



wwPDB X-ray Structure Validation Summary Report

Jun 16, 2014 – 06:41 PM BST

PDB ID : 4V7G
Title : Crystal Structure of Lumazine Synthase from Bacillus Anthracis
Authors : Morgunova, E.; Illarionov, B.; Saller, S.; Popov, A.; Sambaiah, T.; Bacher, A.;
Cushman, M.; Fischer, M.; Ladenstein, R.
Deposited on : 2009-09-16
Resolution : 3.50 Å(reported)

This is a wwPDB validation summary report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

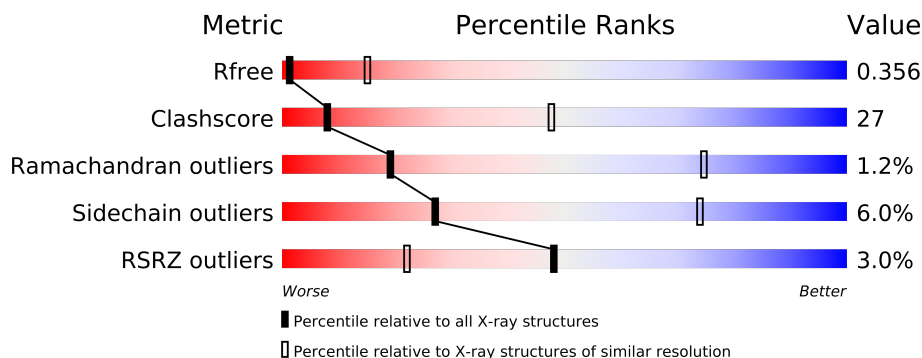
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1323
EDS : stable23397
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable23397

1 Overall quality at a glance

The reported resolution of this entry is 3.50 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}	66092	1243 (3.70-3.30)
Clashscore	79885	1039 (3.66-3.34)
Ramachandran outliers	78287	1000 (3.66-3.34)
Sidechain outliers	78261	1000 (3.66-3.34)
RSRZ outliers	66119	1243 (3.70-3.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A1	153	
1	A2	153	
1	A3	153	
1	A4	153	
1	AA	153	
1	AB	153	
1	AC	153	
1	AD	153	
1	AE	153	
1	AF	153	
1	AG	153	
1	AH	153	
1	AI	153	
1	AJ	153	

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Mol	Chain	Length	Quality of chain
1	AK	153	
1	AL	153	
1	AM	153	
1	AN	153	
1	AO	153	
1	AP	153	
1	AQ	153	
1	AR	153	
1	AS	153	
1	AT	153	
1	AU	153	
1	AV	153	
1	AW	153	
1	AX	153	
1	AY	153	
1	AZ	153	
1	B1	153	
1	B2	153	
1	B3	153	
1	B4	153	
1	BA	153	
1	BB	153	
1	BC	153	
1	BD	153	
1	BE	153	
1	BF	153	
1	BG	153	
1	BH	153	
1	BI	153	
1	BJ	153	
1	BK	153	
1	BL	153	
1	BM	153	
1	BN	153	
1	BO	153	
1	BP	153	
1	BQ	153	
1	BR	153	
1	BS	153	
1	BT	153	
1	BU	153	
1	BV	153	

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Mol	Chain	Length	Quality of chain
1	BW	153	
1	BX	153	
1	BY	153	
1	BZ	153	
1	C1	153	
1	C2	153	
1	C3	153	
1	C4	153	
1	CA	153	
1	CB	153	
1	CC	153	
1	CD	153	
1	CE	153	
1	CF	153	
1	CG	153	
1	CH	153	
1	CI	153	
1	CJ	153	
1	CK	153	
1	CL	153	
1	CM	153	
1	CN	153	
1	CO	153	
1	CP	153	
1	CQ	153	
1	CR	153	
1	CS	153	
1	CT	153	
1	CU	153	
1	CV	153	
1	CW	153	
1	CX	153	
1	CY	153	
1	CZ	153	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	PO4	AI	1001	-	X
2	PO4	AX	1001	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 102534 atoms, of which 0 are hydrogen and 0 are deuterium.

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 6,7-dimethyl-8-ribityllumazinesynthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	AB	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	AC	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AD	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AE	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AF	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	AG	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	AH	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AI	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AJ	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AK	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	AL	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	AM	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AN	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AO	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AP	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AQ	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	AR	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AS	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AT	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AU	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	AV	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	AW	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AX	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AY	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	AZ	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	A1	152	Total	C	N	O	S	0	0	0
			1129	723	193	210	3			
1	A2	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	A3	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	A4	152	Total	C	N	O	S	0	0	0
			1132	724	193	212	3			
1	BA	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BB	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BC	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BD	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BE	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BF	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BG	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	BH	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BI	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BJ	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BK	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BL	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BM	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BN	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BO	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BP	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BQ	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BR	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BS	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BT	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BU	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BV	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BW	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BX	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BY	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	BZ	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	B1	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	B2	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			

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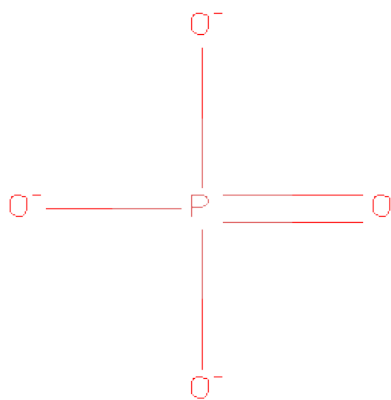
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B3	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	B4	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CA	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CB	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CC	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CD	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CE	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CF	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CG	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CH	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CI	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CJ	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CK	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CL	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CM	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CN	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CO	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CP	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CQ	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CR	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CS	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	CT	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CU	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CV	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CW	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CX	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CY	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	CZ	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	C1	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	C2	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	C3	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			
1	C4	152	Total	C	N	O	S	0	0	0
			1136	726	193	214	3			

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	AA	1	Total	O	P	0	0
			5	4	1		
2	AB	1	Total	O	P	0	0
			5	4	1		
2	AC	1	Total	O	P	0	0
			5	4	1		
2	AD	1	Total	O	P	0	0
			5	4	1		
2	AE	1	Total	O	P	0	0
			5	4	1		
2	AF	1	Total	O	P	0	0
			5	4	1		
2	AG	1	Total	O	P	0	0
			5	4	1		
2	AH	1	Total	O	P	0	0
			5	4	1		
2	AI	1	Total	O	P	0	0
			5	4	1		
2	AJ	1	Total	O	P	0	0
			5	4	1		
2	AK	1	Total	O	P	0	0
			5	4	1		
2	AL	1	Total	O	P	0	0
			5	4	1		
2	AM	1	Total	O	P	0	0
			5	4	1		
2	AN	1	Total	O	P	0	0
			5	4	1		
2	AO	1	Total	O	P	0	0
			5	4	1		
2	AP	1	Total	O	P	0	0
			5	4	1		
2	AQ	1	Total	O	P	0	0
			5	4	1		
2	AR	1	Total	O	P	0	0
			5	4	1		
2	AS	1	Total	O	P	0	0
			5	4	1		
2	AT	1	Total	O	P	0	0
			5	4	1		
2	AU	1	Total	O	P	0	0
			5	4	1		
2	AV	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	AW	1	Total	O	P	0	0
			5	4	1		
2	AX	1	Total	O	P	0	0
			5	4	1		
2	AY	1	Total	O	P	0	0
			5	4	1		
2	AZ	1	Total	O	P	0	0
			5	4	1		
2	A1	1	Total	O	P	0	0
			5	4	1		
2	A2	1	Total	O	P	0	0
			5	4	1		
2	A3	1	Total	O	P	0	0
			5	4	1		
2	A4	1	Total	O	P	0	0
			5	4	1		
2	BA	1	Total	O	P	0	0
			5	4	1		
2	BB	1	Total	O	P	0	0
			5	4	1		
2	BC	1	Total	O	P	0	0
			5	4	1		
2	BD	1	Total	O	P	0	0
			5	4	1		
2	BE	1	Total	O	P	0	0
			5	4	1		
2	BF	1	Total	O	P	0	0
			5	4	1		
2	BG	1	Total	O	P	0	0
			5	4	1		
2	BH	1	Total	O	P	0	0
			5	4	1		
2	BI	1	Total	O	P	0	0
			5	4	1		
2	BJ	1	Total	O	P	0	0
			5	4	1		
2	BK	1	Total	O	P	0	0
			5	4	1		
2	BL	1	Total	O	P	0	0
			5	4	1		
2	BM	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	BN	1	Total	O	P	0	0
			5	4	1		
2	BO	1	Total	O	P	0	0
			5	4	1		
2	BP	1	Total	O	P	0	0
			5	4	1		
2	BQ	1	Total	O	P	0	0
			5	4	1		
2	BR	1	Total	O	P	0	0
			5	4	1		
2	BS	1	Total	O	P	0	0
			5	4	1		
2	BT	1	Total	O	P	0	0
			5	4	1		
2	BU	1	Total	O	P	0	0
			5	4	1		
2	BV	1	Total	O	P	0	0
			5	4	1		
2	BW	1	Total	O	P	0	0
			5	4	1		
2	BX	1	Total	O	P	0	0
			5	4	1		
2	BY	1	Total	O	P	0	0
			5	4	1		
2	BZ	1	Total	O	P	0	0
			5	4	1		
2	B1	1	Total	O	P	0	0
			5	4	1		
2	B2	1	Total	O	P	0	0
			5	4	1		
2	B3	1	Total	O	P	0	0
			5	4	1		
2	B4	1	Total	O	P	0	0
			5	4	1		
2	CA	1	Total	O	P	0	0
			5	4	1		
2	CB	1	Total	O	P	0	0
			5	4	1		
2	CC	1	Total	O	P	0	0
			5	4	1		
2	CD	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	CE	1	Total	O	P	0	0
			5	4	1		
2	CF	1	Total	O	P	0	0
			5	4	1		
2	CG	1	Total	O	P	0	0
			5	4	1		
2	CH	1	Total	O	P	0	0
			5	4	1		
2	CI	1	Total	O	P	0	0
			5	4	1		
2	CJ	1	Total	O	P	0	0
			5	4	1		
2	CK	1	Total	O	P	0	0
			5	4	1		
2	CL	1	Total	O	P	0	0
			5	4	1		
2	CM	1	Total	O	P	0	0
			5	4	1		
2	CN	1	Total	O	P	0	0
			5	4	1		
2	CO	1	Total	O	P	0	0
			5	4	1		
2	CP	1	Total	O	P	0	0
			5	4	1		
2	CQ	1	Total	O	P	0	0
			5	4	1		
2	CR	1	Total	O	P	0	0
			5	4	1		
2	CS	1	Total	O	P	0	0
			5	4	1		
2	CT	1	Total	O	P	0	0
			5	4	1		
2	CU	1	Total	O	P	0	0
			5	4	1		
2	CV	1	Total	O	P	0	0
			5	4	1		
2	CW	1	Total	O	P	0	0
			5	4	1		
2	CX	1	Total	O	P	0	0
			5	4	1		
2	CY	1	Total	O	P	0	0
			5	4	1		

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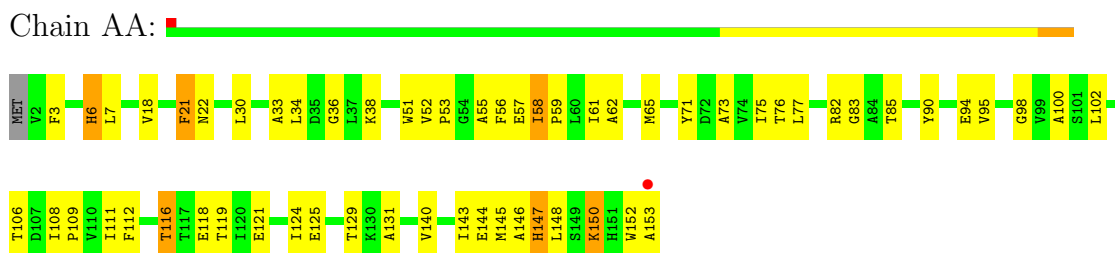
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	CZ	1	Total	O	P	0	0
			5	4	1		
2	C1	1	Total	O	P	0	0
			5	4	1		
2	C2	1	Total	O	P	0	0
			5	4	1		
2	C3	1	Total	O	P	0	0
			5	4	1		
2	C4	1	Total	O	P	0	0
			5	4	1		

