718:LEU:HD22	1.92	0.52
1:A:176:SER:O	1:A:179:SER:HB3	2.10	0.52
1:B:729:ARG:O	1:B:733:ALA:HB2	2.10	0.52
1:A:517:TYR:CE1	1:A:521:LEU:HD11	2.44	0.52
1:A:9:PRO:HB2	1:A:316:VAL:HG22	1.92	0.52
1:B:46:VAL:HG21	1:B:136:PHE:CE2	2.44	0.51
1:A:667:GLY:HA2	1:A:701:ASN:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:589:LYS:HA	1:A:592:TRP:CD1	2.46	0.51
1:A:671:ILE:HD12	1:A:671:ILE:C	2.31	0.51
1:B:395:GLN:OE1	1:B:614:GLN:HG3	2.10	0.51
1:A:154:LEU:HD13	1:A:177:LEU:HD22	1.93	0.51
1:A:232:TYR:O	1:A:233:TYR:CB	2.58	0.51
1:B:220:ARG:O	1:B:223:SER:HB2	2.11	0.51
1:B:164:ILE:HG22	1:B:170:MSE:HG2	1.93	0.51
1:B:68:LEU:HD22	1:B:113:LEU:CD1	2.40	0.50
1:A:30:THR:OG1	1:A:33:GLN:HG3	2.11	0.50
1:B:666:VAL:HG21	1:B:689:VAL:HG11	1.92	0.50
1:A:5:ALA:HB2	1:A:53:VAL:HG11	1.93	0.50
1:A:178:VAL:CG2	1:A:218:VAL:HG13	2.41	0.50
1:A:557:LEU:HD22	1:A:605:LEU:HD21	1.93	0.50
1:A:619:MSE:HG3	1:A:639:ILE:HD11	1.92	0.50
1:A:720:ASN:N	1:A:720:ASN:HD22	2.09	0.50
1:B:283:ARG:HG2	1:B:309:CYS:SG	2.51	0.50
1:A:701:ASN:HB2	1:A:702:PRO:HD2	1.93	0.50
1:A:728:MSE:HE2	1:A:728:MSE:HA	1.94	0.50
1:B:443:ILE:O	1:B:447:ILE:HG12	2.12	0.50
1:B:668:VAL:HG23	1:B:682:MSE:HE2	1.94	0.50
1:A:168:ASN:HB3	1:B:589:LYS:O	2.11	0.49
1:A:371:VAL:CG1	1:A:375:ARG:HD2	2.43	0.49
1:A:413:PRO:HG2	1:A:711:TRP:CE2	2.47	0.49
1:B:9:PRO:HB2	1:B:316:VAL:HG22	1.94	0.49
1:A:436:GLU:OE2	1:A:513:LYS:NZ	2.44	0.49
1:A:414:GLN:HE21	1:A:463:PHE:HA	1.77	0.49
1:A:644:SER:HA	1:A:685:ILE:HD13	1.94	0.49
1:B:47:GLU:HG2	1:B:50:ARG:NH2	2.27	0.49
1:B:296:LEU:O	1:B:300:LEU:HB2	2.12	0.49
1:A:624:PHE:HD2	1:A:631:ILE:CD1	2.26	0.49
1:B:686:VAL:HG21	1:B:724:LEU:HD11	1.93	0.49
1:A:247:VAL:O	1:A:273:LYS:HE3	2.13	0.49
1:A:284:GLN:HE22	1:A:541:TYR:HA	1.78	0.49
1:A:582:ARG:CG	1:A:627:ILE:HG21	2.43	0.49
1:B:509:PRO:HB3	1:B:512:LEU:HD23	1.94	0.49
1:A:446:ALA:O	1:A:450:GLN:HG3	2.12	0.49
1:A:450:GLN:HB3	1:A:458:LEU:HD22	1.94	0.49
1:B:5:ALA:HB2	1:B:53:VAL:HG11	1.95	0.48
1:B:21:LEU:HD12	1:B:25:TRP:CZ2	2.48	0.48
1:B:403:LEU:HD22	1:B:404:PHE:CD2	2.48	0.48
1:A:231:THR:H	1:A:251:HIS:HD2	1.60	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:470:GLU:O	1:A:474:GLU:HG3	2.13	0.48
1:A:669:TRP:CE3	1:A:702:PRO:HB3	2.48	0.48
