



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:08 AM UTC

PDB ID : 4KQ6 / pdb_00004kq6
Title : Product complex of lumazine synthase from candida glabrata
Authors : Shankar, M.; Wilbanks, S.M.; Nakatani, Y.; Monk, B.C.; Tyndall, J.D.A.
Deposited on : 2013-05-14
Resolution : 2.24 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

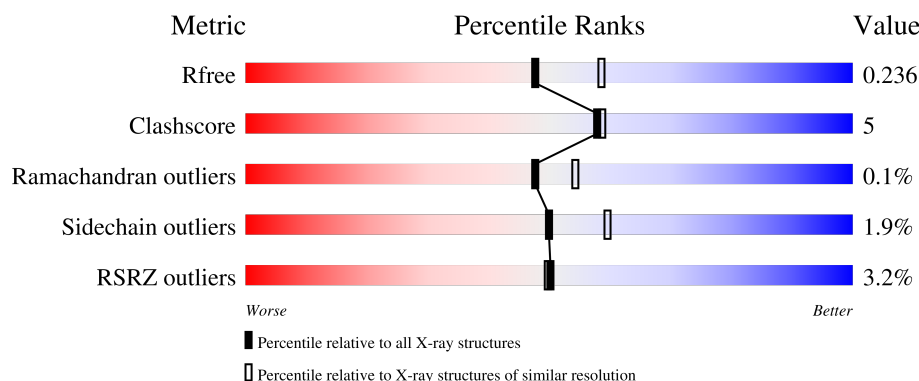
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.24 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	179	 3% 82% 7% 12%
1	B	179	 2% 76% 12% 12%
1	C	179	 3% 75% 12% 12%
1	D	179	 4% 79% 12% 8%
1	E	179	 2% 77% 11% 12%

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Mol	Chain	Length	Quality of chain
1	F	179	2%79%8%12%
1	G	179	2%77%12%11%
1	H	179	4%70%18%12%
1	I	179	3%75%12%12%
1	J	179	2%74%13%12%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13313 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 6,7-dimethyl-8-ribityllumazine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	158	Total	C	N	O	S	0	0	0
			1212	769	209	228	6			
1	B	158	Total	C	N	O	S	0	0	0
			1212	769	209	228	6			
1	C	158	Total	C	N	O	S	0	0	0
			1213	769	209	229	6			
1	D	165	Total	C	N	O	S	0	0	0
			1264	802	218	238	6			
1	E	157	Total	C	N	O	S	0	0	0
			1207	766	208	227	6			
1	F	157	Total	C	N	O	S	0	0	0
			1207	766	208	227	6			
1	G	160	Total	C	N	O	S	0	0	0
			1232	781	212	233	6			
1	H	158	Total	C	N	O	S	0	1	0
			1217	773	209	228	7			
1	I	158	Total	C	N	O	S	0	0	0
			1213	769	209	229	6			
1	J	157	Total	C	N	O	S	0	0	0
			1207	766	208	227	6			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	initiating methionine	UNP Q6FXA8
A	-9	ARG	-	expression tag	UNP Q6FXA8
A	-8	GLY	-	expression tag	UNP Q6FXA8
A	-7	SER	-	expression tag	UNP Q6FXA8
A	-6	HIS	-	expression tag	UNP Q6FXA8
A	-5	HIS	-	expression tag	UNP Q6FXA8
A	-4	HIS	-	expression tag	UNP Q6FXA8
A	-3	HIS	-	expression tag	UNP Q6FXA8
A	-2	HIS	-	expression tag	UNP Q6FXA8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	HIS	-	expression tag	UNP Q6FXA8
A	0	GLY	-	expression tag	UNP Q6FXA8
A	1	SER	-	expression tag	UNP Q6FXA8
B	-10	MET	-	initiating methionine	UNP Q6FXA8
B	-9	ARG	-	expression tag	UNP Q6FXA8
B	-8	GLY	-	expression tag	UNP Q6FXA8
B	-7	SER	-	expression tag	UNP Q6FXA8
B	-6	HIS	-	expression tag	UNP Q6FXA8
B	-5	HIS	-	expression tag	UNP Q6FXA8
B	-4	HIS	-	expression tag	UNP Q6FXA8
B	-3	HIS	-	expression tag	UNP Q6FXA8
B	-2	HIS	-	expression tag	UNP Q6FXA8
B	-1	HIS	-	expression tag	UNP Q6FXA8
B	0	GLY	-	expression tag	UNP Q6FXA8
B	1	SER	-	expression tag	UNP Q6FXA8
C	-10	MET	-	initiating methionine	UNP Q6FXA8
C	-9	ARG	-	expression tag	UNP Q6FXA8
C	-8	GLY	-	expression tag	UNP Q6FXA8
C	-7	SER	-	expression tag	UNP Q6FXA8
C	-6	HIS	-	expression tag	UNP Q6FXA8
C	-5	HIS	-	expression tag	UNP Q6FXA8
C	-4	HIS	-	expression tag	UNP Q6FXA8
C	-3	HIS	-	expression tag	UNP Q6FXA8
C	-2	HIS	-	expression tag	UNP Q6FXA8
C	-1	HIS	-	expression tag	UNP Q6FXA8
C	0	GLY	-	expression tag	UNP Q6FXA8
C	1	SER	-	expression tag	UNP Q6FXA8
D	-10	MET	-	initiating methionine	UNP Q6FXA8
D	-9	ARG	-	expression tag	UNP Q6FXA8
D	-8	GLY	-	expression tag	UNP Q6FXA8
D	-7	SER	-	expression tag	UNP Q6FXA8
D	-6	HIS	-	expression tag	UNP Q6FXA8
D	-5	HIS	-	expression tag	UNP Q6FXA8
D	-4	HIS	-	expression tag	UNP Q6FXA8
D	-3	HIS	-	expression tag	UNP Q6FXA8
D	-2	HIS	-	expression tag	UNP Q6FXA8
D	-1	HIS	-	expression tag	UNP Q6FXA8
D	0	GLY	-	expression tag	UNP Q6FXA8
D	1	SER	-	expression tag	UNP Q6FXA8
E	-10	MET	-	initiating methionine	UNP Q6FXA8
E	-9	ARG	-	expression tag	UNP Q6FXA8
E	-8	GLY	-	expression tag	UNP Q6FXA8

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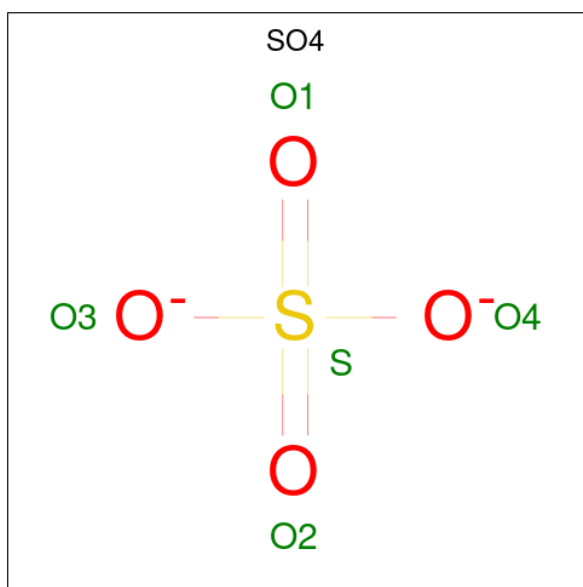
Chain	Residue	Modelled	Actual	Comment	Reference
E	-7	SER	-	expression tag	UNP Q6FXA8
E	-6	HIS	-	expression tag	UNP Q6FXA8
E	-5	HIS	-	expression tag	UNP Q6FXA8
E	-4	HIS	-	expression tag	UNP Q6FXA8
E	-3	HIS	-	expression tag	UNP Q6FXA8
E	-2	HIS	-	expression tag	UNP Q6FXA8
E	-1	HIS	-	expression tag	UNP Q6FXA8
E	0	GLY	-	expression tag	UNP Q6FXA8
E	1	SER	-	expression tag	UNP Q6FXA8
F	-10	MET	-	initiating methionine	UNP Q6FXA8
F	-9	ARG	-	expression tag	UNP Q6FXA8
F	-8	GLY	-	expression tag	UNP Q6FXA8
F	-7	SER	-	expression tag	UNP Q6FXA8
F	-6	HIS	-	expression tag	UNP Q6FXA8
F	-5	HIS	-	expression tag	UNP Q6FXA8
F	-4	HIS	-	expression tag	UNP Q6FXA8
F	-3	HIS	-	expression tag	UNP Q6FXA8
F	-2	HIS	-	expression tag	UNP Q6FXA8
F	-1	HIS	-	expression tag	UNP Q6FXA8
F	0	GLY	-	expression tag	UNP Q6FXA8
F	1	SER	-	expression tag	UNP Q6FXA8
G	-10	MET	-	initiating methionine	UNP Q6FXA8
G	-9	ARG	-	expression tag	UNP Q6FXA8
G	-8	GLY	-	expression tag	UNP Q6FXA8
G	-7	SER	-	expression tag	UNP Q6FXA8
G	-6	HIS	-	expression tag	UNP Q6FXA8
G	-5	HIS	-	expression tag	UNP Q6FXA8
G	-4	HIS	-	expression tag	UNP Q6FXA8
G	-3	HIS	-	expression tag	UNP Q6FXA8
G	-2	HIS	-	expression tag	UNP Q6FXA8
G	-1	HIS	-	expression tag	UNP Q6FXA8
G	0	GLY	-	expression tag	UNP Q6FXA8
G	1	SER	-	expression tag	UNP Q6FXA8
H	-10	MET	-	initiating methionine	UNP Q6FXA8
H	-9	ARG	-	expression tag	UNP Q6FXA8
H	-8	GLY	-	expression tag	UNP Q6FXA8
H	-7	SER	-	expression tag	UNP Q6FXA8
H	-6	HIS	-	expression tag	UNP Q6FXA8
H	-5	HIS	-	expression tag	UNP Q6FXA8
H	-4	HIS	-	expression tag	UNP Q6FXA8
H	-3	HIS	-	expression tag	UNP Q6FXA8
H	-2	HIS	-	expression tag	UNP Q6FXA8