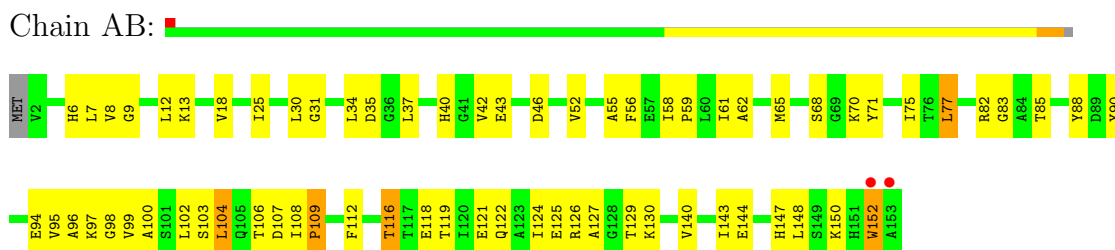
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

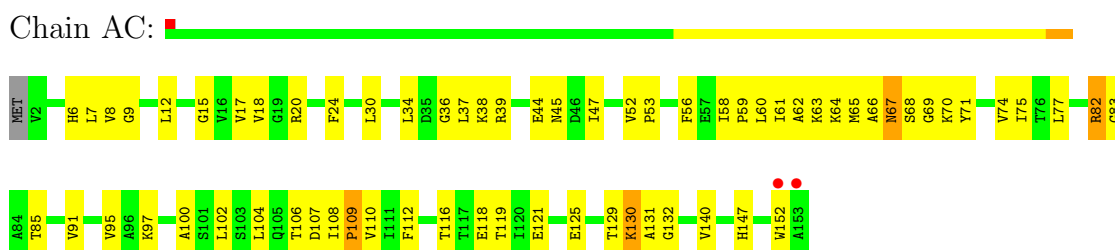
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase



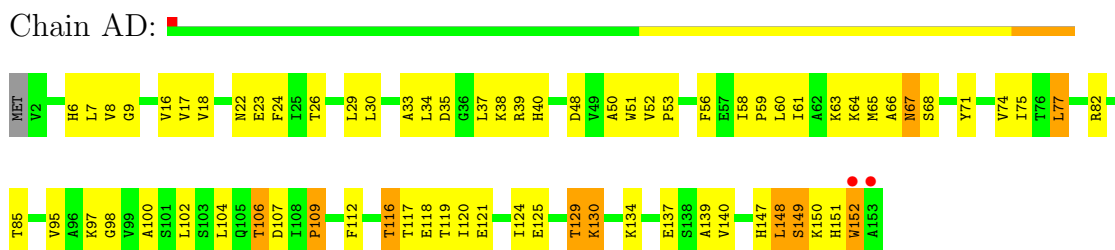
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

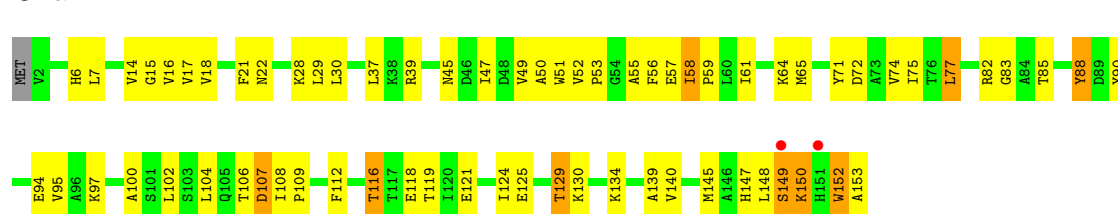


- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase



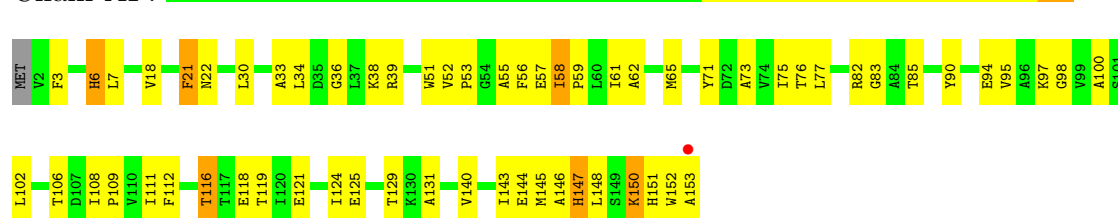
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AE:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AF:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AG:



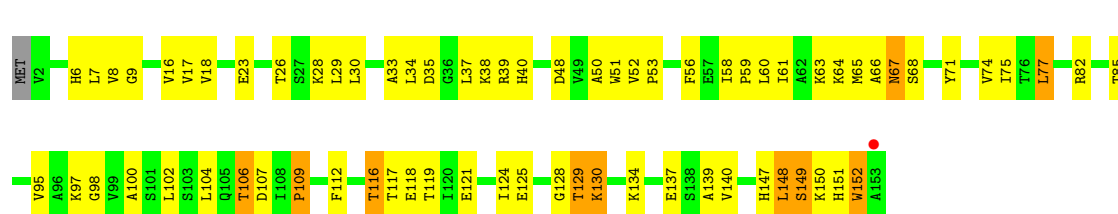
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AH:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

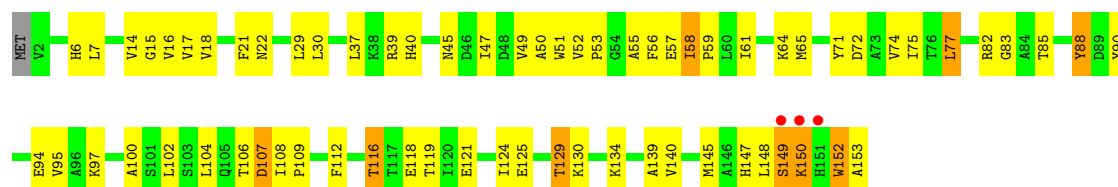
Chain AI:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

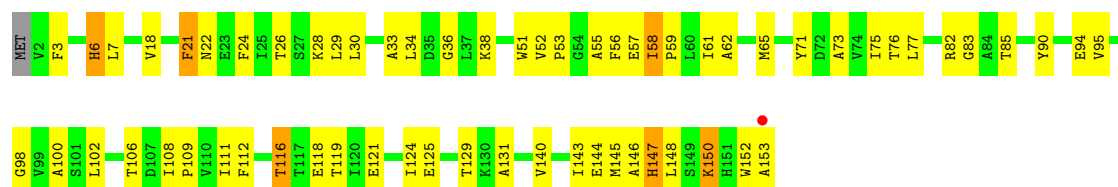
Chain AJ:





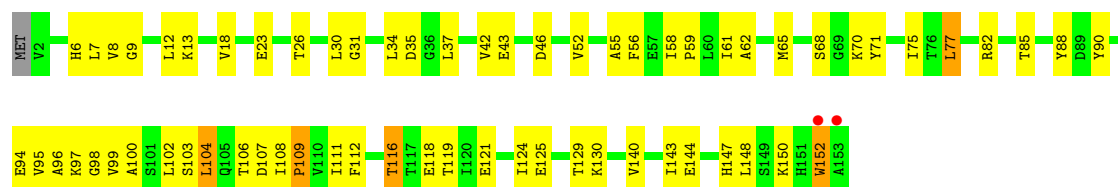
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AK:



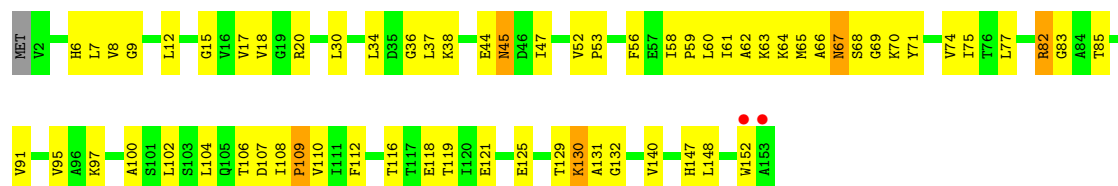
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AL:



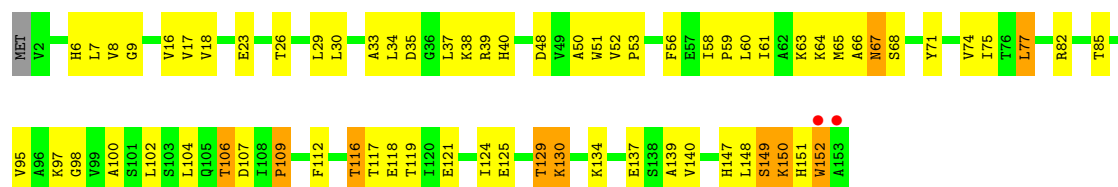
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AM:



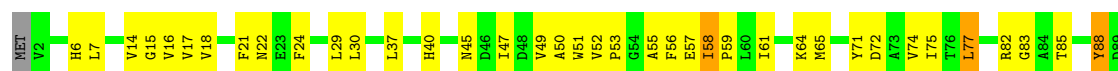
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AN:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AO:





- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AP:



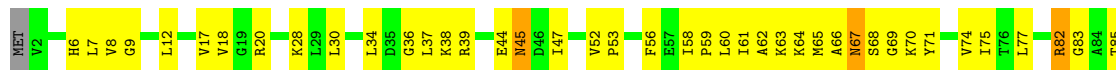
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AQ:



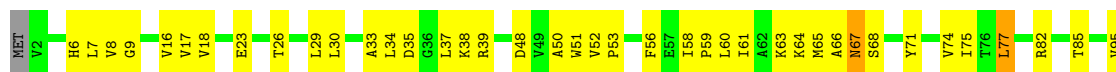
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AR:



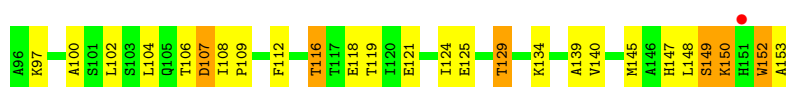
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AS:



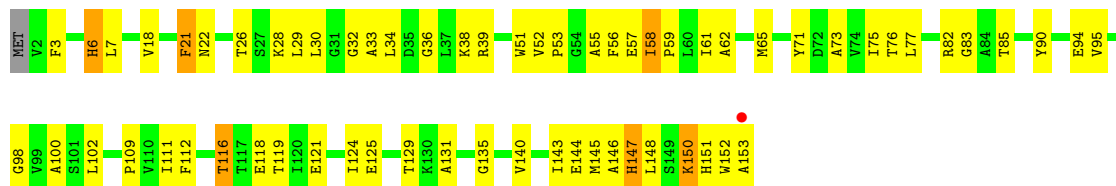
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AT:



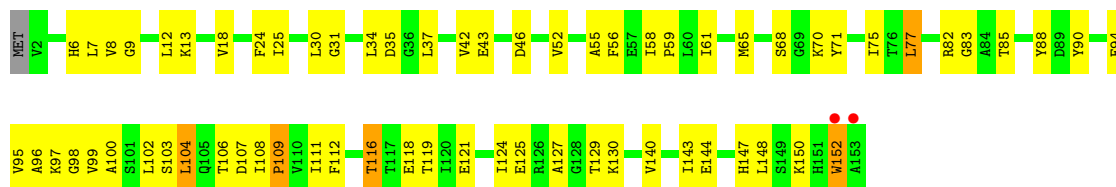
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AU:



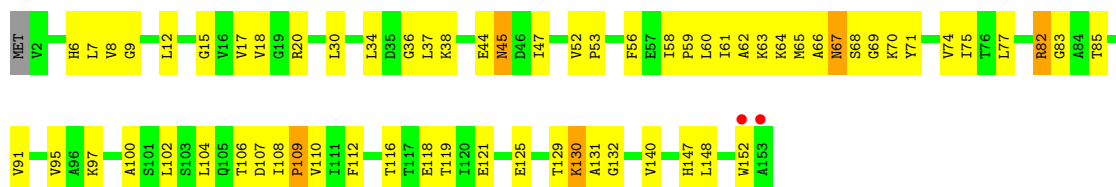
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AV:



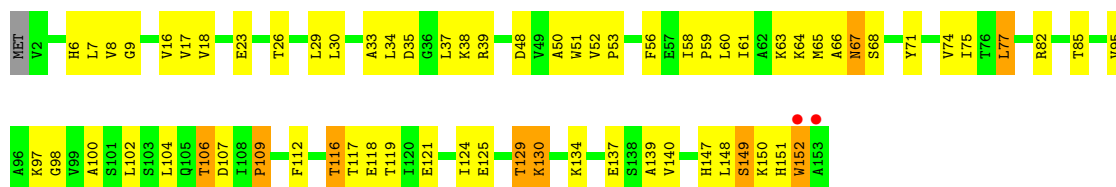
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AW:



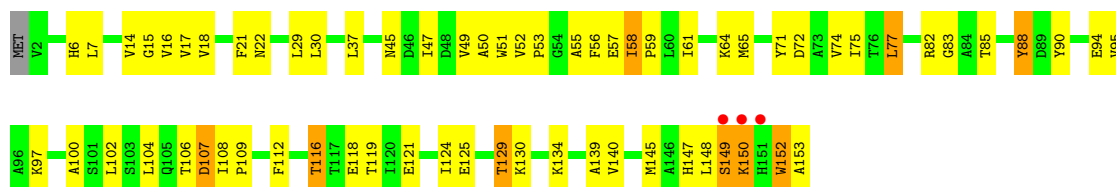
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain AX:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

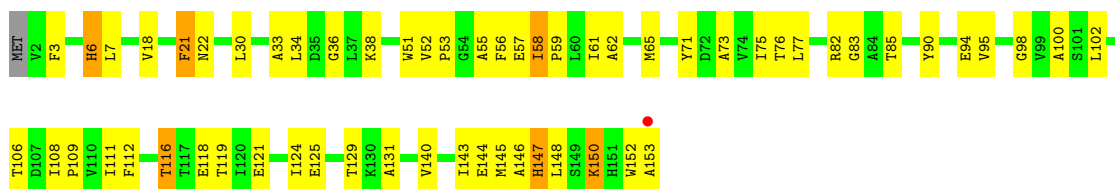
Chain AY:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

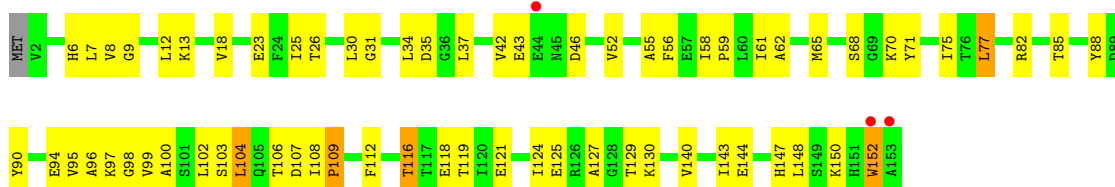
Chain AZ:





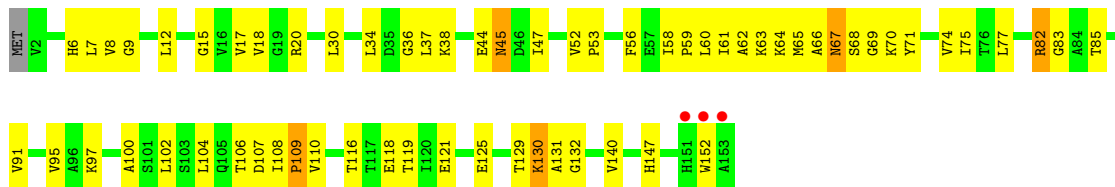
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain A1:



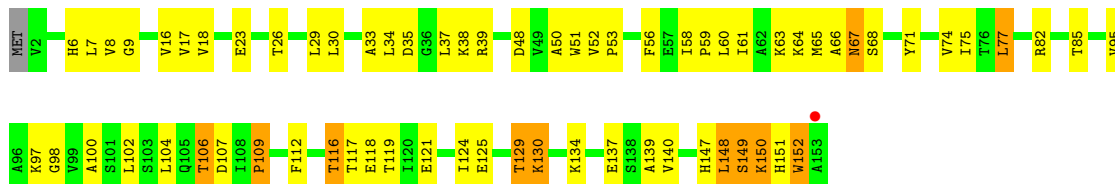
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain A2:



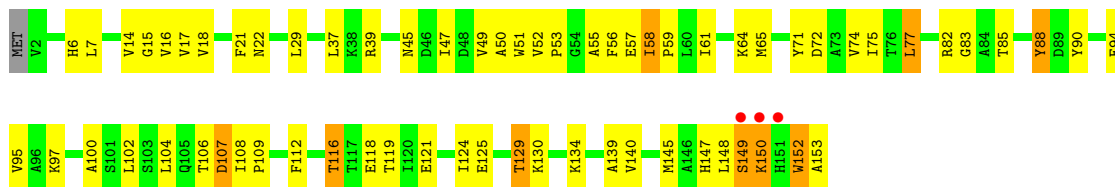
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain A3:



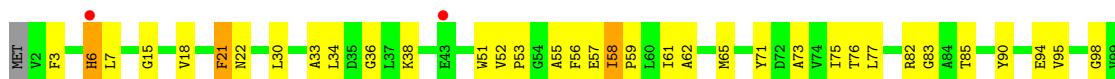
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain A4:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

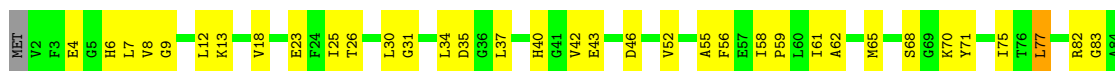
Chain BA:





- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BB:



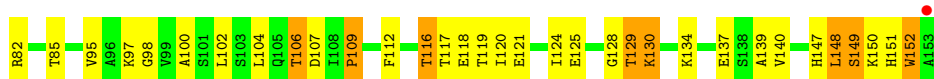
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BC:



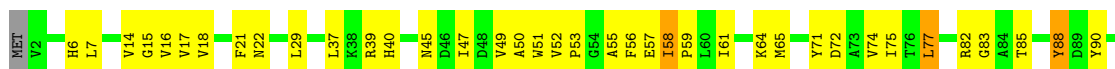
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BD:



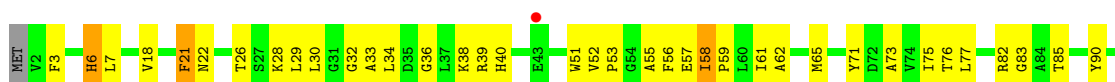
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BE:



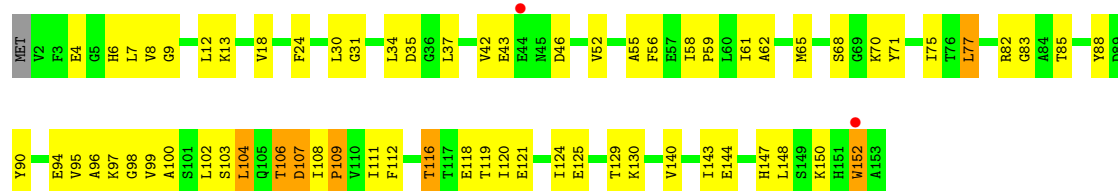
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BF:



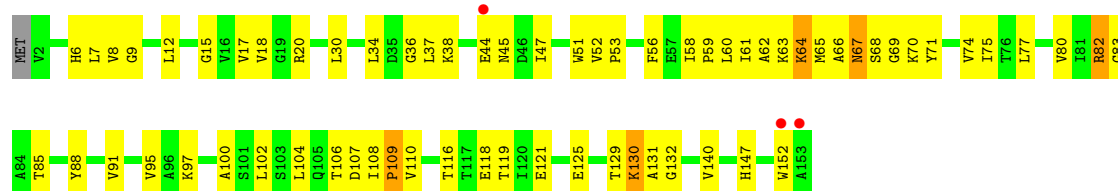
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BG:



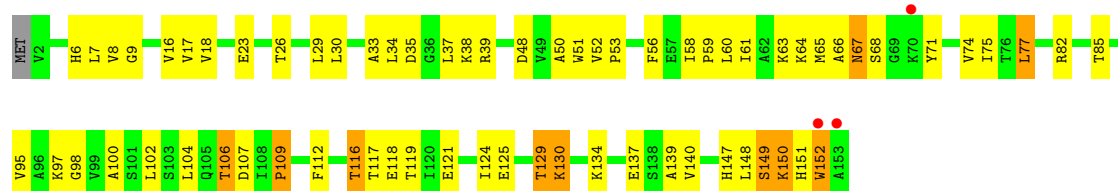
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BH:



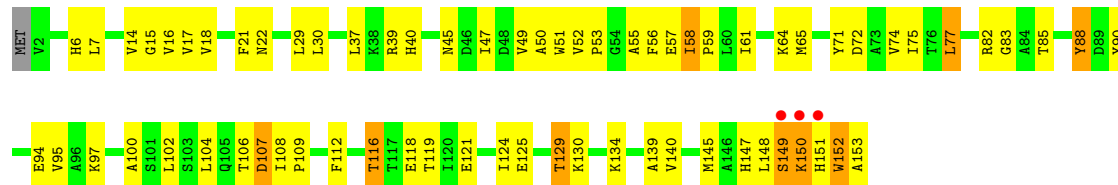
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BI:



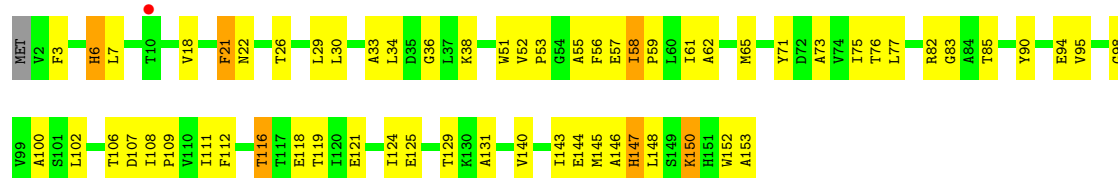
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BJ:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

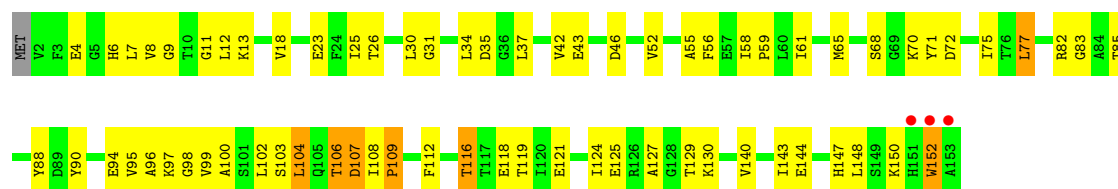
Chain BK:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

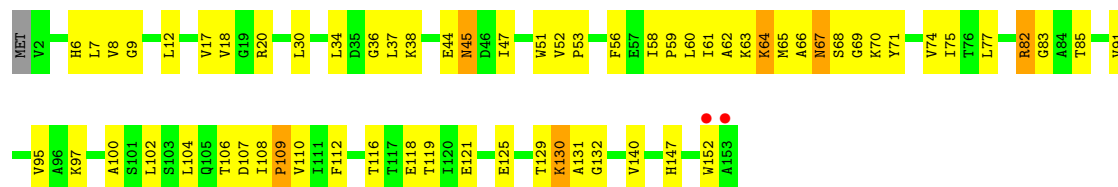
Chain BL:





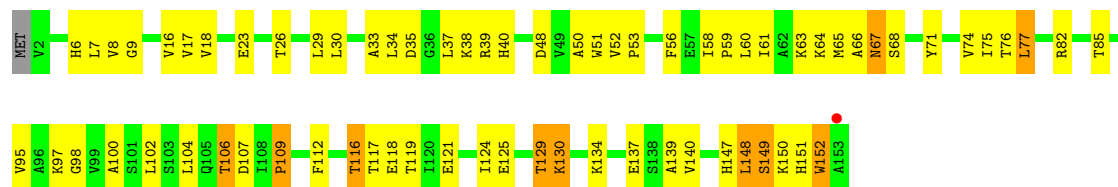
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BM:



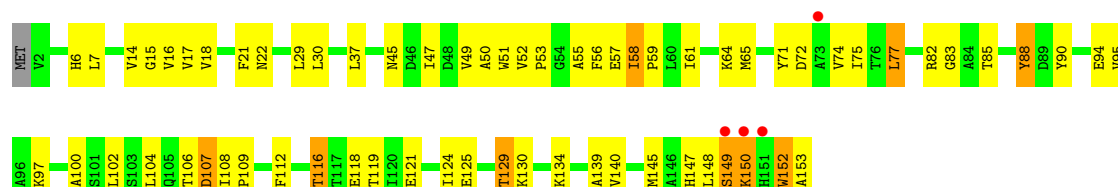
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BN:



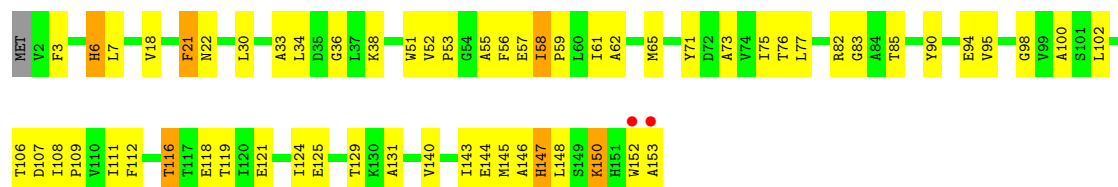
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BO:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BP:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

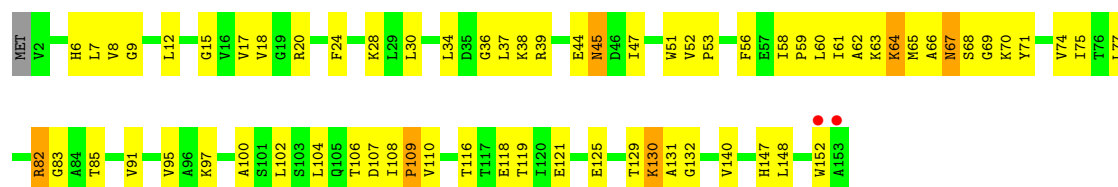
Chain BQ:





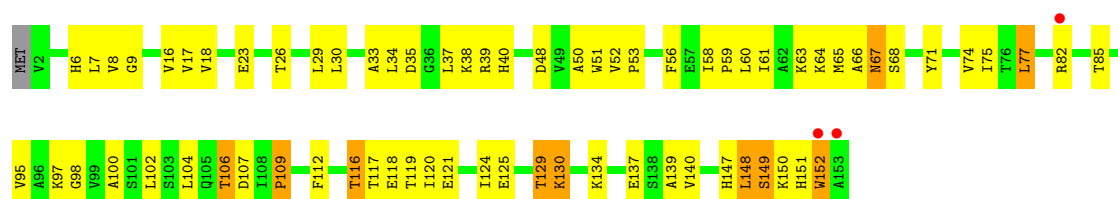
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BR:



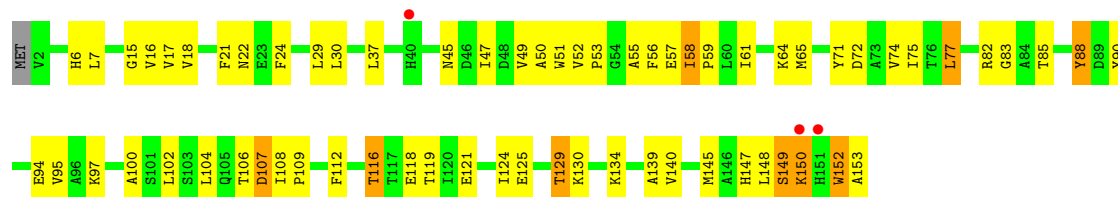
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BS:



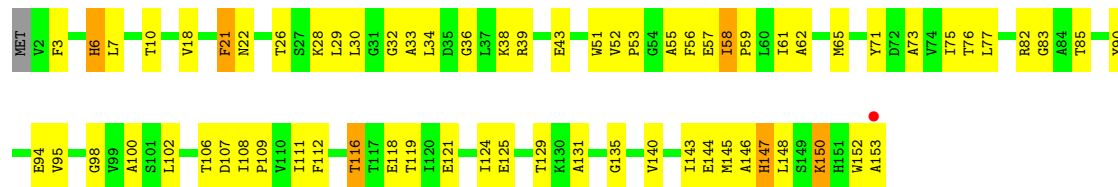
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BT:



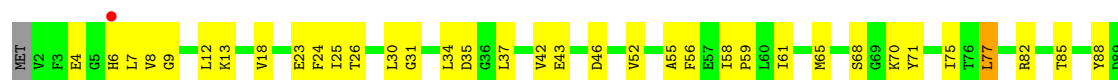
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BU:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BV:





- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BW:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BX:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BY:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain BZ:

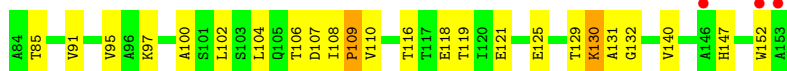
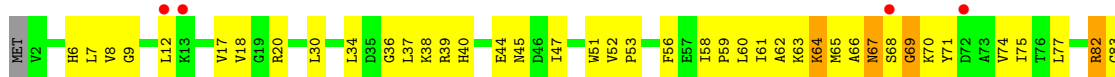


- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain B1:



Chain CC:



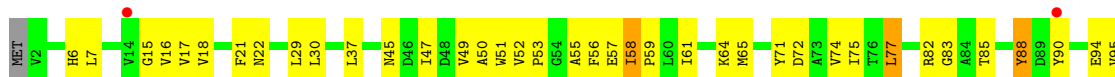
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CD:



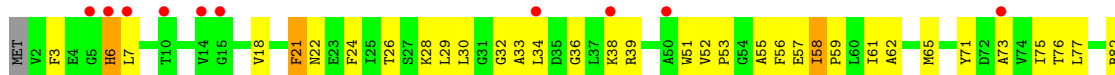
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CE:



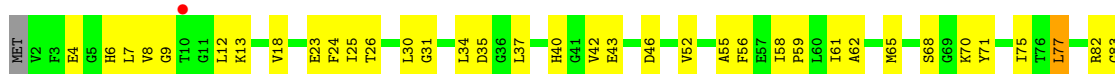
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CF:



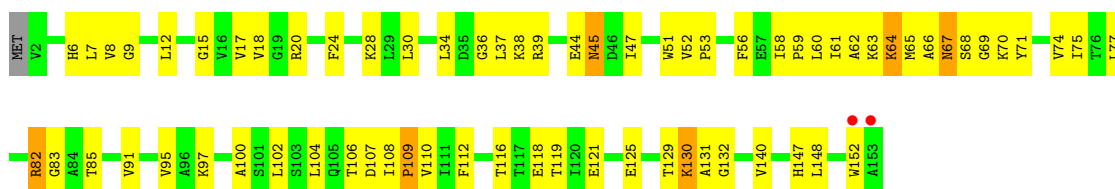
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CG:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CH:



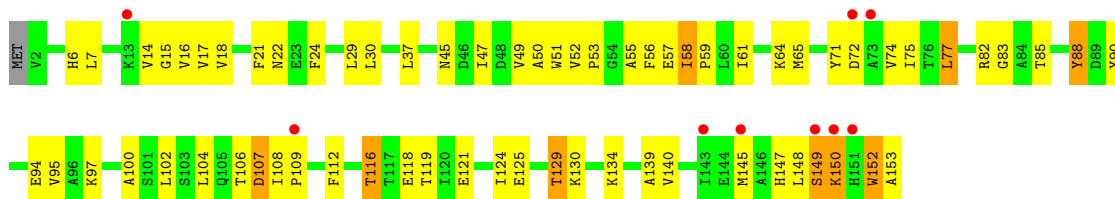
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CI:



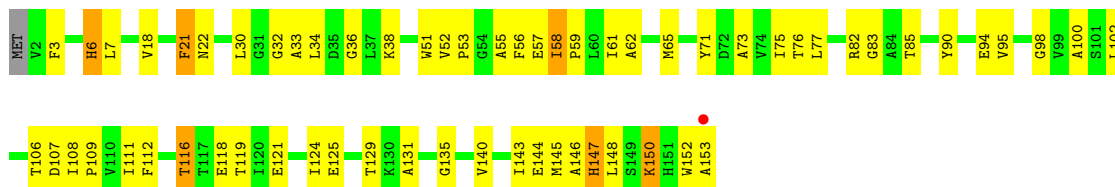
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CJ:



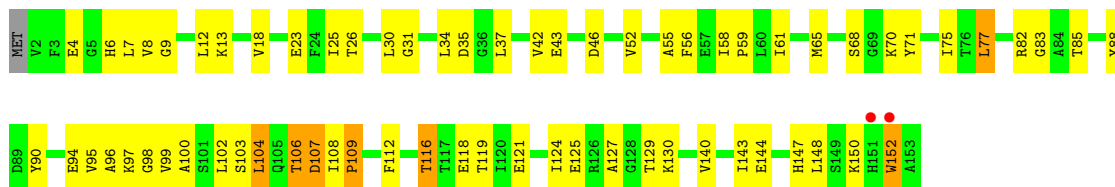
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CK:



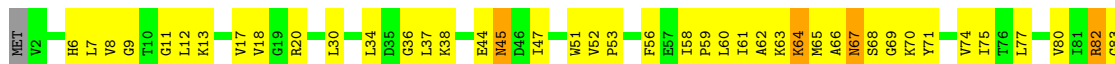
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CL:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CM:





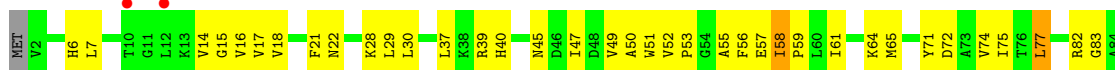
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CN:



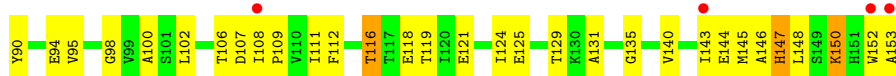
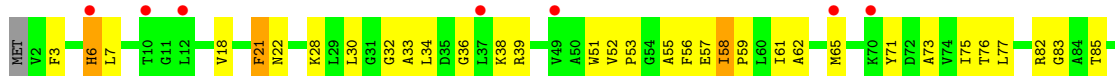
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CO:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CP:



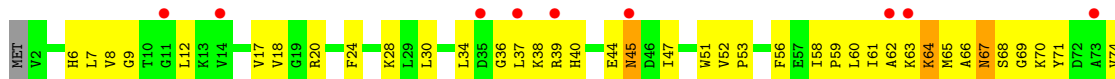
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CQ:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

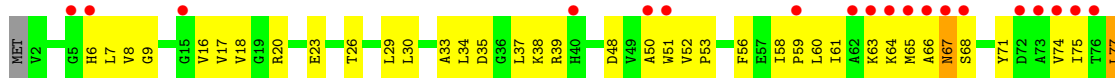
Chain CR:





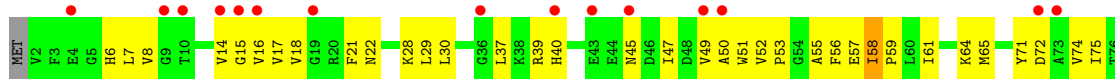
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CS:



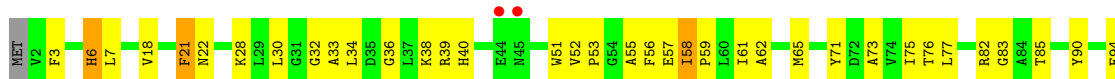
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CT:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CU:



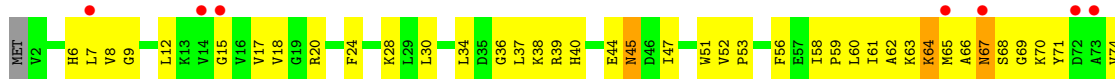
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CV:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

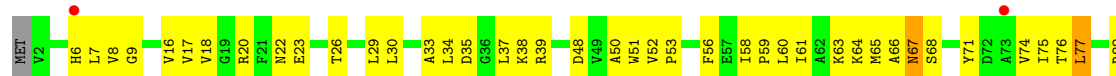
Chain CW:





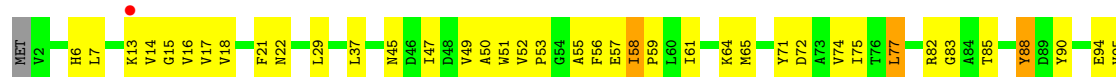
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CX:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CY:



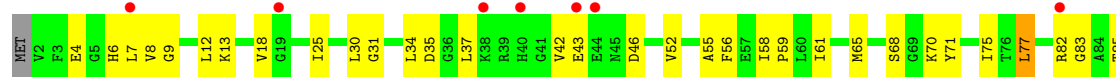
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain CZ:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain C1:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

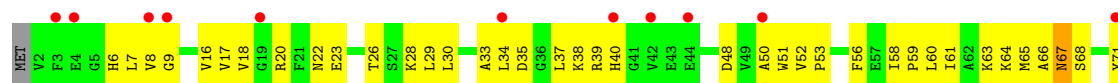
Chain C2:





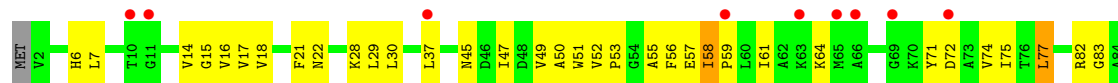
- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain C3:



- Molecule 1: 6,7-dimethyl-8-ribityllumazinesynthase

Chain C4:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	157.20Å 222.24Å 473.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 3.50 39.90 – 3.50	Depositor EDS
% Data completeness (in resolution range)	89.8 (15.00-3.50) 93.0 (39.90-3.50)	Depositor EDS
R_{merge}	0.29	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 3.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.235 , 0.286 0.353 , 0.356	Depositor DCC
R_{free} test set	3896 reflections (2.05%)	DCC
Wilson B-factor (Å ²)	67.0	Xtriage
Anisotropy	0.104	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 44.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Outliers	1 of 194062 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.72	EDS
Total number of atoms	102534	wwPDB-VP
Average B, all atoms (Å ²)	104.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	A1	0.44	0/1149	0.71	2/1558 (0.1%)
1	A2	0.32	0/1152	0.62	1/1562 (0.1%)
1	A3	0.31	0/1152	0.85	5/1562 (0.3%)
1	A4	0.34	0/1152	0.58	1/1562 (0.1%)
1	AA	0.33	0/1149	0.49	1/1558 (0.1%)
1	AB	0.44	0/1149	0.71	2/1558 (0.1%)
1	AC	0.32	0/1152	0.62	1/1562 (0.1%)
1	AD	0.31	0/1152	0.85	5/1562 (0.3%)
1	AE	0.34	0/1152	0.58	1/1562 (0.1%)
1	AF	0.33	0/1149	0.49	1/1558 (0.1%)
1	AG	0.44	0/1149	0.71	2/1558 (0.1%)
1	AH	0.32	0/1152	0.63	1/1562 (0.1%)
1	AI	0.31	0/1152	0.85	5/1562 (0.3%)
1	AJ	0.34	0/1152	0.58	1/1562 (0.1%)
1	AK	0.33	0/1149	0.49	1/1558 (0.1%)
1	AL	0.44	0/1149	0.71	2/1558 (0.1%)
1	AM	0.32	0/1152	0.63	1/1562 (0.1%)
1	AN	0.31	0/1152	0.85	5/1562 (0.3%)
1	AO	0.34	0/1152	0.58	1/1562 (0.1%)
1	AP	0.33	0/1149	0.49	1/1558 (0.1%)
1	AQ	0.44	0/1149	0.71	2/1558 (0.1%)
1	AR	0.32	0/1152	0.62	1/1562 (0.1%)
1	AS	0.31	0/1152	0.85	5/1562 (0.3%)
1	AT	0.34	0/1152	0.58	1/1562 (0.1%)
1	AU	0.33	0/1149	0.49	1/1558 (0.1%)
1	AV	0.44	0/1149	0.71	2/1558 (0.1%)
1	AW	0.32	0/1152	0.63	1/1562 (0.1%)
1	AX	0.31	0/1152	0.85	5/1562 (0.3%)
1	AY	0.34	0/1152	0.58	1/1562 (0.1%)
1	AZ	0.33	0/1149	0.49	1/1558 (0.1%)
1	B1	0.44	0/1156	0.72	3/1567 (0.2%)
1	B2	0.32	0/1156	0.62	1/1567 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	B3	0.31	0/1156	0.85	5/1567 (0.3%)
1	B4	0.34	0/1156	0.58	1/1567 (0.1%)
1	BA	0.32	0/1156	0.51	2/1567 (0.1%)
1	BB	0.44	0/1156	0.72	3/1567 (0.2%)
1	BC	0.32	0/1156	0.62	1/1567 (0.1%)
1	BD	0.31	0/1156	0.85	5/1567 (0.3%)
1	BE	0.34	0/1156	0.58	1/1567 (0.1%)
1	BF	0.32	0/1156	0.51	2/1567 (0.1%)
1	BG	0.44	0/1156	0.72	3/1567 (0.2%)
1	BH	0.32	0/1156	0.62	1/1567 (0.1%)
1	BI	0.31	0/1156	0.85	5/1567 (0.3%)
1	BJ	0.34	0/1156	0.58	1/1567 (0.1%)
1	BK	0.32	0/1156	0.51	2/1567 (0.1%)
1	BL	0.44	0/1156	0.72	3/1567 (0.2%)
1	BM	0.32	0/1156	0.62	1/1567 (0.1%)
1	BN	0.31	0/1156	0.85	5/1567 (0.3%)
1	BO	0.34	0/1156	0.58	1/1567 (0.1%)
1	BP	0.33	0/1156	0.51	2/1567 (0.1%)
1	BQ	0.44	0/1156	0.72	3/1567 (0.2%)
1	BR	0.32	0/1156	0.62	1/1567 (0.1%)
1	BS	0.31	0/1156	0.85	5/1567 (0.3%)
1	BT	0.34	0/1156	0.58	1/1567 (0.1%)
1	BU	0.33	0/1156	0.51	2/1567 (0.1%)
1	BV	0.44	0/1156	0.72	3/1567 (0.2%)
1	BW	0.32	0/1156	0.62	1/1567 (0.1%)
1	BX	0.31	0/1156	0.85	5/1567 (0.3%)
1	BY	0.34	0/1156	0.58	1/1567 (0.1%)
1	BZ	0.33	0/1156	0.51	2/1567 (0.1%)
1	C1	0.44	0/1156	0.72	3/1567 (0.2%)
1	C2	0.32	0/1156	0.62	1/1567 (0.1%)
1	C3	0.31	0/1156	0.85	5/1567 (0.3%)
1	C4	0.34	0/1156	0.58	1/1567 (0.1%)
1	CA	0.33	0/1156	0.51	2/1567 (0.1%)
1	CB	0.44	0/1156	0.72	3/1567 (0.2%)
1	CC	0.32	0/1156	0.62	1/1567 (0.1%)
1	CD	0.31	0/1156	0.85	5/1567 (0.3%)
1	CE	0.34	0/1156	0.58	1/1567 (0.1%)
1	CF	0.33	0/1156	0.51	2/1567 (0.1%)
1	CG	0.44	0/1156	0.72	3/1567 (0.2%)
1	CH	0.32	0/1156	0.62	1/1567 (0.1%)
1	CI	0.31	0/1156	0.85	5/1567 (0.3%)
1	CJ	0.34	0/1156	0.58	1/1567 (0.1%)
1	CK	0.33	0/1156	0.51	2/1567 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	CL	0.44	0/1156	0.72	3/1567 (0.2%)
1	CM	0.32	0/1156	0.62	1/1567 (0.1%)
1	CN	0.31	0/1156	0.85	5/1567 (0.3%)
1	CO	0.34	0/1156	0.58	1/1567 (0.1%)
1	CP	0.32	0/1156	0.51	2/1567 (0.1%)
1	CQ	0.44	0/1156	0.72	3/1567 (0.2%)
1	CR	0.32	0/1156	0.62	1/1567 (0.1%)
1	CS	0.31	0/1156	0.85	5/1567 (0.3%)
1	CT	0.34	0/1156	0.58	1/1567 (0.1%)
1	CU	0.33	0/1156	0.51	2/1567 (0.1%)
1	CV	0.44	0/1156	0.72	3/1567 (0.2%)
1	CW	0.32	0/1156	0.62	1/1567 (0.1%)
1	CX	0.31	0/1156	0.85	5/1567 (0.3%)
1	CY	0.34	0/1156	0.58	1/1567 (0.1%)
1	CZ	0.33	0/1156	0.51	2/1567 (0.1%)
All	All	0.35	0/103884	0.67	204/140832 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A1	0	1
1	AB	0	1
1	AG	0	1
1	AL	0	1
1	AQ	0	1
1	AV	0	1
1	B1	0	1
1	BB	0	1
1	BG	0	1
1	BL	0	1
1	BQ	0	1
1	BV	0	1
1	C1	0	1
1	CB	0	1
1	CG	0	1
1	CL	0	1
1	CQ	0	1
1	CV	0	1
All	All	0	18

There are no bond length outliers.

The worst 5 of 204 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BD	149	SER	CB-CA-C	-14.91	81.78	110.10
1	CD	149	SER	CB-CA-C	-14.90	81.79	110.10
1	C3	149	SER	CB-CA-C	-14.89	81.80	110.10
1	BI	149	SER	CB-CA-C	-14.89	81.81	110.10
1	CX	149	SER	CB-CA-C	-14.89	81.81	110.10

There are no chirality outliers.