1:B:63:TYR:HD2	1:B:87:LEU:HD21	1.78	0.48
1:A:59:ASN:HB2	6:A:781:HOH:O	2.12	0.48
1:A:293:VAL:O	1:A:297:VAL:HG23	2.13	0.48
1:B:619:MSE:HE2	1:B:619:MSE:HA	1.96	0.48
1:B:652:SER:HA	1:B:655:GLU:HG2	1.95	0.48
1:B:584:LYS:O	1:B:584:LYS:HG3	2.13	0.48
1:A:511:THR:O	1:A:515:ILE:HG12	2.13	0.47
1:B:477:ASN:O	1:B:504:VAL:HA	2.14	0.47
1:A:284:GLN:HG3	1:A:285:PRO:HD2	1.96	0.47
1:A:315:PRO:HB3	1:A:485:GLY:HA2	1.97	0.47
1:B:127:ASN:O	1:B:131:GLU:HG3	2.14	0.47
1:B:700:ILE:HD13	1:B:724:LEU:CD1	2.44	0.47
1:B:589:LYS:HA	1:B:592:TRP:CD1	2.49	0.47
1:B:266:LYS:HZ3	1:B:267:HIS:CE1	2.33	0.47
1:B:494:CYS:O	4:B:1883:C2F:H17	2.15	0.47
1:B:233:TYR:O	1:B:234:ASP:HB2	2.15	0.47
1:B:619:MSE:HE3	6:B:850:HOH:O	2.14	0.47
1:B:68:LEU:HD22	1:B:113:LEU:HD12	1.97	0.47
1:B:711:TRP:O	1:B:715:ILE:HG12	2.15	0.47
1:B:163:TRP:HB3	1:B:165:ARG:HG3	1.97	0.46
1:A:284:GLN:HE21	1:A:541:TYR:HB3	1.80	0.46
1:A:547:PRO:HG2	1:A:550:GLU:HB2	1.97	0.46
1:B:715:ILE:HB	1:B:716:PRO:CD	2.44	0.46
1:A:506:ARG:HB2	1:A:558:ALA:HB1	1.97	0.46
1:B:56:ILE:N	1:B:56:ILE:CD1	2.78	0.46
1:B:450:GLN:CB	1:B:458:LEU:HD22	2.43	0.46
1:B:696:GLU:CD	1:B:696:GLU:H	2.18	0.46
1:B:228:THR:HG23	1:B:249:ARG:CB	2.46	0.46
1:A:498:PRO:CD	1:A:540:SER:HB3	2.44	0.46
1:B:449:LEU:HD11	1:B:715:ILE:HD12	1.96	0.46
1:B:511:THR:O	1:B:515:ILE:HG12	2.15	0.46
1:A:284:GLN:HG3	1:A:286:TRP:CD1	2.51	0.46
1:B:677:PRO:HB2	1:B:682:MSE:HE3	1.98	0.46
1:A:21:LEU:HD12	1:A:34:PHE:CE1	2.51	0.45
1:B:228:THR:HG23	1:B:249:ARG:HB3	1.98	0.45
1:B:403:LEU:O	1:B:729:ARG:HD2	2.16	0.45
1:A:309:CYS:HB2	1:A:310:PRO:CD	2.46	0.45
1:A:395:GLN:OE1	1:A:614:GLN:HG3	2.16	0.45
1:A:497:PRO:HA	1:A:498:PRO:HD3	1.78	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:HIS:HB2	1:B:490:TYR:HB3	1.97	0.45
1:A:137:LYS:HE3	1:A:142:GLU:OE2	2.16	0.45
1:A:640:SER:HB2	6:A:833:HOH:O	2.15	0.45
1:B:5:ALA:HA	1:B:307:ASN:OD1	2.16	0.45
1:B:232:TYR:O	1:B:233:TYR:CB	2.64	0.45
1:B:80:ARG:HD2	1:B:98:LYS:HB2	1.99	0.45
1:B:513:LYS:HG3	1:B:514:GLU:OE2	2.16	0.45
1:A:489:SER:OG	4:A:1882:C2F:H92	2.16	0.45
1:A:95:ARG:CB	1:A:95:ARG:HH21	2.29	0.45
1:A:647:LYS:HA	1:A:685:ILE:HG12	1.99	0.45
1:A:232:TYR:HA	1:A:253:ASP:HB2	1.97	0.45
1:A:531:THR:O	1:A:559:ILE:HD13	2.15	0.45