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Chain	Residue	Modelled	Actual	Comment	Reference
H	-1	HIS	-	expression tag	UNP Q6FXA8
H	0	GLY	-	expression tag	UNP Q6FXA8
H	1	SER	-	expression tag	UNP Q6FXA8
I	-10	MET	-	initiating methionine	UNP Q6FXA8
I	-9	ARG	-	expression tag	UNP Q6FXA8
I	-8	GLY	-	expression tag	UNP Q6FXA8
I	-7	SER	-	expression tag	UNP Q6FXA8
I	-6	HIS	-	expression tag	UNP Q6FXA8
I	-5	HIS	-	expression tag	UNP Q6FXA8
I	-4	HIS	-	expression tag	UNP Q6FXA8
I	-3	HIS	-	expression tag	UNP Q6FXA8
I	-2	HIS	-	expression tag	UNP Q6FXA8
I	-1	HIS	-	expression tag	UNP Q6FXA8
I	0	GLY	-	expression tag	UNP Q6FXA8
I	1	SER	-	expression tag	UNP Q6FXA8
J	-10	MET	-	initiating methionine	UNP Q6FXA8
J	-9	ARG	-	expression tag	UNP Q6FXA8
J	-8	GLY	-	expression tag	UNP Q6FXA8
J	-7	SER	-	expression tag	UNP Q6FXA8
J	-6	HIS	-	expression tag	UNP Q6FXA8
J	-5	HIS	-	expression tag	UNP Q6FXA8
J	-4	HIS	-	expression tag	UNP Q6FXA8
J	-3	HIS	-	expression tag	UNP Q6FXA8
J	-2	HIS	-	expression tag	UNP Q6FXA8
J	-1	HIS	-	expression tag	UNP Q6FXA8
J	0	GLY	-	expression tag	UNP Q6FXA8
J	1	SER	-	expression tag	UNP Q6FXA8

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



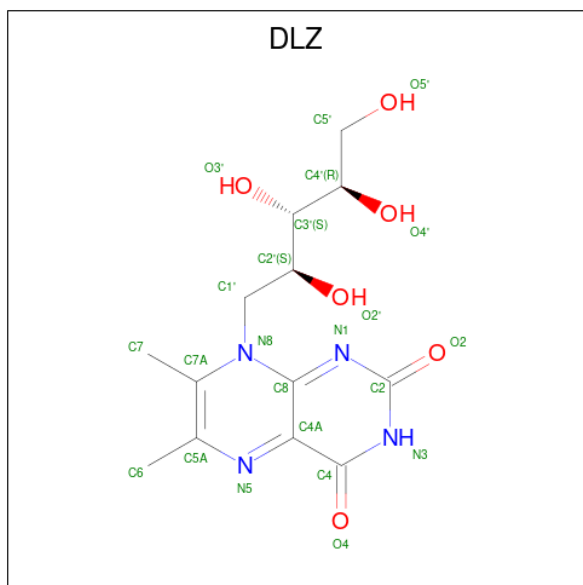
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	I	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		
2	J	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 1-deoxy-1-(6,7-dimethyl-2,4-dioxo-3,4-dihydropteridin-8(2H)-yl)-D-ribose (CCD ID: DLZ) (formula: C₁₃H₁₈N₄O₆).



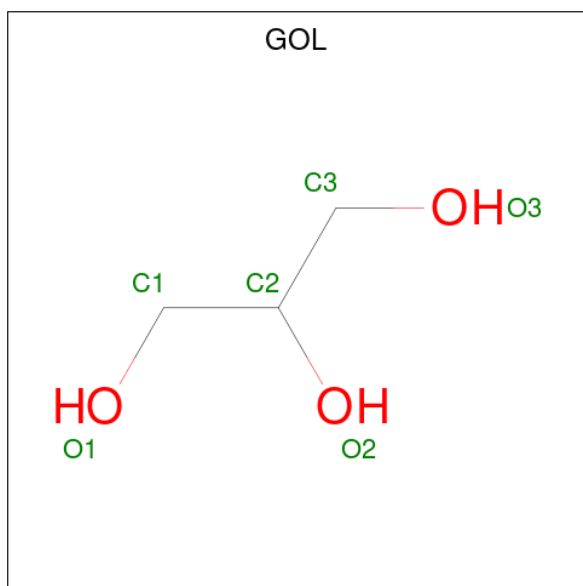
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			23	13	4	6		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	I	1	Total	C	N	O	0	0
			23	13	4	6		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		
4	I	1	Total	C	O	0	0
			6	3	3		
4	J	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	101	Total	O	0	0
			101	101		
5	B	115	Total	O	0	0
			115	115		
5	C	105	Total	O	0	0
			105	105		

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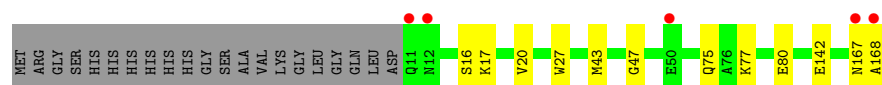
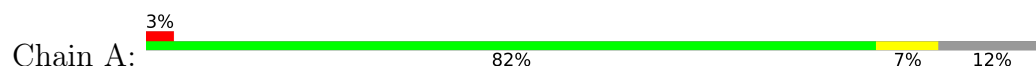
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	103	Total 103	O 103	0	0
5	E	107	Total 107	O 107	0	0
5	F	103	Total 103	O 103	0	0
5	G	70	Total 70	O 70	0	0
5	H	66	Total 66	O 66	0	0
5	I	75	Total 75	O 75	0	0
5	J	88	Total 88	O 88	0	0

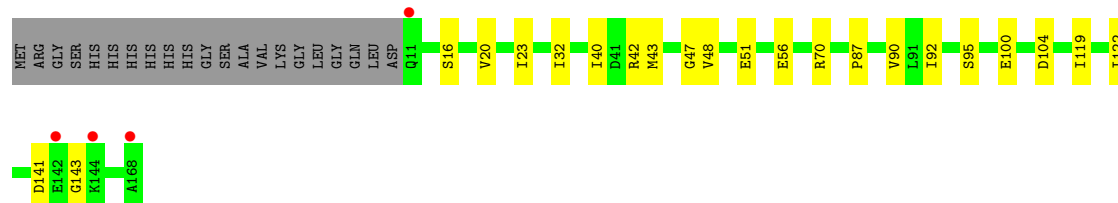
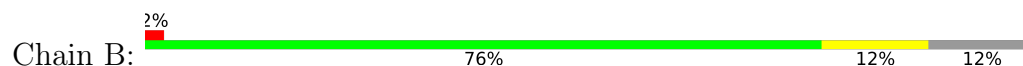
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