5 of 18 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AB	99	VAL	Peptide
1	AG	99	VAL	Peptide
1	AL	99	VAL	Peptide
1	AQ	99	VAL	Peptide
1	AV	99	VAL	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A1	1129	0	1143	68	0
1	A2	1132	0	1145	77	0
1	A3	1132	0	1145	64	0
1	A4	1132	0	1145	75	0
1	AA	1129	0	1143	52	0
1	AB	1129	0	1143	73	0
1	AC	1132	0	1145	81	0
1	AD	1132	0	1145	71	0
1	AE	1132	0	1145	82	0
1	AF	1129	0	1143	57	0
1	AG	1129	0	1143	68	2
1	AH	1132	0	1145	76	3
1	AI	1132	0	1145	70	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AJ	1132	0	1145	81	0
1	AK	1129	0	1143	55	0
1	AL	1129	0	1143	68	0
1	AM	1132	0	1145	78	0
1	AN	1132	0	1145	64	0
1	AO	1132	0	1145	80	0
1	AP	1129	0	1143	53	0
1	AQ	1129	0	1143	90	0
1	AR	1132	0	1145	85	0
1	AS	1132	0	1145	65	0
1	AT	1132	0	1145	75	0
1	AU	1129	0	1143	58	0
1	AV	1129	0	1143	67	0
1	AW	1132	0	1145	78	0
1	AX	1132	0	1145	62	0
1	AY	1132	0	1145	75	0
1	AZ	1129	0	1143	50	0
1	B1	1136	0	1149	73	0
1	B2	1136	0	1149	78	0
1	B3	1136	0	1149	60	0
1	B4	1136	0	1149	73	0
1	BA	1136	0	1149	49	0
1	BB	1136	0	1149	81	0
1	BC	1136	0	1149	88	0
1	BD	1136	0	1149	75	0
1	BE	1136	0	1149	74	0
1	BF	1136	0	1149	64	0
1	BG	1136	0	1149	76	0
1	BH	1136	0	1149	80	0
1	BI	1136	0	1149	63	0
1	BJ	1136	0	1149	78	1
1	BK	1136	0	1149	51	0
1	BL	1136	0	1149	74	3
1	BM	1136	0	1149	78	0
1	BN	1136	0	1149	65	0
1	BO	1136	0	1149	75	0
1	BP	1136	0	1149	49	0
1	BQ	1136	0	1149	81	0
1	BR	1136	0	1149	90	0
1	BS	1136	0	1149	67	0
1	BT	1136	0	1149	74	0
1	BU	1136	0	1149	77	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	BV	1136	0	1149	75	0
1	BW	1136	0	1149	80	0
1	BX	1136	0	1149	63	0
1	BY	1136	0	1149	77	0
1	BZ	1136	0	1149	49	0
1	C1	1136	0	1149	74	0
1	C2	1136	0	1149	91	0
1	C3	1136	0	1149	76	0
1	C4	1136	0	1149	74	1
1	CA	1136	0	1149	55	0
1	CB	1136	0	1149	75	2
1	CC	1136	0	1149	96	2
1	CD	1136	0	1149	75	0
1	CE	1136	0	1149	76	0
1	CF	1136	0	1149	64	0
1	CG	1136	0	1149	84	3
1	CH	1136	0	1149	86	3
1	CI	1136	0	1149	68	0
1	CJ	1136	0	1149	77	0
1	CK	1136	0	1149	52	0
1	CL	1136	0	1149	74	0
1	CM	1136	0	1149	86	3
1	CN	1136	0	1149	68	0
1	CO	1136	0	1149	80	0
1	CP	1136	0	1149	59	0
1	CQ	1136	0	1149	81	0
1	CR	1136	0	1149	96	0
1	CS	1136	0	1149	71	0
1	CT	1136	0	1149	83	0
1	CU	1136	0	1149	61	0
1	CV	1136	0	1149	77	0
1	CW	1136	0	1149	92	0
1	CX	1136	0	1149	69	0
1	CY	1136	0	1149	73	1
1	CZ	1136	0	1149	52	1
2	A1	5	0	0	0	0
2	A2	5	0	0	3	0
2	A3	5	0	0	0	0
2	A4	5	0	0	2	0
2	AA	5	0	0	1	0
2	AB	5	0	0	1	0
2	AC	5	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	AD	5	0	0	0	0
2	AE	5	0	0	2	0
2	AF	5	0	0	1	0
2	AG	5	0	0	1	0
2	AH	5	0	0	3	0
2	AI	5	0	0	0	0
2	AJ	5	0	0	2	0
2	AK	5	0	0	1	0
2	AL	5	0	0	0	0
2	AM	5	0	0	3	0
2	AN	5	0	0	0	0
2	AO	5	0	0	2	0
2	AP	5	0	0	1	0
2	AQ	5	0	0	1	0
2	AR	5	0	0	3	0
2	AS	5	0	0	0	0
2	AT	5	0	0	2	0
2	AU	5	0	0	1	0
2	AV	5	0	0	1	0
2	AW	5	0	0	3	0
2	AX	5	0	0	0	0
2	AY	5	0	0	2	0
2	AZ	5	0	0	1	0
2	B1	5	0	0	1	0
2	B2	5	0	0	3	0
2	B3	5	0	0	0	0
2	B4	5	0	0	2	0
2	BA	5	0	0	1	0
2	BB	5	0	0	1	0
2	BC	5	0	0	3	0
2	BD	5	0	0	0	0
2	BE	5	0	0	2	0
2	BF	5	0	0	1	0
2	BG	5	0	0	1	0
2	BH	5	0	0	3	0
2	BI	5	0	0	0	0
2	BJ	5	0	0	2	0
2	BK	5	0	0	1	0
2	BL	5	0	0	1	0
2	BM	5	0	0	3	0
2	BN	5	0	0	0	0
2	BO	5	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	BP	5	0	0	1	0
2	BQ	5	0	0	1	0
2	BR	5	0	0	3	0
2	BS	5	0	0	0	0
2	BT	5	0	0	2	0
2	BU	5	0	0	1	0
2	BV	5	0	0	0	0
2	BW	5	0	0	3	0
2	BX	5	0	0	0	0
2	BY	5	0	0	2	0
2	BZ	5	0	0	1	0
2	C1	5	0	0	1	0
2	C2	5	0	0	3	0
2	C3	5	0	0	0	0
2	C4	5	0	0	2	0
2	CA	5	0	0	1	0
2	CB	5	0	0	1	0
2	CC	5	0	0	3	0
2	CD	5	0	0	0	0
2	CE	5	0	0	2	0
2	CF	5	0	0	1	0
2	CG	5	0	0	1	0
2	CH	5	0	0	3	0
2	CI	5	0	0	0	0
2	CJ	5	0	0	2	0
2	CK	5	0	0	1	0
2	CL	5	0	0	1	0
2	CM	5	0	0	3	0
2	CN	5	0	0	0	0
2	CO	5	0	0	2	0
2	CP	5	0	0	1	0
2	CQ	5	0	0	0	0
2	CR	5	0	0	3	0
2	CS	5	0	0	0	0
2	CT	5	0	0	2	0
2	CU	5	0	0	1	0
2	CV	5	0	0	0	0
2	CW	5	0	0	3	0
2	CX	5	0	0	0	0
2	CY	5	0	0	2	0
2	CZ	5	0	0	1	0
All	All	102534	0	103266	5521	13

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 27.

The worst 5 of 5521 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:CT:53:PRO:CG	1:CT:57:GLU:HG3	1.49	1.43
1:CJ:53:PRO:CG	1:CJ:57:GLU:HG3	1.49	1.42
1:AE:53:PRO:CG	1:AE:57:GLU:HG3	1.49	1.42
1:B4:53:PRO:CG	1:B4:57:GLU:HG3	1.49	1.42
1:C4:53:PRO:CG	1:C4:57:GLU:HG3	1.49	1.42

The worst 5 of 13 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	Distance(Å)	Clash(Å)
1:AG:153:ALA:O	1:AH:46:ASP:OD1[2_565]	1.84	0.36
1:AH:69:GLY:O	1:AH:69:GLY:O[2_565]	1.90	0.30
1:BJ:151:HIS:CE1	1:CY:13:LYS:CE[2_665]	1.91	0.29
1:CB:153:ALA:O	1:CC:69:GLY:CA[2_565]	1.94	0.26
1:CH:39:ARG:NH2	1:CM:20:ARG:NH1[2_665]	1.96	0.24

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	A1	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	A2	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	A3	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	A4	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	AA	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	AB	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	AC	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	AD	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	AE	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	AF	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	AG	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	AH	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	AI	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	AJ	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	AK	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	AL	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	AM	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	AN	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	AO	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	AP	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	AQ	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	AR	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	AS	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	AT	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	AU	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	AV	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	AW	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	AX	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	AY	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	AZ	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	B1	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	B2	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	B3	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	B4	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	BA	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	BB	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	BC	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	BD	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	BE	150/153 (98%)	133 (89%)	17 (11%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	BF	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	BG	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	BH	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	BI	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	BJ	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	BK	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	BL	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	BM	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	BN	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	BO	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	BP	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	BQ	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	BR	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	BS	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	BT	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	BU	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	BV	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	BW	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	BX	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	BY	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	BZ	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	C1	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	C2	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	C3	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	C4	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	CA	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	CB	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	CC	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	CD	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	CE	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	CF	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	CG	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	CH	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	CI	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	CJ	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	CK	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	CL	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	CM	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	CN	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	CO	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	CP	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	CQ	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	CR	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	CS	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	CT	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	CU	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
1	CV	150/153 (98%)	132 (88%)	18 (12%)	0	100	100
1	CW	150/153 (98%)	126 (84%)	19 (13%)	5 (3%)	6	51
1	CX	150/153 (98%)	137 (91%)	11 (7%)	2 (1%)	18	72
1	CY	150/153 (98%)	133 (89%)	17 (11%)	0	100	100
1	CZ	150/153 (98%)	129 (86%)	19 (13%)	2 (1%)	18	72
All	All	13500/13770 (98%)	11826 (88%)	1512 (11%)	162 (1%)	19	75

5 of 162 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	AC	130	LYS
1	AH	130	LYS
1	AM	130	LYS
1	AR	130	LYS
1	AW	130	LYS

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	A1	115/118 (98%)	105 (91%)	10 (9%)	15	57
1	A2	116/118 (98%)	112 (97%)	4 (3%)	49	88
1	A3	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	A4	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AA	115/118 (98%)	110 (96%)	5 (4%)	40	84
1	AB	115/118 (98%)	105 (91%)	10 (9%)	15	57
1	AC	116/118 (98%)	112 (97%)	4 (3%)	49	88
1	AD	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AE	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AF	115/118 (98%)	110 (96%)	5 (4%)	40	84
1	AG	115/118 (98%)	105 (91%)	10 (9%)	15	57
1	AH	116/118 (98%)	112 (97%)	4 (3%)	49	88
1	AI	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AJ	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AK	115/118 (98%)	110 (96%)	5 (4%)	40	84
1	AL	115/118 (98%)	105 (91%)	10 (9%)	15	57
1	AM	116/118 (98%)	112 (97%)	4 (3%)	49	88
1	AN	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AO	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AP	115/118 (98%)	110 (96%)	5 (4%)	40	84
1	AQ	115/118 (98%)	105 (91%)	10 (9%)	15	57
1	AR	116/118 (98%)	112 (97%)	4 (3%)	49	88
1	AS	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AT	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AU	115/118 (98%)	110 (96%)	5 (4%)	40	84
1	AV	115/118 (98%)	105 (91%)	10 (9%)	15	57
1	AW	116/118 (98%)	112 (97%)	4 (3%)	49	88
1	AX	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AY	116/118 (98%)	108 (93%)	8 (7%)	22	69
1	AZ	115/118 (98%)	110 (96%)	5 (4%)	40	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B1	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	B2	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	B3	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	B4	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BA	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	BB	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	BC	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	BD	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BE	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BF	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	BG	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	BH	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	BI	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BJ	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BK	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	BL	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	BM	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	BN	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BO	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BP	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	BQ	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	BR	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	BS	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BT	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BU	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	BV	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	BW	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	BX	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BY	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	BZ	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	C1	117/118 (99%)	107 (92%)	10 (8%)	15	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C2	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	C3	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	C4	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CA	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	CB	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	CC	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	CD	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CE	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CF	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	CG	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	CH	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	CI	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CJ	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CK	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	CL	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	CM	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	CN	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CO	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CP	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	CQ	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	CR	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	CS	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CT	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CU	117/118 (99%)	112 (96%)	5 (4%)	40	84
1	CV	117/118 (99%)	107 (92%)	10 (8%)	15	58
1	CW	117/118 (99%)	113 (97%)	4 (3%)	49	88
1	CX	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CY	117/118 (99%)	109 (93%)	8 (7%)	22	70
1	CZ	117/118 (99%)	112 (96%)	5 (4%)	40	84
All	All	10488/10620 (99%)	9858 (94%)	630 (6%)	27	74

5 of 630 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	BK	150	LYS
1	BU	116	THR
1	CX	109	PRO
1	BL	116	THR
1	BP	150	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 252 such sidechains are listed below:

Mol	Chain	Res	Type
1	BK	67	ASN
1	BU	45	ASN
1	CW	93	ASN
1	BL	93	ASN
1	BP	67	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

