



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 09:43 pm BST

PDB ID : 4KQ6
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

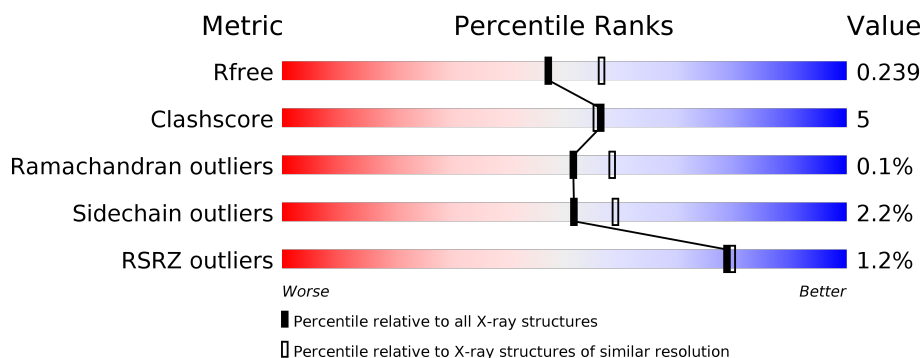
1 Overall quality at a glance

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}	130704	2391 (2.26-2.22)
Clashscore	141614	2539 (2.26-2.22)
Ramachandran outliers	138981	2489 (2.26-2.22)
Sidechain outliers	138945	2490 (2.26-2.22)
RSRZ outliers	127900	2353 (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 on 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	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 80%, grey 12%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 80% 7% • 12% </div> </div>
1	B	179	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 2%, yellow 11%, green 77%, grey 12%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 77% 11% • 12% </div> </div>
1	C	179	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 13%, green 74%, grey 12%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 74% 13% • 12% </div> </div>
1	D	179	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, orange 2%, yellow 11%, green 80%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 80% 11% •• 8% </div> </div>
1	E	179	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 7%, green 80%, grey 12%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 80% 7% 12% </div> </div>
1	F	179	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 7%, green 81%, grey 12%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 81% 7% 12% </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	179	<div><div></div><div>78%</div><div>11%</div><div>•</div><div>11%</div></div>
1	H	179	<div><div>%</div><div></div><div>72%</div><div>15%</div><div>•</div><div>12%</div></div>
1	I	179	<div><div>2%</div><div></div><div>77%</div><div>11%</div><div>•</div><div>12%</div></div>
1	J	179	<div><div></div><div>73%</div><div>14%</div><div>•</div><div>12%</div></div>

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 (three-letter code: SO4) (formula: O₄S).



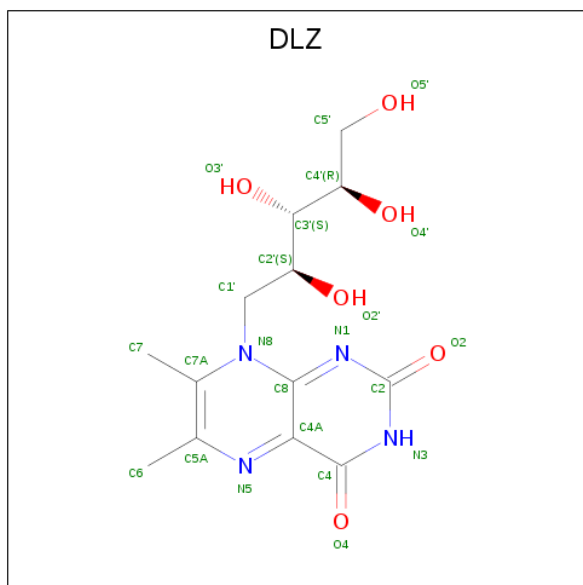
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	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-ribitol (three-letter code: DLZ) (formula: C₁₃H₁₈N₄O₆).