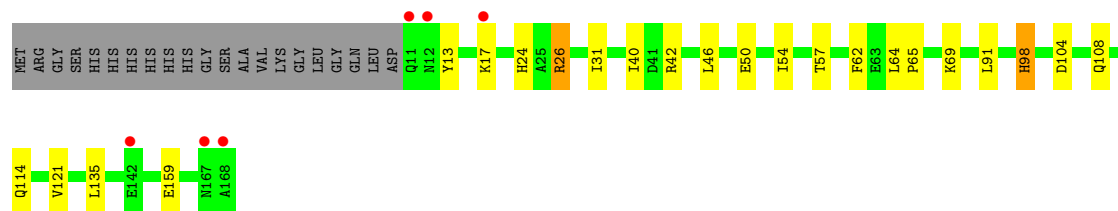
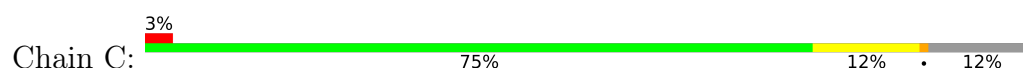
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



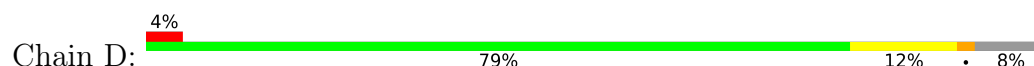
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



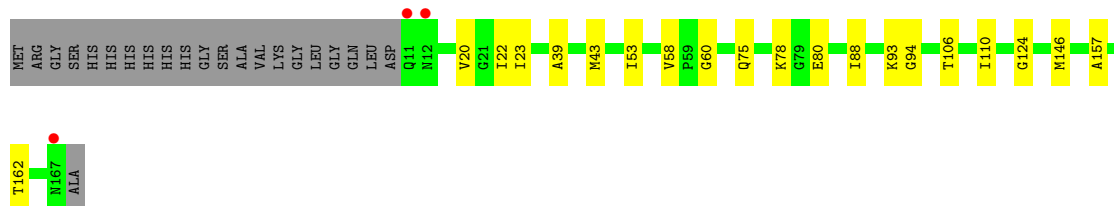
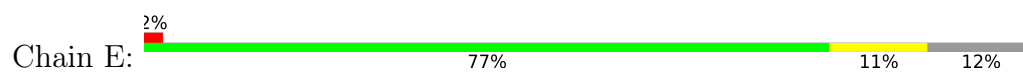
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



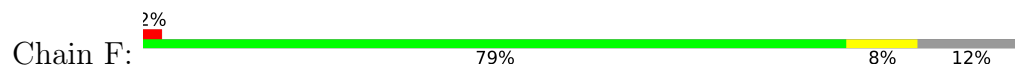
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



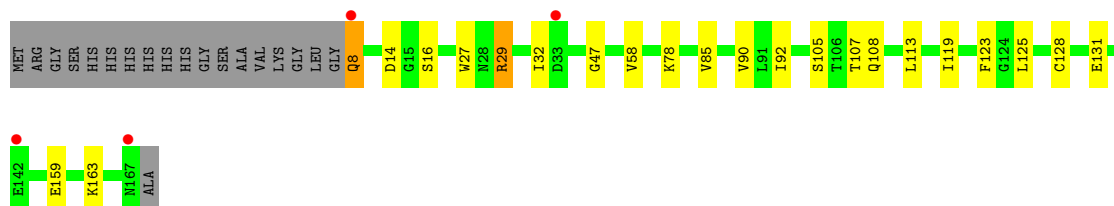
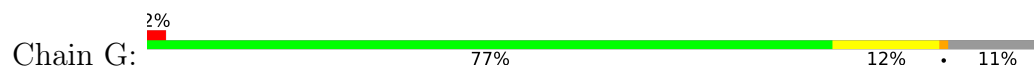
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



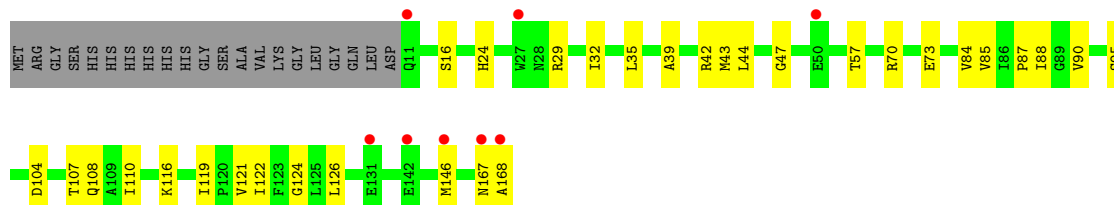
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



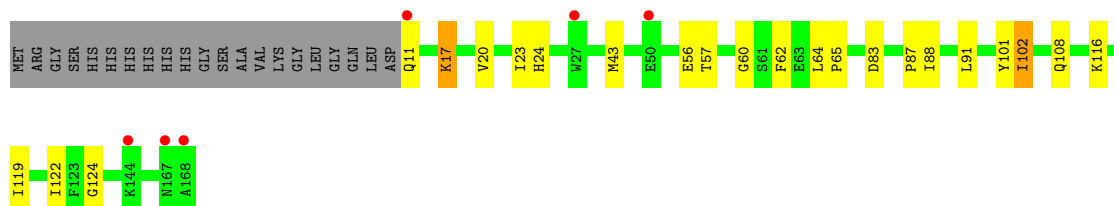
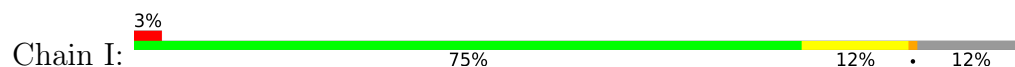
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



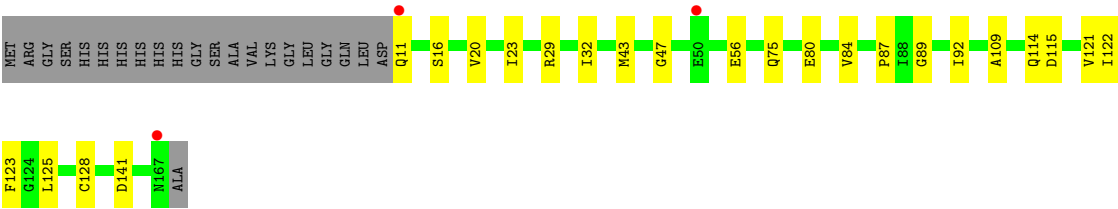
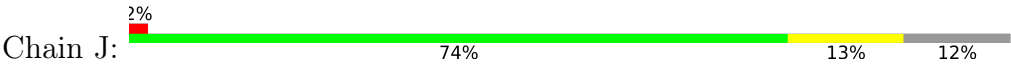
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	84.75Å 84.84Å 310.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.43 – 2.24 47.43 – 2.24	Depositor EDS
% Data completeness (in resolution range)	95.2 (47.43-2.24) 95.2 (47.43-2.24)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	11.63 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.200 , 0.252 (Not available) , 0.236	Depositor DCC
R_{free} test set	2012 reflections (1.95%)	wwPDB-VP
Wilson B-factor (Å ²)	15.1	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.054 for k,h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13313	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4, DLZ

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.28	0/1232	1.25	2/1663 (0.1%)
1	B	1.26	3/1232 (0.2%)	1.30	6/1663 (0.4%)
1	C	1.26	3/1233 (0.2%)	1.33	5/1663 (0.3%)
1	D	1.36	2/1284 (0.2%)	1.29	7/1732 (0.4%)
1	E	1.39	4/1227 (0.3%)	1.25	5/1656 (0.3%)
1	F	1.36	3/1227 (0.2%)	1.29	6/1656 (0.4%)
1	G	1.23	2/1252 (0.2%)	1.25	4/1690 (0.2%)
1	H	1.21	2/1240 (0.2%)	1.19	2/1673 (0.1%)
1	I	1.25	0/1233	1.22	4/1663 (0.2%)
1	J	1.24	0/1227	1.24	1/1656 (0.1%)
All	All	1.28	19/12387 (0.2%)	1.26	42/16715 (0.3%)