90 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	PO4	A1	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	A2	1001	-	4,4,4	0.40	0	6,6,6	0.28	0
2	PO4	A3	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	A4	1001	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	AA	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	AB	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	AC	1001	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	AD	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	AE	1001	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	AF	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	AG	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	AH	1001	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	AI	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	AJ	1001	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	AK	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	AL	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	AM	1001	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	AN	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	AO	1001	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	AP	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	AQ	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	AR	1001	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	AS	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	AT	1001	-	4,4,4	0.39	0	6,6,6	0.29	0
2	PO4	AU	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	AV	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	AW	201	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	AX	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	AY	1001	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	AZ	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	B1	1001	-	4,4,4	0.30	0	6,6,6	0.28	0
2	PO4	B2	1001	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	B3	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	B4	1001	-	4,4,4	0.37	0	6,6,6	0.29	0
2	PO4	BA	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	BB	1001	-	4,4,4	0.30	0	6,6,6	0.28	0
2	PO4	BC	1001	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	BD	1001	-	4,4,4	0.35	0	6,6,6	0.28	0
2	PO4	BE	1001	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	BF	1001	-	4,4,4	0.30	0	6,6,6	0.29	0
2	PO4	BG	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	BH	1001	-	4,4,4	0.40	0	6,6,6	0.28	0
2	PO4	BI	1001	-	4,4,4	0.37	0	6,6,6	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	BJ	1001	-	4,4,4	0.37	0	6,6,6	0.29	0
2	PO4	BK	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	BL	1001	-	4,4,4	0.30	0	6,6,6	0.28	0
2	PO4	BM	1001	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	BN	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	BO	1001	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	BP	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	BQ	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	BR	1001	-	4,4,4	0.40	0	6,6,6	0.28	0
2	PO4	BS	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	BT	1001	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	BU	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	BV	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	BW	1001	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	BX	1001	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	BY	1001	-	4,4,4	0.37	0	6,6,6	0.29	0
2	PO4	BZ	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	C1	1201	-	4,4,4	0.30	0	6,6,6	0.28	0
2	PO4	C2	2201	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	C3	3201	-	4,4,4	0.37	0	6,6,6	0.28	0
2	PO4	C4	4201	-	4,4,4	0.37	0	6,6,6	0.29	0
2	PO4	CA	1001	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	CB	1001	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	CC	1001	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	CD	1001	-	4,4,4	0.37	0	6,6,6	0.28	0
2	PO4	CE	1001	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	CF	201	-	4,4,4	0.28	0	6,6,6	0.29	0
2	PO4	CG	1201	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	CH	2201	-	4,4,4	0.40	0	6,6,6	0.28	0
2	PO4	CI	3201	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	CJ	4201	-	4,4,4	0.38	0	6,6,6	0.29	0
2	PO4	CK	201	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	CL	1201	-	4,4,4	0.30	0	6,6,6	0.28	0
2	PO4	CM	2201	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	CN	3201	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	CO	4201	-	4,4,4	0.37	0	6,6,6	0.29	0
2	PO4	CP	201	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	CQ	1201	-	4,4,4	0.29	0	6,6,6	0.28	0
2	PO4	CR	2201	-	4,4,4	0.39	0	6,6,6	0.28	0
2	PO4	CS	3201	-	4,4,4	0.36	0	6,6,6	0.28	0
2	PO4	CT	4201	-	4,4,4	0.39	0	6,6,6	0.29	0
2	PO4	CU	201	-	4,4,4	0.29	0	6,6,6	0.29	0
2	PO4	CV	1201	-	4,4,4	0.29	0	6,6,6	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	CW	2201	-	4,4,4	0.39	0	6,6,6	0.29	0
2	PO4	CX	3201	-	4,4,4	0.35	0	6,6,6	0.28	0
2	PO4	CY	4201	-	4,4,4	0.37	0	6,6,6	0.29	0
2	PO4	CZ	201	-	4,4,4	0.28	0	6,6,6	0.29	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
2	PO4	A1	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	A2	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	A3	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	A4	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AA	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AB	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AC	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AD	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AE	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AF	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AG	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AH	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AI	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AJ	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AK	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AL	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AM	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AN	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AO	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AP	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AQ	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AR	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AS	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AT	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AU	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AV	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AW	201	-	-	0/0/0/0	0/0/0/0
2	PO4	AX	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AY	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	AZ	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	B1	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	B2	1001	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PO4	B3	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	B4	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BA	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BB	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BC	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BD	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BE	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BF	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BG	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BH	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BI	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BJ	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BK	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BL	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BM	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BN	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BO	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BP	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BQ	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BR	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BS	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BT	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BU	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BV	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BW	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BX	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BY	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	BZ	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	C1	1201	-	-	0/0/0/0	0/0/0/0
2	PO4	C2	2201	-	-	0/0/0/0	0/0/0/0
2	PO4	C3	3201	-	-	0/0/0/0	0/0/0/0
2	PO4	C4	4201	-	-	0/0/0/0	0/0/0/0
2	PO4	CA	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	CB	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	CC	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	CD	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	CE	1001	-	-	0/0/0/0	0/0/0/0
2	PO4	CF	201	-	-	0/0/0/0	0/0/0/0
2	PO4	CG	1201	-	-	0/0/0/0	0/0/0/0
2	PO4	CH	2201	-	-	0/0/0/0	0/0/0/0
2	PO4	CI	3201	-	-	0/0/0/0	0/0/0/0
2	PO4	CJ	4201	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PO4	CK	201	-	-	0/0/0/0	0/0/0/0
2	PO4	CL	1201	-	-	0/0/0/0	0/0/0/0
2	PO4	CM	2201	-	-	0/0/0/0	0/0/0/0
2	PO4	CN	3201	-	-	0/0/0/0	0/0/0/0
2	PO4	CO	4201	-	-	0/0/0/0	0/0/0/0
2	PO4	CP	201	-	-	0/0/0/0	0/0/0/0
2	PO4	CQ	1201	-	-	0/0/0/0	0/0/0/0
2	PO4	CR	2201	-	-	0/0/0/0	0/0/0/0
2	PO4	CS	3201	-	-	0/0/0/0	0/0/0/0
2	PO4	CT	4201	-	-	0/0/0/0	0/0/0/0
2	PO4	CU	201	-	-	0/0/0/0	0/0/0/0
2	PO4	CV	1201	-	-	0/0/0/0	0/0/0/0
2	PO4	CW	2201	-	-	0/0/0/0	0/0/0/0
2	PO4	CX	3201	-	-	0/0/0/0	0/0/0/0
2	PO4	CY	4201	-	-	0/0/0/0	0/0/0/0
2	PO4	CZ	201	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

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, 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	A1	152/153 (99%)	0.25	3 (1%) 62 30	65, 95, 151, 347	0
1	A2	152/153 (99%)	0.14	3 (1%) 62 30	56, 97, 184, 226	0
1	A3	152/153 (99%)	0.13	1 (0%) 84 56	57, 96, 147, 259	0
1	A4	152/153 (99%)	0.16	3 (1%) 62 30	56, 94, 153, 393	0
1	AA	152/153 (99%)	0.09	1 (0%) 84 56	55, 97, 150, 271	0
1	AB	152/153 (99%)	0.18	2 (1%) 74 40	65, 95, 151, 347	0
1	AC	152/153 (99%)	0.15	2 (1%) 74 40	56, 97, 184, 226	0
1	AD	152/153 (99%)	0.14	2 (1%) 74 40	57, 96, 147, 259	0
1	AE	152/153 (99%)	0.10	2 (1%) 74 40	56, 94, 153, 393	0
1	AF	152/153 (99%)	0.18	1 (0%) 84 56	55, 97, 150, 271	0
1	AG	152/153 (99%)	0.30	3 (1%) 62 30	65, 95, 151, 347	0
1	AH	152/153 (99%)	0.32	2 (1%) 74 40	56, 97, 184, 226	0
1	AI	152/153 (99%)	0.21	1 (0%) 84 56	57, 96, 147, 259	0
1	AJ	152/153 (99%)	0.27	3 (1%) 62 30	56, 94, 153, 393	0
1	AK	152/153 (99%)	0.15	1 (0%) 84 56	55, 97, 150, 271	0
1	AL	152/153 (99%)	0.28	2 (1%) 74 40	65, 95, 151, 347	0
1	AM	152/153 (99%)	0.18	2 (1%) 74 40	56, 97, 184, 226	0
1	AN	152/153 (99%)	0.20	2 (1%) 74 40	57, 96, 147, 259	0
1	AO	152/153 (99%)	0.19	3 (1%) 62 30	56, 94, 153, 393	0
1	AP	152/153 (99%)	0.10	0 100 100	55, 97, 150, 271	0
1	AQ	152/153 (99%)	0.16	1 (0%) 84 56	65, 95, 151, 347	0
1	AR	152/153 (99%)	0.21	2 (1%) 74 40	56, 97, 184, 226	0
1	AS	152/153 (99%)	0.19	1 (0%) 84 56	57, 96, 147, 259	0
1	AT	152/153 (99%)	0.10	1 (0%) 84 56	56, 94, 153, 393	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	AU	152/153 (99%)	0.21	1 (0%) 84 56	55, 97, 150, 271	0
1	AV	152/153 (99%)	0.28	2 (1%) 74 40	65, 95, 151, 347	0
1	AW	152/153 (99%)	0.35	2 (1%) 74 40	56, 97, 184, 226	0
1	AX	152/153 (99%)	0.24	2 (1%) 74 40	57, 96, 147, 259	0
1	AY	152/153 (99%)	0.17	3 (1%) 62 30	56, 94, 153, 393	0
1	AZ	152/153 (99%)	0.14	1 (0%) 84 56	55, 97, 150, 271	0
1	B1	152/153 (99%)	0.29	3 (1%) 62 30	60, 95, 151, 347	0
1	B2	152/153 (99%)	0.30	3 (1%) 62 30	56, 97, 184, 226	0
1	B3	152/153 (99%)	0.32	4 (2%) 53 24	57, 96, 147, 259	0
1	B4	152/153 (99%)	0.26	2 (1%) 74 40	55, 93, 153, 393	0
1	BA	152/153 (99%)	0.22	4 (2%) 53 24	55, 97, 147, 271	0
1	BB	152/153 (99%)	0.39	3 (1%) 62 30	60, 95, 151, 347	0
1	BC	152/153 (99%)	0.29	4 (2%) 53 24	56, 97, 184, 226	0
1	BD	152/153 (99%)	0.19	1 (0%) 84 56	57, 96, 147, 259	0
1	BE	152/153 (99%)	0.37	2 (1%) 74 40	55, 93, 153, 393	0
1	BF	152/153 (99%)	0.32	3 (1%) 62 30	55, 97, 147, 271	0
1	BG	152/153 (99%)	0.25	2 (1%) 74 40	60, 95, 151, 347	0
1	BH	152/153 (99%)	0.20	3 (1%) 62 30	56, 97, 184, 226	0
1	BI	152/153 (99%)	0.27	3 (1%) 62 30	57, 96, 147, 259	0
1	BJ	152/153 (99%)	0.24	3 (1%) 62 30	55, 93, 153, 393	0
1	BK	152/153 (99%)	0.07	1 (0%) 84 56	55, 97, 147, 271	0
1	BL	152/153 (99%)	0.31	3 (1%) 62 30	60, 95, 151, 347	0
1	BM	152/153 (99%)	0.31	2 (1%) 74 40	56, 97, 184, 226	0
1	BN	152/153 (99%)	0.14	1 (0%) 84 56	57, 96, 147, 259	0
1	BO	152/153 (99%)	0.19	4 (2%) 53 24	55, 93, 153, 393	0
1	BP	152/153 (99%)	0.24	2 (1%) 74 40	55, 97, 147, 271	0
1	BQ	152/153 (99%)	0.25	2 (1%) 74 40	60, 95, 151, 347	0
1	BR	152/153 (99%)	0.25	2 (1%) 74 40	56, 97, 184, 226	0
1	BS	152/153 (99%)	0.30	3 (1%) 62 30	57, 96, 147, 259	0
1	BT	152/153 (99%)	0.29	3 (1%) 62 30	55, 93, 153, 393	0
1	BU	152/153 (99%)	0.27	1 (0%) 84 56	55, 97, 147, 271	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	BV	152/153 (99%)	0.39	4 (2%) 53 24	60, 95, 151, 347	0
1	BW	152/153 (99%)	0.26	2 (1%) 74 40	56, 97, 184, 226	0
1	BX	152/153 (99%)	0.19	0 100 100	57, 96, 147, 259	0
1	BY	152/153 (99%)	0.26	3 (1%) 62 30	55, 93, 153, 393	0
1	BZ	152/153 (99%)	0.20	2 (1%) 74 40	55, 97, 147, 271	0
1	C1	152/153 (99%)	0.96	12 (7%) 13 7	60, 95, 151, 347	0
1	C2	152/153 (99%)	0.80	9 (5%) 22 9	56, 97, 184, 226	0
1	C3	152/153 (99%)	0.98	15 (9%) 8 5	57, 96, 147, 259	0
1	C4	152/153 (99%)	1.05	17 (11%) 6 4	55, 93, 153, 393	0
1	CA	152/153 (99%)	0.48	6 (3%) 37 16	55, 97, 147, 271	0
1	CB	152/153 (99%)	0.69	10 (6%) 18 8	60, 95, 151, 347	0
1	CC	152/153 (99%)	0.75	7 (4%) 31 14	56, 97, 184, 226	0
1	CD	152/153 (99%)	0.83	12 (7%) 13 7	57, 96, 147, 259	0
1	CE	152/153 (99%)	0.41	5 (3%) 44 20	55, 93, 153, 393	0
1	CF	152/153 (99%)	0.86	12 (7%) 13 7	55, 97, 147, 271	0
1	CG	152/153 (99%)	0.50	3 (1%) 62 30	60, 95, 151, 347	0
1	CH	152/153 (99%)	0.20	2 (1%) 74 40	56, 97, 184, 226	0
1	CI	152/153 (99%)	0.23	3 (1%) 62 30	57, 96, 147, 259	0
1	CJ	152/153 (99%)	0.62	9 (5%) 22 9	55, 93, 153, 393	0
1	CK	152/153 (99%)	0.07	1 (0%) 84 56	55, 97, 147, 271	0
1	CL	152/153 (99%)	0.24	2 (1%) 74 40	60, 95, 151, 347	0
1	CM	152/153 (99%)	0.15	2 (1%) 74 40	56, 97, 184, 226	0
1	CN	152/153 (99%)	0.30	2 (1%) 74 40	57, 96, 147, 259	0
1	CO	152/153 (99%)	0.22	4 (2%) 53 24	55, 93, 153, 393	0
1	CP	152/153 (99%)	0.88	11 (7%) 15 7	55, 97, 147, 271	0
1	CQ	152/153 (99%)	1.07	16 (10%) 7 5	60, 95, 151, 347	0
1	CR	152/153 (99%)	1.18	15 (9%) 8 5	56, 97, 184, 226	0
1	CS	152/153 (99%)	1.30	30 (19%) 2 2	57, 96, 147, 259	0
1	CT	152/153 (99%)	1.18	18 (11%) 5 4	55, 93, 153, 393	0
1	CU	152/153 (99%)	0.42	5 (3%) 44 20	55, 97, 147, 271	0
1	CV	152/153 (99%)	0.72	8 (5%) 25 10	60, 95, 151, 347	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	CW	152/153 (99%)	0.74	11 (7%) 15 7	56, 97, 184, 226	0
1	CX	152/153 (99%)	0.46	3 (1%) 62 30	57, 96, 147, 259	0
1	CY	152/153 (99%)	0.24	4 (2%) 53 24	55, 93, 153, 393	0
1	CZ	152/153 (99%)	1.17	21 (13%) 4 3	55, 97, 147, 271	0
All	All	13680/13770 (99%)	0.37	405 (2%) 48 22	55, 96, 157, 393	0