1:B:639:ILE:O	1:B:639:ILE:HG23	2.16	0.45
1:B:679:ILE:H	1:B:679:ILE:CD1	2.30	0.45
1:A:538:SER:HB2	1:A:584:LYS:HG3	1.98	0.45
1:B:646:SER:O	1:B:647:LYS:HB2	2.17	0.45
1:B:31:GLU:HG3	1:B:87:LEU:HD12	1.99	0.45
1:B:671:ILE:C	1:B:671:ILE:HD12	2.37	0.45
1:B:186:GLU:O	1:B:190:GLU:HG3	2.17	0.44
1:A:84:TYR:HA	1:A:89:THR:HG21	1.99	0.44
1:A:584:LYS:O	1:A:584:LYS:HD3	2.17	0.44
1:A:315:PRO:HB2	1:A:334:PHE:CD2	2.52	0.44
1:A:288:VAL:HG23	1:A:290:LEU:CD1	2.47	0.44
1:B:20:ALA:HB2	1:B:37:GLU:HG2	2.00	0.44
1:B:154:LEU:HD12	1:B:154:LEU:HA	1.84	0.44
1:A:104:LYS:NZ	1:A:583:GLU:OE2	2.50	0.43
1:A:178:VAL:HG22	1:A:218:VAL:HG13	2.00	0.43
1:A:644:SER:HA	1:A:685:ILE:CD1	2.48	0.43
1:A:91:PHE:O	1:A:95:ARG:HG2	2.18	0.43
1:B:367:ASP:O	1:B:371:VAL:HG23	2.18	0.43
1:A:327:GLY:O	1:A:331:LYS:HD3	2.18	0.43
1:A:477:ASN:ND2	1:A:507:PRO:HD3	2.33	0.43
1:A:548:GLU:OE2	1:A:584:LYS:HE3	2.19	0.43
1:B:384:ARG:HG3	1:B:635:GLU:OE2	2.17	0.43
1:A:11:ILE:HA	1:A:17:PHE:HB3	2.01	0.43
1:A:321:GLU:O	1:A:329:LYS:HD2	2.18	0.43
1:A:438:PHE:O	1:A:442:GLN:HG2	2.18	0.43
1:A:669:TRP:HB2	1:A:682:MSE:CE	2.46	0.43
1:B:650:ILE:HG12	1:B:650:ILE:O	2.18	0.43
1:A:432:LYS:O	1:A:436:GLU:HG3	2.19	0.43
1:A:371:VAL:O	1:A:375:ARG:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:413:PRO:HB3	1:A:709:ARG:O	2.19	0.43
1:B:159:ARG:HG3	1:B:159:ARG:NH2	2.34	0.43
1:B:497:PRO:HA	1:B:498:PRO:HD3	1.96	0.43
1:A:7:GLY:O	1:A:312:PHE:N	2.52	0.42
1:B:264:LEU:C	1:B:264:LEU:HD23	2.39	0.42
1:A:582:ARG:HG3	1:A:627:ILE:HG21	2.00	0.42
1:A:398:ARG:NH1	1:A:572:LYS:HG2	2.34	0.42
1:A:650:ILE:O	1:A:650:ILE:HG12	2.20	0.42
1:B:368:LEU:C	1:B:368:LEU:HD13	2.39	0.42
1:B:489:SER:OG	4:B:1883:C2F:H92	2.20	0.42
1:A:128:LYS:N	1:A:129:PRO:CD	2.83	0.42
1:B:661:ILE:O	1:B:697:LEU:HD21	2.20	0.42
1:B:719:ARG:HG3	1:B:719:ARG:NH1	2.31	0.42
1:B:407:THR:HB	1:B:457:VAL:HB	2.02	0.42
1:A:14:LYS:HB2	1:A:16:GLU:OE1	2.19	0.42
1:A:330:GLU:OE1	1:A:330:GLU:N	2.53	0.42
1:B:459:VAL:HB	6:B:922:HOH:O	2.19	0.42
1:B:253:ASP:OD2	1:B:308:SER:HB2	2.20	0.42
1:B:578:GLU:O	1:B:619:MSE:HE1	2.18	0.42
1:B:678:SER:OG	1:B:681:GLU:HG3	2.19	0.42
1:A:112:TYR:CE2	1:A:156:LEU:HB3	2.55	0.41
1:A:235:SER:CB	1:A:259:GLU:HB2	2.50	0.41
1:B:378:PRO:C	1:B:380:ASP:N	2.73	0.41