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	D	0	1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	54	ILE	C-O	6.86	1.31	1.24
1	G	119	ILE	C-O	-6.54	1.19	1.24
1	E	60	GLY	N-CA	6.36	1.51	1.44
1	E	53	ILE	C-O	-6.20	1.17	1.24
1	G	113	LEU	CA-C	5.82	1.60	1.52
1	H	104	ASP	N-CA	-5.77	1.39	1.46
1	E	23	ILE	N-CA	-5.64	1.39	1.46
1	C	98	HIS	C-O	-5.46	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	141	ASP	CA-C	5.42	1.59	1.52
1	F	133	GLN	N-CA	5.35	1.52	1.46
1	D	31	ILE	C-O	5.34	1.30	1.24
1	F	105	SER	CA-C	5.34	1.59	1.52
1	C	98	HIS	N-CA	-5.24	1.40	1.46
1	F	72	ALA	N-CA	-5.15	1.40	1.46
1	E	146	MET	N-CA	5.15	1.52	1.45
1	B	92	ILE	C-O	-5.13	1.18	1.24
1	H	110	ILE	CA-C	5.09	1.59	1.52
1	D	86	ILE	CA-CB	5.02	1.58	1.54
1	B	70	ARG	CA-C	5.02	1.59	1.52

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	119	ILE	CA-C-N	-9.68	110.45	120.03
1	B	119	ILE	C-N-CA	-9.68	110.45	120.03
1	C	64	LEU	CA-C-N	-8.90	109.15	118.85
1	C	64	LEU	C-N-CA	-8.90	109.15	118.85
1	D	6	LEU	CB-CA-C	-8.73	99.31	113.37
1	D	6	LEU	N-CA-C	7.74	122.96	109.80
1	F	86	ILE	CA-C-N	-7.38	112.01	119.99
1	F	86	ILE	C-N-CA	-7.38	112.01	119.99
1	F	58	VAL	CA-C-N	-5.91	112.45	119.84
1	F	58	VAL	C-N-CA	-5.91	112.45	119.84
1	D	39	ALA	N-CA-C	-5.87	104.81	111.14
1	G	58	VAL	CA-C-N	-5.80	112.58	119.84
1	G	58	VAL	C-N-CA	-5.80	112.58	119.84
1	E	58	VAL	CA-C-N	-5.78	112.62	119.84
1	E	58	VAL	C-N-CA	-5.78	112.62	119.84
1	G	14	ASP	CB-CA-C	-5.74	100.94	109.90
1	I	119	ILE	CA-C-N	-5.74	114.05	119.85
1	I	119	ILE	C-N-CA	-5.74	114.05	119.85
1	I	101	TYR	N-CA-C	5.70	117.57	111.36
1	D	44	LEU	N-CA-C	-5.68	105.09	111.28
1	B	48	VAL	N-CA-C	-5.66	102.28	109.58
1	J	109	ALA	N-CA-C	-5.65	105.11	112.23
1	C	69	LYS	N-CA-C	-5.64	104.82	110.97
1	H	29	ARG	N-CA-C	5.54	117.33	111.28
1	E	22	ILE	CA-C-N	-5.48	115.37	122.99
1	E	22	ILE	C-N-CA	-5.48	115.37	122.99
1	C	26	ARG	CA-CB-CG	-5.47	103.15	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	51	GLU	N-CA-C	-5.39	106.47	113.16
1	B	143	GLY	N-CA-C	5.33	120.53	114.67
1	E	157	ALA	N-CA-C	-5.31	105.58	111.36
1	G	29	ARG	N-CA-C	5.28	117.44	111.11
1	D	20	VAL	CA-C-N	-5.28	117.68	122.80
1	D	20	VAL	C-N-CA	-5.28	117.68	122.80
1	C	26	ARG	NE-CZ-NH2	-5.26	114.47	119.20
1	D	36	VAL	N-CA-C	5.23	115.85	110.36
1	B	40	ILE	N-CA-C	5.20	115.82	110.36
1	H	119	ILE	N-CA-C	-5.14	104.97	109.19
1	A	77	LYS	CA-C-N	-5.13	114.14	122.65
1	A	77	LYS	C-N-CA	-5.13	114.14	122.65
1	F	37	LYS	CA-C-N	-5.08	113.77	120.22
1	F	37	LYS	C-N-CA	-5.08	113.77	120.22
1	I	83	ASP	N-CA-C	-5.04	107.30	113.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	5	GLY	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1212	0	1229	5	0
1	B	1212	0	1229	7	0
1	C	1213	0	1229	20	0
1	D	1264	0	1286	13	0
1	E	1207	0	1224	7	0
1	F	1207	0	1224	3	0
1	G	1232	0	1247	10	0
1	H	1217	0	1238	16	0
1	I	1213	0	1229	15	0
1	J	1207	0	1224	16	0
2	A	10	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	10	0	0	0	0
2	C	15	0	0	0	0
2	D	10	0	0	0	0
2	E	10	0	0	0	0
2	F	15	0	0	0	0
2	G	15	0	0	1	0
2	H	10	0	0	0	0
2	I	10	0	0	0	0
2	J	15	0	0	0	0
3	C	23	0	18	4	0
3	I	23	0	18	3	0
4	D	6	0	8	2	0
4	E	6	0	8	2	0
4	G	6	0	8	0	0
4	I	6	0	8	1	0
4	J	6	0	8	3	0
5	A	101	0	0	0	0
5	B	115	0	0	1	0
5	C	105	0	0	4	0
5	D	103	0	0	3	0
5	E	107	0	0	3	0
5	F	103	0	0	0	0
5	G	70	0	0	2	0
5	H	66	0	0	3	0
5	I	75	0	0	2	0
5	J	88	0	0	3	0
All	All	13313	0	12435	111	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 5.