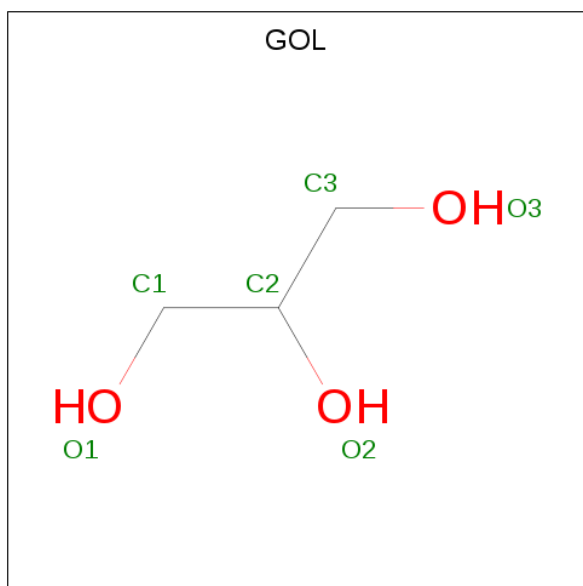
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			23	13	4	6		

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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 (three-letter code: 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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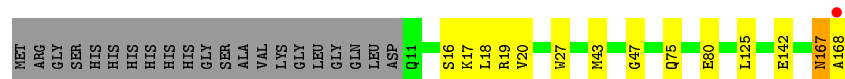
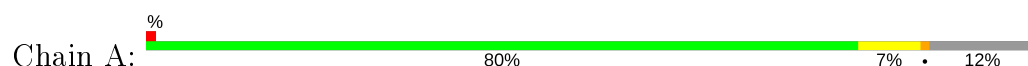
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	103	Total 103	O 103	0	0
5	E	107	Total 107	O 107	0	0
5	F	103	Total 103	O 103	0	0
5	G	70	Total 70	O 70	0	0
5	H	66	Total 66	O 66	0	0
5	I	75	Total 75	O 75	0	0
5	J	88	Total 88	O 88	0	0

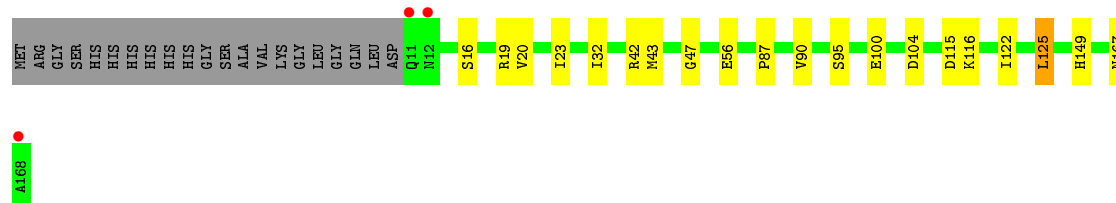
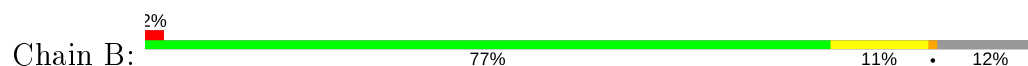
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

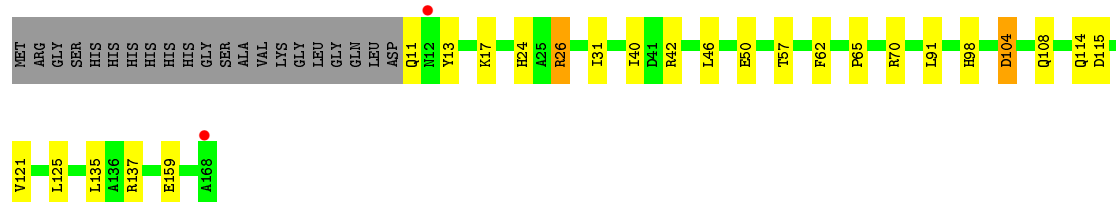
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