1:B:449:LEU:HD21	1:B:715:ILE:HD11	2.01	0.41
1:A:348:PHE:CG	1:A:356:LEU:HD11	2.54	0.41
1:B:517:TYR:CE2	1:B:521:LEU:HD11	2.55	0.41
1:A:439:ILE:HG22	1:A:440:LYS:N	2.36	0.41
1:A:448:GLU:O	1:A:452:GLU:HG3	2.20	0.41
1:A:467:ASP:OD2	1:A:495:TYR:HE2	2.03	0.41
1:B:486:TRP:CE3	1:B:496:ARG:HG3	2.55	0.41
1:B:231:THR:H	1:B:251:HIS:HD2	1.68	0.41
1:B:669:TRP:HB2	1:B:682:MSE:CE	2.51	0.41
1:B:0:PHE:HA	1:B:350:GLU:HG2	2.02	0.41
1:A:46:VAL:HG11	1:A:141:ILE:HD13	2.02	0.41
1:A:600:ILE:HD13	1:A:633:GLN:HB2	2.03	0.41
1:A:104:LYS:HD2	1:A:111:HIS:CD2	2.56	0.41
1:A:515:ILE:HG21	1:A:571:ILE:HG12	2.02	0.40
1:A:571:ILE:N	1:A:571:ILE:CD1	2.84	0.40
1:B:4:TYR:CE1	1:B:55:VAL:HG21	2.57	0.40
1:A:18:LYS:NZ	4:A:1882:C2F:O2	2.41	0.40
1:A:580:ALA:O	1:A:581:PHE:C	2.60	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:PRO:O	1:B:204:VAL:HG22	2.20	0.40
1:A:107:ASN:CG	1:A:283:ARG:HH12	2.25	0.40
1:A:453:ILE:HG22	1:A:722:VAL:HG21	2.02	0.40
1:B:7:GLY:O	1:B:312:PHE:N	2.54	0.40
1:B:154:LEU:HB3	1:B:170:MSE:HE3	2.04	0.40
1:B:573:ILE:HG12	1:B:614:GLN:HB3	2.03	0.40
1:B:710:ASN:HB3	1:B:712:ASP:OD1	2.22	0.40
4:B:1883:C2F:O4	4:B:1883:C2F:C11	2.69	0.40
1:A:2:LYS:O	1:A:304:ALA:HA	2.21	0.40
1:B:37:GLU:HA	1:B:40:LYS:HD3	2.03	0.40
1:B:214:LEU:O	1:B:218:VAL:HG23	2.21	0.40
1:A:431:SER:O	1:A:434:GLU:N	2.54	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	A	732/765 (96%)	695 (95%)	34 (5%)	3 (0%)	34	57
1	B	732/765 (96%)	693 (95%)	36 (5%)	3 (0%)	34	57
All	All	1464/1530 (96%)	1388 (95%)	70 (5%)	6 (0%)	34	57

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	64	TYR
1	A	64	TYR
1	A	359	VAL
1	A	233	TYR
1	B	359	VAL
1	B	410	GLY

### 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	597/668 (89%)	565 (95%)	32 (5%)	22	44
1	B	604/668 (90%)	580 (96%)	24 (4%)	31	57
All	All	1201/1336 (90%)	1145 (95%)	56 (5%)	26	50

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ARG
1	A	21	LEU
1	A	31	GLU
1	A	50	ARG
1	A	61	LEU
1	A	95	ARG
1	A	104	LYS
1	A	154	LEU
1	A	214	LEU
1	A	217	ASN
1	A	245	LEU
1	A	285	PRO
1	A	290	LEU
1	A	300	LEU
1	A	312	PHE
1	A	323	ASN
1	A	324	LEU
1	A	339	LEU
1	A	398	ARG
1	A	399	LEU
1	A	403	LEU
1	A	412	PHE
1	A	476	LEU
1	A	488	LEU
1	A	540	SER
1	A	557	LEU
1	A	582	ARG