All (111) 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:I:87:PRO:HD2	1:I:122:ILE:O	1.84	0.78
1:I:91:LEU:HD23	3:I:204:DLZ:N3	2.02	0.75
4:E:203:GOL:H32	5:E:316:HOH:O	1.89	0.73
1:C:104:ASP:HB3	5:C:343:HOH:O	1.90	0.70
5:H:352:HOH:O	4:I:203:GOL:H32	1.90	0.70
1:C:24:HIS:CE1	1:C:57:THR:HG22	2.27	0.70
1:I:91:LEU:HD23	3:I:204:DLZ:C4	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:HIS:CD2	3:C:204:DLZ:C6	2.78	0.67
1:F:87:PRO:HD2	1:F:122:ILE:O	1.94	0.67
1:H:87:PRO:HD2	1:H:122:ILE:O	1.94	0.67
4:D:203:GOL:H32	5:D:372:HOH:O	1.95	0.66
1:C:98:HIS:CD2	3:C:204:DLZ:H6	2.32	0.65
1:C:17:LYS:HB2	5:C:377:HOH:O	1.97	0.63
1:I:60:GLY:HA3	3:I:204:DLZ:H1'	1.81	0.62
1:G:8:GLN:N	5:G:353:HOH:O	2.32	0.61
1:E:93:LYS:HE3	1:E:94:GLY:O	2.03	0.59
1:A:167:ASN:O	1:A:168:ALA:C	2.45	0.58
1:F:88:ILE:HA	1:F:124:GLY:O	2.04	0.57
1:C:24:HIS:NE2	1:C:57:THR:HG22	2.20	0.56
4:J:204:GOL:C3	5:J:309:HOH:O	2.54	0.56
1:C:42:ARG:HD2	5:C:341:HOH:O	2.06	0.55
4:J:204:GOL:H31	5:J:309:HOH:O	2.07	0.55
1:J:11:GLN:N	1:J:11:GLN:CD	2.65	0.55
1:H:42:ARG:CD	5:H:313:HOH:O	2.54	0.55
1:H:42:ARG:HD2	5:H:313:HOH:O	2.06	0.55
1:C:24:HIS:CE1	1:C:57:THR:CG2	2.90	0.54
1:C:98:HIS:NE2	3:C:204:DLZ:H7A	2.23	0.54
1:C:50:GLU:HG2	1:D:4:LYS:HD2	1.90	0.54
1:A:16:SER:HA	1:A:47:GLY:O	2.08	0.53
1:H:88:ILE:HA	1:H:124:GLY:O	2.09	0.53
1:D:114:GLN:HG2	1:D:121:VAL:HG23	1.90	0.53
1:E:88:ILE:HA	1:E:124:GLY:O	2.09	0.53
1:D:20:VAL:HG11	1:D:43:MET:SD	2.49	0.53
1:D:64:LEU:N	1:D:65:PRO:CD	2.72	0.52
1:B:42:ARG:HD2	5:B:415:HOH:O	2.09	0.52
1:G:32:ILE:HD13	1:G:90:VAL:HG23	1.92	0.52
1:B:87:PRO:HD2	1:B:122:ILE:O	2.10	0.51
4:D:203:GOL:C3	5:D:372:HOH:O	2.55	0.51
1:F:20:VAL:HG11	1:F:43:MET:SD	2.50	0.51
1:B:20:VAL:HG11	1:B:43:MET:SD	2.51	0.50
1:J:16:SER:HA	1:J:47:GLY:O	2.12	0.50
1:A:75:GLN:HB3	1:A:80:GLU:O	2.12	0.50
1:C:13:TYR:CZ	1:C:159:GLU:HG3	2.47	0.50
1:C:91:LEU:HD23	3:C:204:DLZ:C4	2.42	0.50
1:G:159:GLU:OE2	1:G:163:LYS:NZ	2.44	0.49
1:I:17:LYS:HE3	5:I:333:HOH:O	2.12	0.49
1:H:167:ASN:O	1:H:168:ALA:HB2	2.12	0.49
1:J:114:GLN:HG2	1:J:121:VAL:HG23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:16:SER:HA	1:G:47:GLY:O	2.13	0.48
1:J:20:VAL:HG11	1:J:43:MET:SD	2.54	0.48
1:C:114:GLN:HG2	1:C:121:VAL:HG23	1.95	0.47
1:G:78:LYS:NZ	5:G:332:HOH:O	2.46	0.47
1:J:20:VAL:HG13	1:J:84:VAL:HG13	1.97	0.47
1:A:20:VAL:HG11	1:A:43:MET:SD	2.54	0.47
1:E:75:GLN:HB3	1:E:80:GLU:O	2.15	0.47
1:I:20:VAL:HG11	1:I:43:MET:SD	2.55	0.47
1:C:31:ILE:HG12	1:C:135:LEU:HD23	1.96	0.47
1:A:27:TRP:HE3	2:A:201:SO4:O4	1.97	0.47
1:J:122:ILE:HG23	1:J:122:ILE:HD12	1.73	0.46
1:H:85:VAL:CG1	1:H:121:VAL:HG22	2.46	0.46
1:D:6:LEU:HB2	1:D:7:GLY:H	1.55	0.46
1:I:88:ILE:HA	1:I:124:GLY:O	2.16	0.46
1:C:46:LEU:HD23	1:C:46:LEU:HA	1.85	0.45
1:I:102:ILE:HD11	1:J:123:PHE:CE2	2.51	0.45
1:C:40:ILE:HD13	1:D:4:LYS:HE3	1.97	0.45
1:D:75:GLN:HB3	1:D:80:GLU:O	2.16	0.45
1:B:32:ILE:HD13	1:B:90:VAL:HG23	1.99	0.45
1:E:78:LYS:HE3	5:E:337:HOH:O	2.16	0.45
1:G:105:SER:HB3	1:H:107:THR:HG22	1.99	0.45
1:I:116:LYS:CE	1:J:115:ASP:OD2	2.65	0.45
1:I:116:LYS:HE2	1:J:115:ASP:OD2	2.17	0.45
1:D:88:ILE:HA	1:D:124:GLY:O	2.17	0.44
1:J:23:ILE:HA	1:J:56:GLU:O	2.18	0.44
1:D:42:ARG:NH1	5:D:386:HOH:O	2.50	0.44
1:J:75:GLN:HB3	1:J:80:GLU:O	2.17	0.44
1:H:16:SER:HA	1:H:47:GLY:O	2.17	0.44
1:C:26:ARG:NH2	1:D:159:GLU:OE1	2.44	0.43
1:G:92:ILE:HA	1:G:128:CYS:O	2.18	0.43
1:I:62:PHE:O	1:I:65:PRO:HD2	2.18	0.43
1:J:29:ARG:HA	1:J:29:ARG:HD2	1.87	0.43
1:B:16:SER:HA	1:B:47:GLY:O	2.18	0.43
1:I:24:HIS:O	1:I:57:THR:HA	2.19	0.43
1:H:70:ARG:HD2	1:H:73:GLU:OE1	2.19	0.43
1:E:39:ALA:O	1:E:43:MET:HG3	2.19	0.43
1:G:27:TRP:HE3	2:G:201:SO4:O1	2.01	0.43
1:H:44:LEU:HD23	1:H:44:LEU:HA	1.70	0.43
1:H:39:ALA:O	1:H:43:MET:HG3	2.19	0.42
1:D:64:LEU:N	1:D:65:PRO:HD3	2.35	0.42
1:G:107:THR:HG23	1:G:123:PHE:CE1	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:GLN:NE2	5:C:375:HOH:O	2.48	0.42
1:H:35:LEU:HD11	1:H:126:LEU:CD1	2.49	0.42
4:J:204:GOL:H32	5:J:309:HOH:O	2.17	0.42
1:H:24:HIS:CE1	1:H:57:THR:HG22	2.55	0.42
1:I:108:GLN:NE2	5:I:338:HOH:O	2.53	0.42
1:J:32:ILE:HD11	1:J:89:GLY:HA2	2.01	0.42
1:B:100:GLU:O	1:B:104:ASP:HB2	2.21	0.41
1:I:64:LEU:HB2	1:I:65:PRO:HD3	2.03	0.41
1:D:141:ASP:OD1	1:D:141:ASP:N	2.53	0.41
1:E:20:VAL:HG11	1:E:43:MET:SD	2.61	0.41
1:E:106:THR:O	1:E:110:ILE:HG13	2.21	0.41
1:G:108:GLN:OE1	1:H:108:GLN:NE2	2.54	0.41
1:I:23:ILE:HA	1:I:56:GLU:O	2.21	0.41
1:J:87:PRO:HD2	1:J:122:ILE:O	2.21	0.41
1:C:62:PHE:O	1:C:65:PRO:HD2	2.21	0.41
1:H:167:ASN:O	1:H:168:ALA:CB	2.70	0.40
1:J:141:ASP:N	1:J:141:ASP:OD1	2.54	0.40
1:C:50:GLU:OE1	1:D:3:VAL:HA	2.21	0.40
1:H:32:ILE:HD13	1:H:90:VAL:HG23	2.03	0.40
1:J:92:ILE:HA	1:J:128:CYS:O	2.21	0.40
4:E:203:GOL:C3	5:E:316:HOH:O	2.58	0.40
1:B:23:ILE:HA	1:B:56:GLU:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	156/179 (87%)	149 (96%)	7 (4%)	0	100	100
1	B	156/179 (87%)	150 (96%)	6 (4%)	0	100	100
1	C	156/179 (87%)	149 (96%)	7 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	163/179 (91%)	155 (95%)	7 (4%)	1 (1%)	21	20
1	E	155/179 (87%)	152 (98%)	3 (2%)	0	100	100
1	F	155/179 (87%)	150 (97%)	5 (3%)	0	100	100
1	G	158/179 (88%)	150 (95%)	8 (5%)	0	100	100
1	H	157/179 (88%)	147 (94%)	10 (6%)	0	100	100
1	I	156/179 (87%)	152 (97%)	4 (3%)	0	100	100
1	J	155/179 (87%)	151 (97%)	4 (3%)	0	100	100
All	All	1567/1790 (88%)	1505 (96%)	61 (4%)	1 (0%)	48	54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	4	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	A	128/144 (89%)	126 (98%)	2 (2%)	55	65
1	B	91/144 (63%)	90 (99%)	1 (1%)	65	74
1	D	99/144 (69%)	98 (99%)	1 (1%)	68	76
1	E	128/144 (89%)	127 (99%)	1 (1%)	73	80
1	F	128/144 (89%)	125 (98%)	3 (2%)	44	53
1	G	131/144 (91%)	126 (96%)	5 (4%)	29	34
1	H	129/144 (90%)	124 (96%)	5 (4%)	28	33
1	I	128/144 (89%)	125 (98%)	3 (2%)	44	53
1	J	128/144 (89%)	127 (99%)	1 (1%)	73	80
All	All	1090/1296 (84%)	1068 (98%)	22 (2%)	50	58