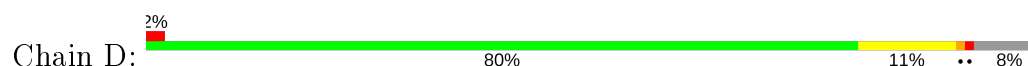
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



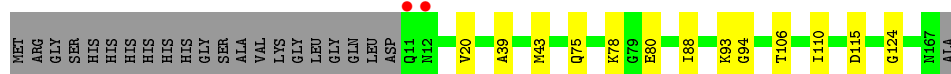
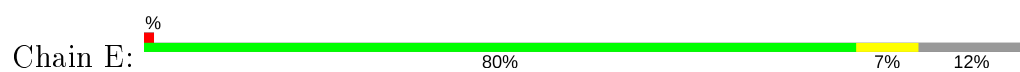
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



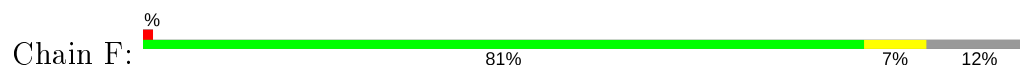
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



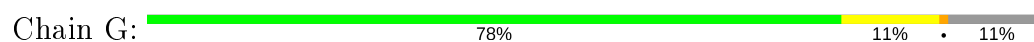
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



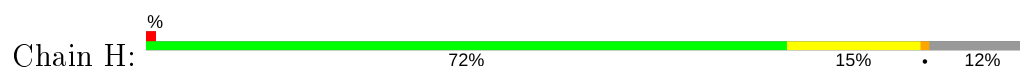
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



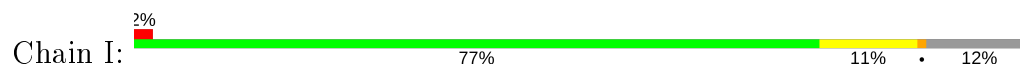
- Molecule 1: 6,7-dimethyl-8-ribityllumazine synthase



- 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 0.190 , 0.239	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)	0.34 , 39.3	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	0.93	0/1232	0.97	4/1663 (0.2%)
1	B	0.94	0/1232	0.99	3/1663 (0.2%)
1	C	0.94	0/1233	1.07	7/1663 (0.4%)
1	D	0.99	0/1284	1.02	3/1732 (0.2%)
1	E	1.03	0/1227	0.97	1/1656 (0.1%)
1	F	1.00	0/1227	0.99	5/1656 (0.3%)
1	G	0.91	0/1252	0.99	3/1690 (0.2%)
1	H	0.89	0/1240	0.95	4/1673 (0.2%)
1	I	0.94	0/1233	0.92	0/1663
1	J	0.92	1/1227 (0.1%)	0.95	3/1656 (0.2%)
All	All	0.95	1/12387 (0.0%)	0.98	33/16715 (0.2%)

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 (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	153	TRP	CE3-CZ3	6.25	1.49	1.38