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Mol	Chain	Res	Type
1	A	680	ASN
1	A	687	GLU
1	A	701	ASN
1	A	718	LEU
1	A	720	ASN
1	B	21	LEU
1	B	124	LEU
1	B	154	LEU
1	B	162	GLU
1	B	194	LYS
1	B	227	LEU
1	B	245	LEU
1	B	285	PRO
1	B	300	LEU
1	B	312	PHE
1	B	324	LEU
1	B	398	ARG
1	B	399	LEU
1	B	403	LEU
1	B	452	GLU
1	B	470	GLU
1	B	493	ARG
1	B	557	LEU
1	B	612	GLU
1	B	619	MSE
1	B	701	ASN
1	B	706	LEU
1	B	718	LEU
1	B	724	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	251	HIS
1	A	284	GLN
1	A	414	GLN
1	A	477	ASN
1	A	720	ASN
1	B	251	HIS
1	B	376	ASN
1	B	477	ASN
1	B	720	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 for Mogul analysis.

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	$\# Z  > 2$
4	C2F	B	1883	-	33,35,35	1.36	2 (6%)	35,49,49	1.48	5 (14%)
5	MRY	A	1890	-	7,7,7	0.31	0	8,8,8	0.48	0
4	C2F	A	1882	-	33,35,35	1.35	2 (6%)	35,49,49	1.51	6 (17%)
2	SO4	A	1884	-	4,4,4	0.65	0	6,6,6	0.06	0
5	MRY	B	1889	-	7,7,7	0.32	0	8,8,8	0.52	0
2	SO4	B	1885	-	4,4,4	0.64	0	6,6,6	0.08	0
2	SO4	A	1886	-	4,4,4	0.72	0	6,6,6	0.07	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
5	MRY	B	1889	-	-	1/8/8/8	-
4	C2F	B	1883	-	-	6/22/35/35	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	MRY	A	1890	-	-	1/8/8/8	-