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	142	GLU
1	B	95	SER
1	D	125	LEU
1	E	162	THR
1	F	11	GLN
1	F	85	VAL
1	F	116	LYS
1	G	8	GLN
1	G	29	ARG
1	G	85	VAL
1	G	125	LEU
1	G	131	GLU
1	H	84	VAL
1	H	95	SER
1	H	116	LYS
1	H	146[A]	MET
1	H	146[B]	MET
1	I	11	GLN
1	I	17	LYS
1	I	102	ILE
1	J	125	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	GLN
1	A	118	ASN
1	A	149	HIS
1	A	167	ASN
1	B	75	GLN
1	E	75	GLN
1	E	108	GLN
1	F	75	GLN
1	G	75	GLN
1	H	108	GLN
1	I	149	HIS
1	J	11	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

31 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	D	202	-	4,4,4	0.55	0	6,6,6	0.45	0
2	SO4	C	203	-	4,4,4	0.60	0	6,6,6	0.26	0
2	SO4	E	202	-	4,4,4	0.64	0	6,6,6	0.72	0
2	SO4	F	201	-	4,4,4	0.57	0	6,6,6	0.94	0
4	GOL	G	204	-	5,5,5	1.50	1 (20%)	5,5,5	1.61	2 (40%)
2	SO4	E	201	-	4,4,4	0.46	0	6,6,6	0.33	0
4	GOL	E	203	-	5,5,5	0.87	0	5,5,5	0.86	0
2	SO4	A	201	-	4,4,4	0.56	0	6,6,6	0.69	0
2	SO4	J	203	-	4,4,4	0.59	0	6,6,6	0.27	0
2	SO4	G	202	-	4,4,4	0.51	0	6,6,6	0.31	0
4	GOL	J	204	-	5,5,5	0.61	0	5,5,5	1.58	1 (20%)
2	SO4	J	201	-	4,4,4	0.58	0	6,6,6	0.48	0
2	SO4	G	203	-	4,4,4	0.57	0	6,6,6	0.72	0
4	GOL	I	203	-	5,5,5	0.59	0	5,5,5	1.01	0
3	DLZ	C	204	-	23,24,24	1.65	6 (26%)	30,35,35	1.90	9 (30%)
2	SO4	J	202	-	4,4,4	0.49	0	6,6,6	0.77	0
2	SO4	F	202	-	4,4,4	0.58	0	6,6,6	0.39	0
2	SO4	H	201	-	4,4,4	0.45	0	6,6,6	1.03	0
4	GOL	D	203	-	5,5,5	0.57	0	5,5,5	0.81	0
2	SO4	B	202	-	4,4,4	0.70	0	6,6,6	0.45	0
2	SO4	C	202	-	4,4,4	0.36	0	6,6,6	0.71	0
3	DLZ	I	204	-	23,24,24	1.73	7 (30%)	30,35,35	1.97	5 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	B	201	-	4,4,4	0.64	0	6,6,6	0.34	0
2	SO4	A	202	-	4,4,4	0.41	0	6,6,6	0.63	0
2	SO4	F	203	-	4,4,4	0.25	0	6,6,6	0.49	0
2	SO4	C	201	-	4,4,4	0.45	0	6,6,6	0.57	0
2	SO4	I	202	-	4,4,4	0.35	0	6,6,6	0.49	0
2	SO4	D	201	-	4,4,4	0.68	0	6,6,6	0.85	0
2	SO4	G	201	-	4,4,4	0.62	0	6,6,6	0.82	0
2	SO4	I	201	-	4,4,4	0.58	0	6,6,6	0.54	0
2	SO4	H	202	-	4,4,4	0.46	0	6,6,6	0.41	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
4	GOL	G	204	-	-	2/4/4/4	-
4	GOL	E	203	-	-	0/4/4/4	-
4	GOL	I	203	-	-	1/4/4/4	-
4	GOL	D	203	-	-	2/4/4/4	-
4	GOL	J	204	-	-	0/4/4/4	-
3	DLZ	I	204	-	-	4/14/14/14	0/2/2/2
3	DLZ	C	204	-	-	2/14/14/14	0/2/2/2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	204	DLZ	C7A-N8	-3.74	1.33	1.39
3	I	204	DLZ	C8-N8	-3.16	1.33	1.39
3	I	204	DLZ	C7A-N8	-2.84	1.34	1.39
3	C	204	DLZ	C8-N8	-2.54	1.34	1.39
3	I	204	DLZ	C4A-C8	2.44	1.51	1.44
3	I	204	DLZ	C2'-C3'	2.42	1.57	1.53
3	I	204	DLZ	C1'-C2'	-2.41	1.49	1.52
3	C	204	DLZ	C5A-N5	-2.30	1.34	1.39
3	C	204	DLZ	O2-C2	-2.30	1.19	1.24
3	C	204	DLZ	C2'-C3'	2.25	1.57	1.53
3	I	204	DLZ	C4'-C3'	2.24	1.57	1.53
3	I	204	DLZ	O4-C4	2.16	1.27	1.23
4	G	204	GOL	O3-C3	2.15	1.51	1.42
3	C	204	DLZ	O4-C4	2.03	1.27	1.23

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	204	DLZ	C6-C5A-N5	5.98	121.38	112.45
3	I	204	DLZ	O2'-C2'-C3'	4.44	119.65	109.25
3	C	204	DLZ	O3'-C3'-C4'	-3.99	99.86	108.93
3	C	204	DLZ	C6-C5A-N5	3.62	117.86	112.45
3	C	204	DLZ	C1'-N8-C7A	-3.12	116.71	121.74
3	I	204	DLZ	C7-C7A-C5A	-3.10	119.18	125.26
3	I	204	DLZ	O4'-C4'-C3'	2.94	116.13	109.25
3	C	204	DLZ	O2'-C2'-C3'	2.78	115.75	109.25
4	G	204	GOL	O3-C3-C2	2.71	122.59	110.38
3	C	204	DLZ	O4'-C4'-C3'	2.54	115.20	109.25
3	C	204	DLZ	C4A-C8-N1	-2.42	118.65	124.59
3	I	204	DLZ	C4-C4A-N5	2.27	121.32	118.18
4	J	204	GOL	C3-C2-C1	-2.23	103.61	111.80
4	G	204	GOL	O2-C2-C3	2.21	118.34	109.18
3	C	204	DLZ	N3-C2-N1	2.12	124.00	119.50
3	C	204	DLZ	O4-C4-C4A	-2.09	121.01	126.53
3	C	204	DLZ	O4'-C4'-C5'	2.09	113.78	109.03

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	I	204	DLZ	C1'-C2'-C3'-O3'
3	I	204	DLZ	C1'-C2'-C3'-C4'
4	G	204	GOL	C1-C2-C3-O3
3	I	204	DLZ	O2'-C2'-C3'-C4'
4	G	204	GOL	O2-C2-C3-O3
3	I	204	DLZ	O2'-C2'-C3'-O3'
4	D	203	GOL	O1-C1-C2-C3
4	D	203	GOL	O1-C1-C2-O2
3	C	204	DLZ	O2'-C2'-C3'-O3'
4	I	203	GOL	O2-C2-C3-O3
3	C	204	DLZ	C1'-C2'-C3'-O3'

There are no ring outliers.

8 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	203	GOL	2	0
2	A	201	SO4	1	0

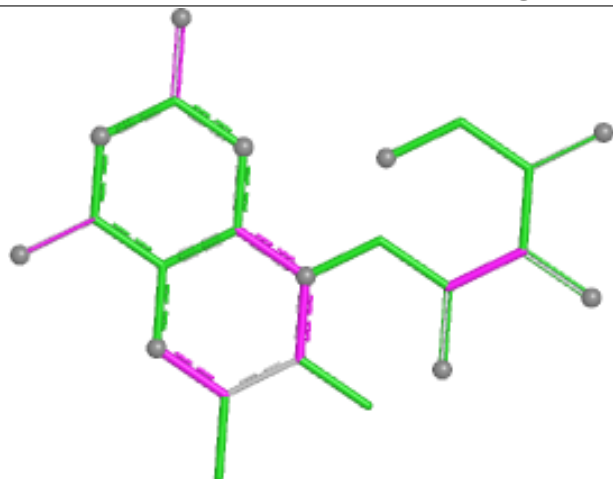
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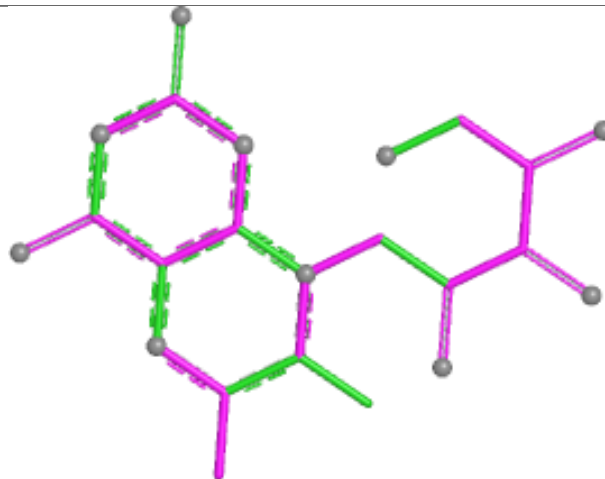
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	204	GOL	3	0
4	I	203	GOL	1	0
3	C	204	DLZ	4	0
4	D	203	GOL	2	0
3	I	204	DLZ	3	0
2	G	201	SO4	1	0

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.