The worst 5 of 405 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BE	151	HIS	13.0
1	AN	153	ALA	11.9
1	BM	153	ALA	11.8
1	CD	153	ALA	10.7
1	AM	153	ALA	10.5

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	PO4	AI	1001	5/5	0.31	2.04	58,65,106,138	0
2	PO4	AX	1001	5/5	0.31	2.03	58,65,106,138	0
2	PO4	BG	1001	5/5	0.27	1.47	74,78,104,213	0
2	PO4	CD	1001	5/5	0.32	1.04	58,65,106,138	0
2	PO4	BD	1001	5/5	0.23	0.39	58,65,106,138	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	PO4	AW	201	5/5	0.24	0.18	34,46,111,116	0
2	PO4	CQ	1201	5/5	0.32	0.13	74,78,104,213	0
2	PO4	AK	1001	5/5	0.23	0.09	80,81,83,167	0
2	PO4	CB	1001	5/5	0.35	0.08	74,78,104,213	0
2	PO4	AL	1001	5/5	0.24	0.06	74,78,104,213	0
2	PO4	AJ	1001	5/5	0.23	-0.03	23,28,53,142	0
2	PO4	AG	1001	5/5	0.23	-0.13	74,78,104,213	0
2	PO4	AY	1001	5/5	0.24	-0.15	23,28,53,142	0
2	PO4	CZ	201	5/5	0.29	-0.20	80,81,83,167	0
2	PO4	CY	4201	5/5	0.32	-0.21	23,28,53,142	0
2	PO4	CJ	4201	5/5	0.25	-0.22	23,28,53,142	0
2	PO4	BR	1001	5/5	0.22	-0.24	34,46,111,116	0
2	PO4	AH	1001	5/5	0.20	-0.26	34,46,111,116	0
2	PO4	BL	1001	5/5	0.26	-0.27	74,78,104,213	0
2	PO4	CT	4201	5/5	0.38	-0.37	23,28,53,142	0
2	PO4	BZ	1001	5/5	0.20	-0.39	80,81,83,167	0
2	PO4	CK	201	5/5	0.18	-0.40	80,81,83,167	0
2	PO4	CN	3201	5/5	0.20	-0.43	58,65,106,138	0
2	PO4	AC	1001	5/5	0.17	-0.48	34,46,111,116	0
2	PO4	CG	1201	5/5	0.24	-0.48	74,78,104,213	0
2	PO4	BX	1001	5/5	0.17	-0.53	58,65,106,138	0
2	PO4	BB	1001	5/5	0.21	-0.53	74,78,104,213	0
2	PO4	B1	1001	5/5	0.20	-0.57	74,78,104,213	0
2	PO4	BP	1001	5/5	0.17	-0.70	80,81,83,167	0
2	PO4	BE	1001	5/5	0.22	-0.71	23,28,53,142	0
2	PO4	C2	2201	5/5	0.28	-0.71	34,46,111,116	0
2	PO4	AM	1001	5/5	0.19	-0.73	34,46,111,116	0
2	PO4	CW	2201	5/5	0.29	-0.74	34,46,111,116	0
2	PO4	CH	2201	5/5	0.24	-0.75	34,46,111,116	0
2	PO4	AV	1001	5/5	0.19	-0.75	74,78,104,213	0
2	PO4	BK	1001	5/5	0.17	-0.76	80,81,83,167	0
2	PO4	CX	3201	5/5	0.20	-0.78	58,65,106,138	0
2	PO4	CL	1201	5/5	0.19	-0.81	74,78,104,213	0
2	PO4	AS	1001	5/5	0.18	-0.83	58,65,106,138	0
2	PO4	AB	1001	5/5	0.19	-0.83	74,78,104,213	0
2	PO4	BA	1001	5/5	0.18	-0.84	80,81,83,167	0
2	PO4	AZ	1001	5/5	0.18	-0.88	80,81,83,167	0
2	PO4	BW	1001	5/5	0.20	-0.89	34,46,111,116	0
2	PO4	CC	1001	5/5	0.30	-0.90	34,46,111,116	0
2	PO4	BQ	1001	5/5	0.17	-0.90	74,78,104,213	0
2	PO4	AA	1001	5/5	0.15	-0.98	80,81,83,167	0
2	PO4	CF	201	5/5	0.22	-0.98	80,81,83,167	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	PO4	BV	1001	5/5	0.22	-1.00	74,78,104,213	0
2	PO4	A3	1001	5/5	0.17	-1.04	58,65,106,138	0
2	PO4	BH	1001	5/5	0.16	-1.07	34,46,111,116	0
2	PO4	BM	1001	5/5	0.16	-1.07	34,46,111,116	0
2	PO4	AP	1001	5/5	0.17	-1.08	80,81,83,167	0
2	PO4	AT	1001	5/5	0.19	-1.12	23,28,53,142	0
2	PO4	A4	1001	5/5	0.16	-1.12	23,28,53,142	0
2	PO4	CE	1001	5/5	0.27	-1.14	23,28,53,142	0
2	PO4	BF	1001	5/5	0.21	-1.18	80,81,83,167	0
2	PO4	B3	1001	5/5	0.17	-1.19	58,65,106,138	0
2	PO4	CA	1001	5/5	0.18	-1.21	80,81,83,167	0
2	PO4	BS	1001	5/5	0.20	-1.30	58,65,106,138	0
2	PO4	A2	1001	5/5	0.16	-1.32	34,46,111,116	0
2	PO4	B4	1001	5/5	0.17	-1.34	23,28,53,142	0
2	PO4	CV	1201	5/5	0.20	-1.37	74,78,104,213	0
2	PO4	B2	1001	5/5	0.20	-1.38	34,46,111,116	0
2	PO4	CU	201	5/5	0.13	-1.41	80,81,83,167	0
2	PO4	A1	1001	5/5	0.17	-1.47	74,78,104,213	0
2	PO4	AU	1001	5/5	0.18	-1.52	80,81,83,167	0
2	PO4	BC	1001	5/5	0.15	-1.53	34,46,111,116	0
2	PO4	CS	3201	5/5	0.30	-1.55	58,65,106,138	0
2	PO4	AN	1001	5/5	0.16	-1.59	58,65,106,138	0
2	PO4	BO	1001	5/5	0.12	-1.60	23,28,53,142	0
2	PO4	CP	201	5/5	0.23	-1.61	80,81,83,167	0
2	PO4	CM	2201	5/5	0.15	-1.63	34,46,111,116	0
2	PO4	BT	1001	5/5	0.12	-1.64	23,28,53,142	0
2	PO4	AD	1001	5/5	0.15	-1.68	58,65,106,138	0
2	PO4	BJ	1001	5/5	0.13	-1.71	23,28,53,142	0
2	PO4	C3	3201	5/5	0.22	-1.76	58,65,106,138	0
2	PO4	BU	1001	5/5	0.12	-1.79	80,81,83,167	0
2	PO4	AO	1001	5/5	0.17	-1.82	23,28,53,142	0
2	PO4	CR	2201	5/5	0.26	-1.90	34,46,111,116	0
2	PO4	AQ	1001	5/5	0.16	-1.99	74,78,104,213	0
2	PO4	AF	1001	5/5	0.14	-2.01	80,81,83,167	0
2	PO4	BY	1001	5/5	0.16	-2.05	23,28,53,142	0
2	PO4	BN	1001	5/5	0.14	-2.15	58,65,106,138	0
2	PO4	C1	1201	5/5	0.23	-2.18	74,78,104,213	0
2	PO4	BI	1001	5/5	0.14	-2.25	58,65,106,138	0
2	PO4	CI	3201	5/5	0.16	-2.30	58,65,106,138	0
2	PO4	AR	1001	5/5	0.13	-2.41	34,46,111,116	0
2	PO4	CO	4201	5/5	0.18	-2.67	23,28,53,142	0
2	PO4	AE	1001	5/5	0.11	-3.87	23,28,53,142	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	PO4	C4	4201	5/5	0.25	-4.53	23,28,53,142	0

6.5 Other polymers ⓘ

There are no such residues in this entry.