4	C2F	A	1882	-	-	4/22/35/35	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1883	C2F	O4-C4	4.95	1.36	1.24
4	A	1882	C2F	O4-C4	4.95	1.36	1.24
4	B	1883	C2F	C4A-N5	-3.09	1.34	1.41
4	A	1882	C2F	C4A-N5	-3.08	1.34	1.41

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1882	C2F	C2-N3-C4	3.82	121.27	115.96
4	B	1883	C2F	C2-N3-C4	3.75	121.18	115.96
4	B	1883	C2F	N1-C2-N3	-3.66	119.88	125.48
4	A	1882	C2F	N1-C2-N3	-3.63	119.92	125.48
4	B	1883	C2F	C8A-C4A-C4	3.22	116.93	114.44
4	A	1882	C2F	C8A-C4A-C4	3.17	116.89	114.44
4	B	1883	C2F	C2-N1-C8A	2.76	121.15	114.59
4	A	1882	C2F	C2-N1-C8A	2.71	121.03	114.59
4	A	1882	C2F	C11-N5-C4A	2.64	116.88	113.30
4	B	1883	C2F	C11-N5-C4A	2.38	116.54	113.30
4	A	1882	C2F	NA2-C2-N3	2.05	120.29	117.22

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1882	C2F	N5-C6-C9-N10
4	A	1882	C2F	C7-C6-C9-N10
4	B	1883	C2F	C7-C6-C9-N10
4	B	1883	C2F	CB-CA-CT-O1
4	B	1883	C2F	CB-CA-CT-O2
4	B	1883	C2F	N5-C6-C9-N10
4	B	1883	C2F	OE2-CD-CG-CB
4	B	1883	C2F	OE1-CD-CG-CB
4	A	1882	C2F	OE2-CD-CG-CB
4	A	1882	C2F	OE1-CD-CG-CB
5	A	1890	MRY	OAD-CAC-CAE-CAG

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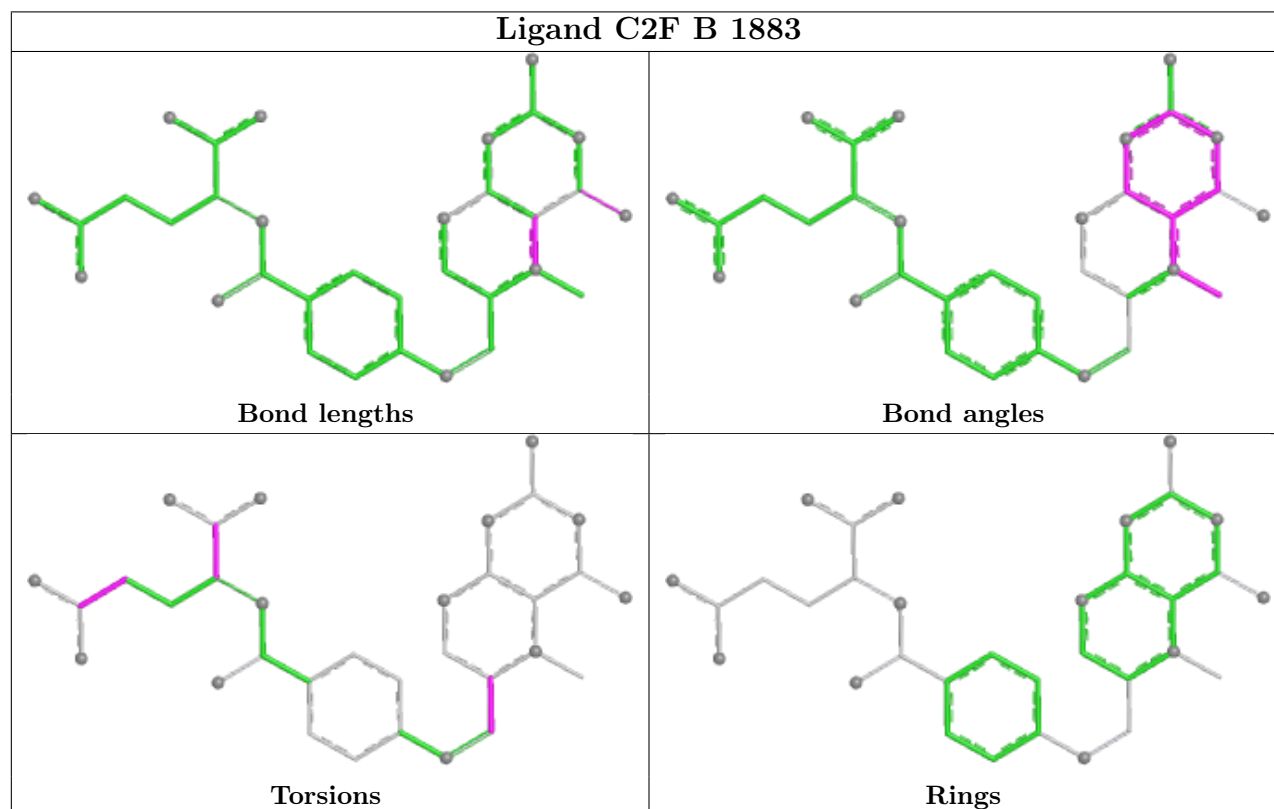
Mol	Chain	Res	Type	Atoms
5	B	1889	MRY	OAD-CAC-CAE-CAG

There are no ring outliers.