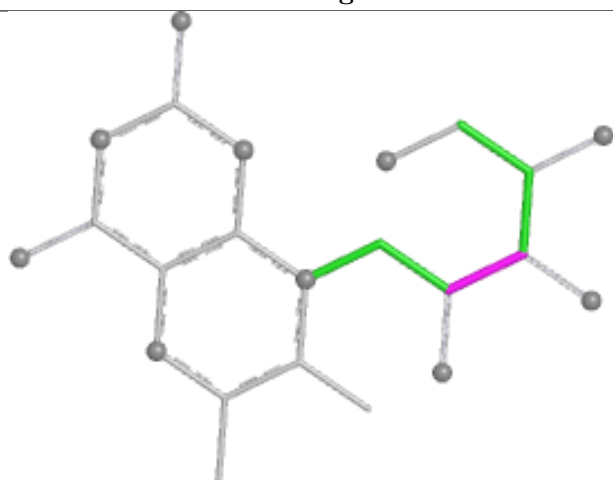
Ligand DLZ C 204



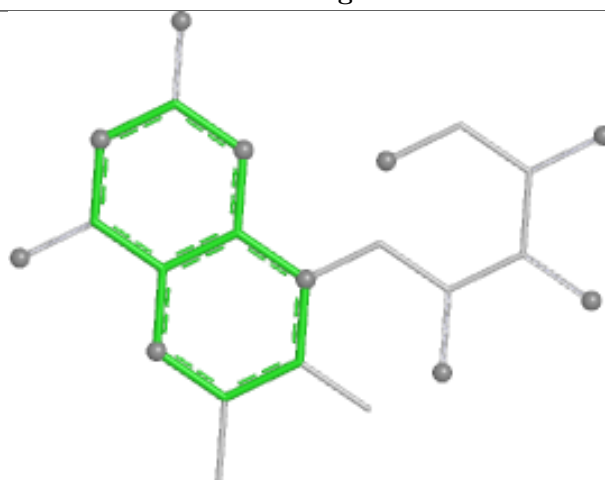
Bond lengths



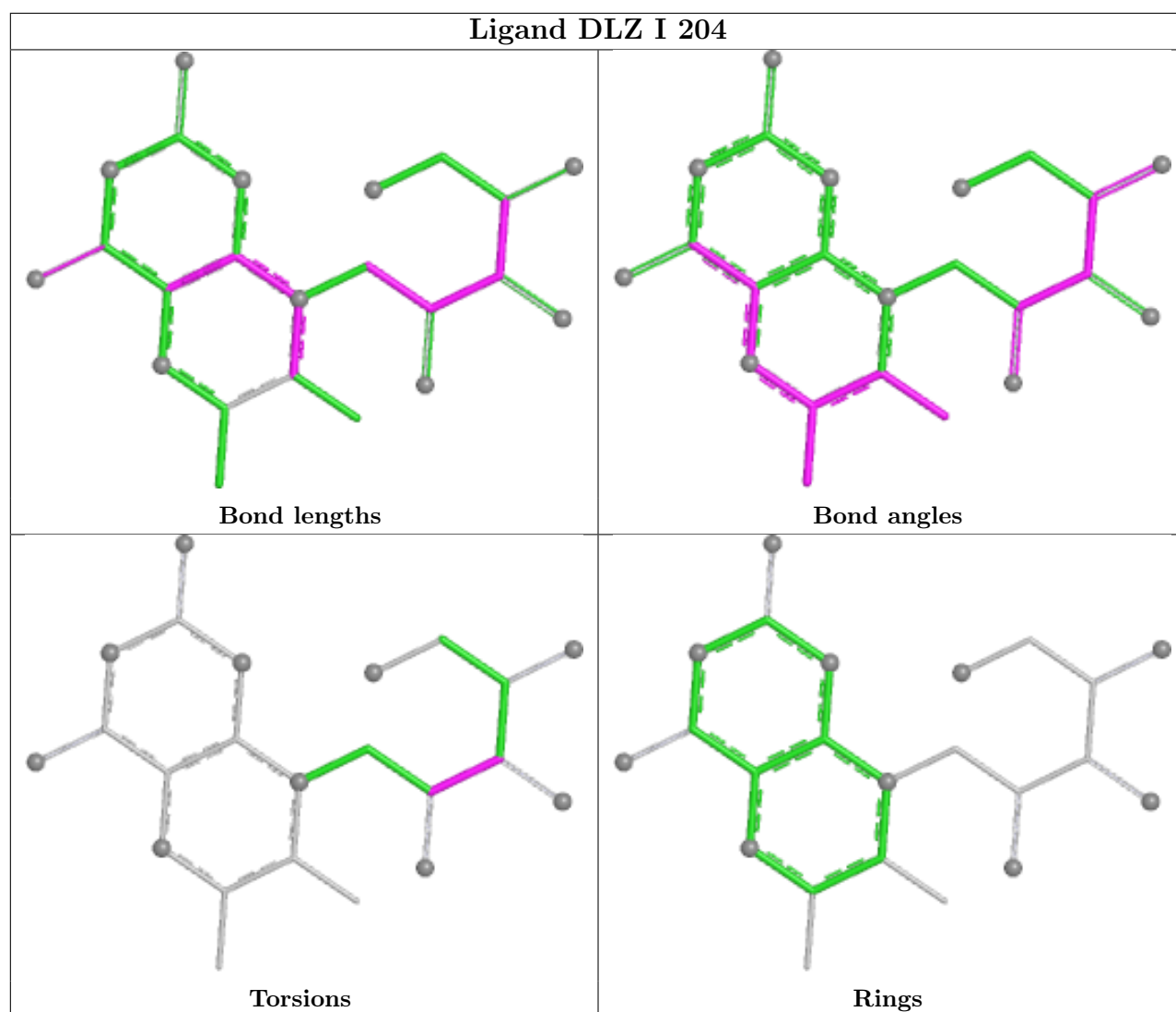
Bond angles



Torsions



Rings



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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	158/179 (88%)	-0.55	5 (3%)	50	50	8, 16, 40, 78	2 (1%)
1	B	158/179 (88%)	-0.52	4 (2%)	58	59	7, 15, 34, 60	6 (3%)
1	C	158/179 (88%)	-0.46	6 (3%)	44	43	9, 15, 36, 71	4 (2%)
1	D	165/179 (92%)	-0.34	7 (4%)	40	39	8, 15, 32, 59	10 (6%)
1	E	157/179 (87%)	-0.59	3 (1%)	66	68	8, 13, 33, 71	2 (1%)
1	F	157/179 (87%)	-0.53	4 (2%)	58	59	9, 14, 32, 41	4 (2%)
1	G	160/179 (89%)	-0.41	4 (2%)	58	59	10, 18, 36, 56	6 (3%)
1	H	158/179 (88%)	-0.14	8 (5%)	33	32	12, 22, 39, 70	8 (5%)
1	I	158/179 (88%)	-0.28	6 (3%)	44	43	11, 20, 44, 72	3 (1%)
1	J	157/179 (87%)	-0.46	3 (1%)	66	68	8, 17, 38, 73	3 (1%)
All	All	1586/1790 (88%)	-0.43	50 (3%)	50	50	7, 17, 37, 78	48 (3%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	12	ASN	12.7
1	F	167	ASN	10.4
1	I	167	ASN	10.4
1	D	167	ASN	10.1
1	F	12	ASN	10.1
1	D	3	VAL	9.9
1	J	167	ASN	9.5
1	H	167	ASN	9.0
1	H	11	GLN	8.5
1	E	12	ASN	7.9
1	B	11	GLN	7.8
1	C	11	GLN	7.3
1	G	167	ASN	7.2