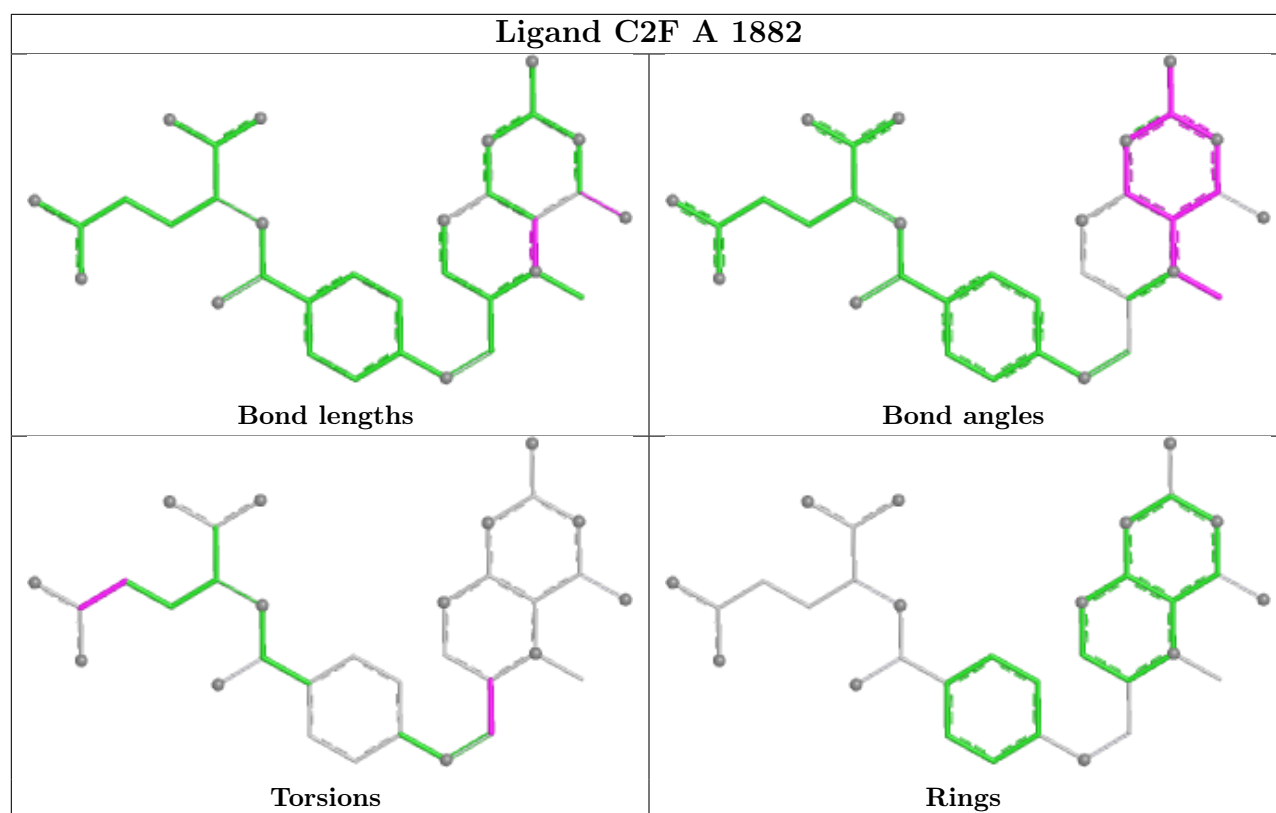
2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1883	C2F	4	0
4	A	1882	C2F	3	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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	717/765 (93%)	-0.12	22 (3%) 49 42	11, 26, 55, 77	0
1	B	717/765 (93%)	-0.21	8 (1%) 80 78	10, 24, 44, 61	0
All	All	1434/1530 (93%)	-0.17	30 (2%) 63 58	10, 25, 50, 77	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	361	PHE	3.9
1	A	428	GLY	3.8
1	A	360	SER	3.7
1	A	423	SER	3.7
1	A	359	VAL	3.5
1	B	358	ASN	3.4
1	A	377	LEU	3.4
1	A	363	ASP	3.2
1	A	732	PHE	3.2
1	A	362	GLU	3.2
1	B	354	PHE	3.1
1	B	361	PHE	3.1
1	B	428	GLY	2.9
1	A	425	TYR	2.9
1	B	732	PHE	2.9
1	B	425	TYR	2.5
1	A	419	ARG	2.4
1	B	377	LEU	2.3
1	B	672	HIS	2.3
1	A	429	GLU	2.3
1	A	733	ALA	2.3
1	A	358	ASN	2.3
1	A	435	TYR	2.3
1	A	123	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	357	PRO	2.1
1	A	427	LYS	2.1
1	A	416	PRO	2.1
1	A	418	VAL	2.1