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Mol	Chain	Res	Type	RSRZ
1	D	4	LYS	7.2
1	E	167	ASN	5.6
1	C	168	ALA	5.4
1	I	168	ALA	5.2
1	D	9	LEU	4.9
1	H	142	GLU	4.7
1	H	168	ALA	4.7
1	G	142	GLU	4.5
1	B	142	GLU	4.4
1	G	33	ASP	4.2
1	A	168	ALA	4.2
1	I	144	LYS	4.0
1	F	11	GLN	3.9
1	B	168	ALA	3.7
1	H	27	TRP	3.6
1	D	6	LEU	3.6
1	H	146[A]	MET	3.4
1	I	50	GLU	3.4
1	H	50	GLU	3.3
1	I	27	TRP	3.3
1	J	11	GLN	3.2
1	D	10	ASP	3.2
1	C	17	LYS	3.1
1	C	142	GLU	3.0
1	A	12	ASN	3.0
1	A	11	GLN	2.9
1	D	5	GLY	2.8
1	G	8	GLN	2.8
1	J	50	GLU	2.8
1	H	131	GLU	2.6
1	B	144	LYS	2.5
1	C	167	ASN	2.2
1	I	11	GLN	2.2
1	A	50	GLU	2.1
1	F	50	GLU	2.1
1	E	11	GLN	2.1
1	A	167	ASN	2.0

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 oligosaccharides in this entry.

6.4 Ligands ⓘ

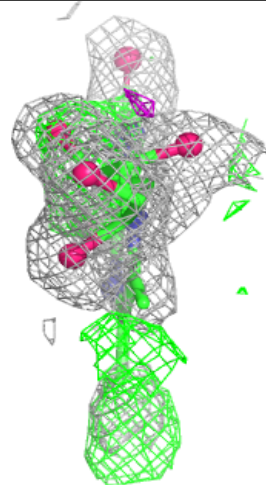
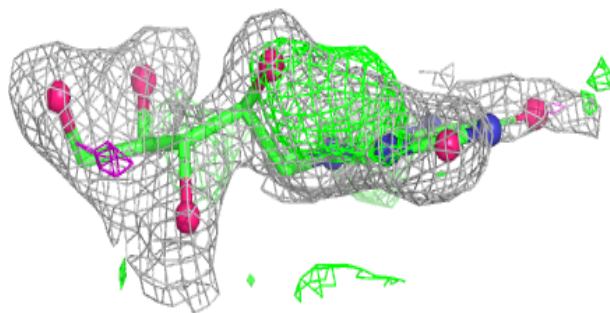
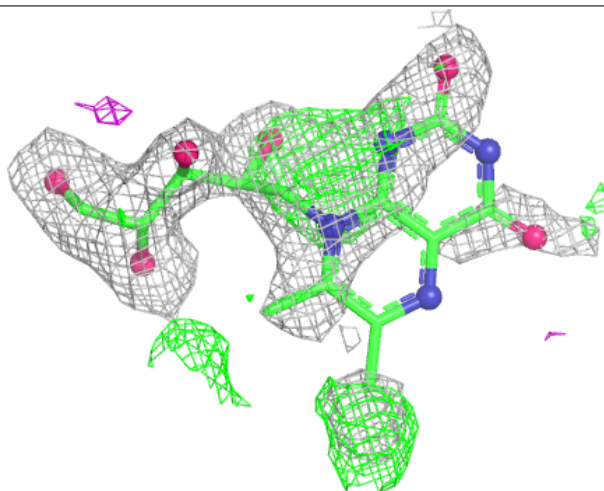
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	J	203	5/5	0.74	0.15	80,89,94,97	0
2	SO4	I	201	5/5	0.75	0.14	66,71,78,81	0
3	DLZ	I	204	23/23	0.75	0.24	16,28,34,39	17
4	GOL	G	204	6/6	0.78	0.16	32,33,36,37	0
3	DLZ	C	204	23/23	0.79	0.19	9,19,29,35	17
2	SO4	G	202	5/5	0.80	0.13	67,74,77,81	0
2	SO4	G	201	5/5	0.84	0.15	56,59,69,70	0
2	SO4	C	201	5/5	0.85	0.13	68,68,73,74	0
2	SO4	B	201	5/5	0.85	0.15	54,57,62,68	0
4	GOL	D	203	6/6	0.87	0.12	29,30,33,35	0
2	SO4	J	201	5/5	0.92	0.15	56,56,64,69	0
2	SO4	H	202	5/5	0.92	0.12	50,54,58,66	0
4	GOL	E	203	6/6	0.92	0.12	28,29,33,33	0
2	SO4	D	201	5/5	0.92	0.12	41,47,54,56	0
4	GOL	I	203	6/6	0.92	0.10	32,33,36,37	0
2	SO4	F	202	5/5	0.93	0.12	41,47,53,58	0
2	SO4	E	202	5/5	0.94	0.13	31,37,47,49	0
2	SO4	F	201	5/5	0.94	0.15	36,43,49,56	0
2	SO4	A	201	5/5	0.94	0.11	39,44,47,49	0
4	GOL	J	204	6/6	0.94	0.07	22,25,26,27	0
2	SO4	H	201	5/5	0.95	0.08	35,40,42,50	0
2	SO4	C	203	5/5	0.96	0.11	42,43,55,57	0
2	SO4	G	203	5/5	0.97	0.07	34,37,40,45	0
2	SO4	B	202	5/5	0.98	0.06	32,33,35,39	0
2	SO4	J	202	5/5	0.98	0.08	24,28,31,33	0
2	SO4	A	202	5/5	0.98	0.05	28,29,34,35	0
2	SO4	E	201	5/5	0.98	0.06	30,32,34,38	0
2	SO4	I	202	5/5	0.98	0.06	30,30,32,34	0
2	SO4	F	203	5/5	0.99	0.06	32,34,36,40	0
2	SO4	C	202	5/5	0.99	0.06	27,27,30,32	0
2	SO4	D	202	5/5	0.99	0.06	25,26,29,30	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.

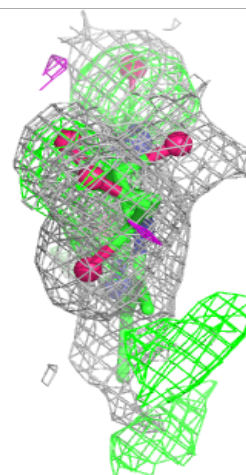
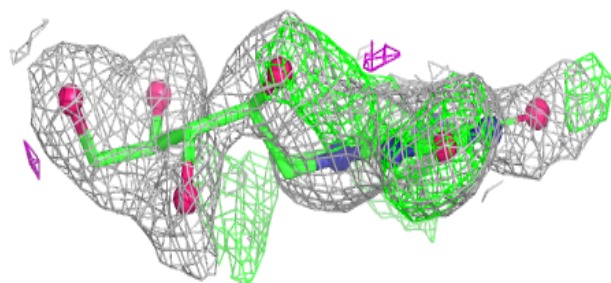
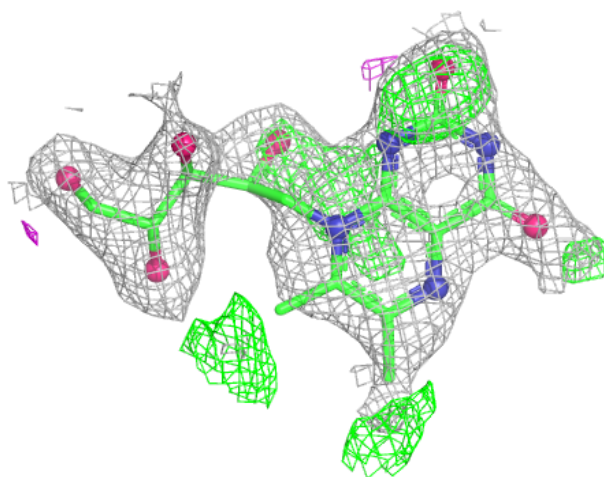
Electron density around DLZ I 204:

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



Electron density around DLZ C 204:

$2mF_o-DF_c$ (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.