1	A	430	ILE	2.1
1	A	415	THR	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

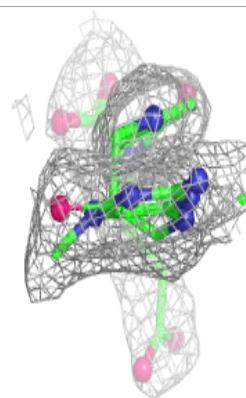
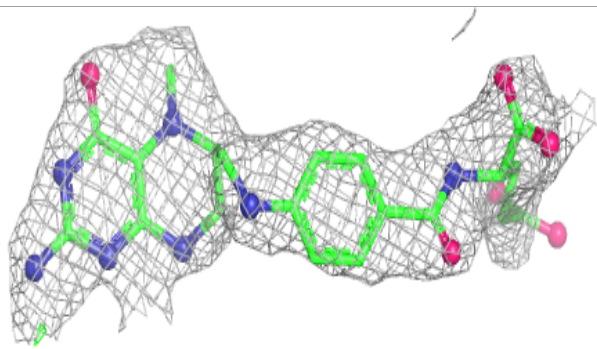
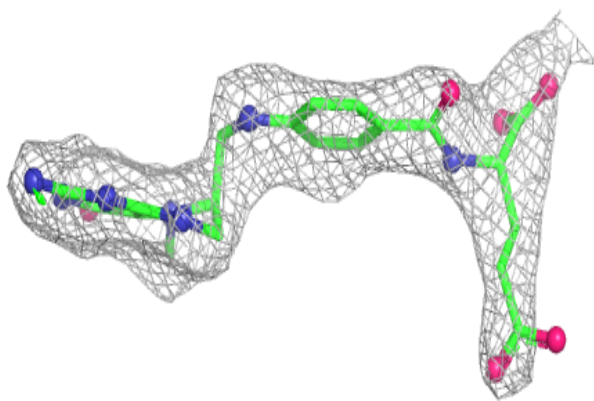
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
4	C2F	A	1882	33/33	0.87	0.24	36,40,55,57	0
2	SO4	A	1886	5/5	0.88	0.21	68,69,70,70	0
5	MRY	B	1889	8/8	0.90	0.20	31,41,44,45	0
4	C2F	B	1883	33/33	0.91	0.22	27,31,51,53	0
5	MRY	A	1890	8/8	0.93	0.23	38,43,45,46	0
2	SO4	A	1884	5/5	0.99	0.10	27,28,30,31	0
2	SO4	B	1885	5/5	0.99	0.09	25,26,28,28	0
3	ZN	A	1887	1/1	0.99	0.02	32,32,32,32	0
3	ZN	B	1888	1/1	0.99	0.03	33,33,33,33	0

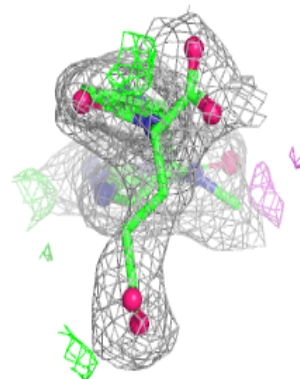
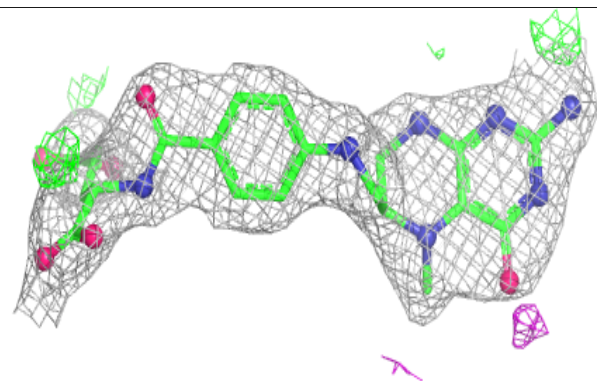
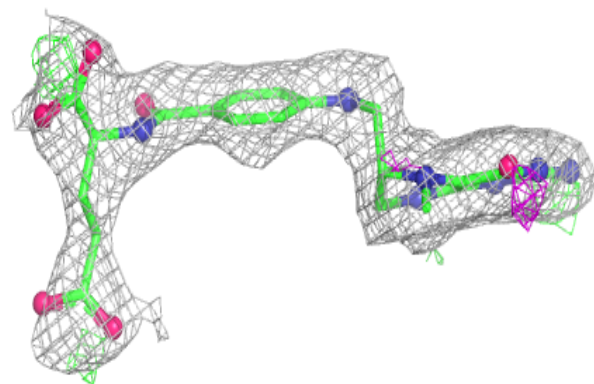
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around C2F A 1882:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around C2F B 1883:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.