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	26	ARG	NE-CZ-NH2	-11.66	114.47	120.30
1	C	26	ARG	NE-CZ-NH1	8.57	124.59	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	19	ARG	NE-CZ-NH1	7.68	124.14	120.30
1	A	125	LEU	CA-CB-CG	7.41	132.34	115.30
1	B	115	ASP	CB-CG-OD1	7.33	124.90	118.30
1	G	19	ARG	NE-CZ-NH2	-7.27	116.67	120.30
1	H	26	ARG	NE-CZ-NH2	-7.00	116.80	120.30
1	E	115	ASP	CB-CG-OD1	6.72	124.35	118.30
1	J	70	ARG	NE-CZ-NH1	6.55	123.57	120.30
1	D	6	LEU	CA-CB-CG	6.18	129.51	115.30
1	B	19	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	F	19	ARG	NE-CZ-NH1	6.06	123.33	120.30
1	F	26	ARG	NE-CZ-NH2	-6.01	117.30	120.30
1	C	104	ASP	CB-CG-OD2	-5.98	112.92	118.30
1	C	115	ASP	CB-CG-OD1	5.87	123.58	118.30
1	D	6	LEU	CB-CA-C	-5.73	99.31	110.20
1	H	125	LEU	CA-CB-CG	5.64	128.27	115.30
1	A	18	LEU	CB-CG-CD1	-5.62	101.45	111.00
1	A	19	ARG	NE-CZ-NH2	-5.58	117.51	120.30
1	J	125	LEU	CA-CB-CG	5.55	128.06	115.30
1	C	125	LEU	CA-CB-CG	5.51	127.98	115.30
1	A	19	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	B	125	LEU	CA-CB-CG	5.47	127.88	115.30
1	J	19	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	C	137	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	F	115	ASP	CB-CG-OD1	5.38	123.15	118.30
1	C	70	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	F	19	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	H	70	ARG	NE-CZ-NH1	5.34	122.97	120.30
1	G	104	ASP	CB-CG-OD2	-5.23	113.59	118.30
1	F	115	ASP	CB-CG-OD2	-5.18	113.64	118.30
1	D	83	ASP	CB-CG-OD2	5.17	122.95	118.30
1	H	44	LEU	CA-CB-CG	-5.10	103.57	115.30

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	22	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	17	0
1	J	1207	0	1224	16	0
2	A	10	0	0	1	0
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	5	0
3	I	23	0	18	5	0
4	D	6	0	8	3	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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	13313	0	12435	123	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 (123) 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:D:4:LYS:N	1:D:4:LYS:HD3	1.57	1.10
1:C:98:HIS:CD2	3:C:204:DLZ:H6A	1.97	0.97
1:D:4:LYS:H	1:D:4:LYS:HD3	1.44	0.79
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
1:D:3:VAL:HG13	1:D:4:LYS:N	2.09	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:C6	2.78	0.66
1:C:98:HIS:NE2	3:C:204:DLZ:H7B	2.10	0.66
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: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: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: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:A:167:ASN:O	1:A:168:ALA:C	2.45	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:D:4:LYS:N	1:D:4:LYS:CD	2.47	0.51
1:F:20:VAL:HG11	1:F:43:MET:SD	2.50	0.51
1:D:3:VAL:HG22	1:D:4:LYS:CE	2.40	0.51
1:I:98:HIS:CD2	3:I:204:DLZ:H6A	2.46	0.51
1:B:20:VAL:HG11	1:B:43:MET:SD	2.51	0.50
1:J:11:GLN:N	1:J:11:GLN:CD	2.65	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:D:3:VAL:HG13	1:D:4:LYS:H	1.76	0.50
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
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: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:A:20:VAL:HG11	1:A:43:MET:SD	2.54	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:HD12	1:J:122:ILE:HG23	1.73	0.46
1:H:85:VAL:CG1	1:H:121:VAL:HG22	2.46	0.46
1:G:159:GLU:OE2	1:G:163:LYS:NZ	2.44	0.46
1:D:3:VAL:C	1:D:4:LYS:HD3	2.31	0.46
1:D:63:GLU:OE2	4:D:203:GOL:O1	2.30	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:98:HIS:NE2	3:I:204:DLZ:H6A	2.31	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
3:C:204:DLZ:H7	3:C:204:DLZ:H1'A	1.57	0.43
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:HA	1:H:44:LEU:HD23	1.70	0.43
1:H:39:ALA:O	1:H:43:MET:HG3	2.19	0.42
1:D:3:VAL:HG22	1:D:4:LYS:NZ	2.34	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
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:N	1:D:141:ASP:OD1	2.53	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:E:106:THR:O	1:E:110:ILE:HG13	2.21	0.41
1:E:20:VAL:HG11	1:E:43:MET:SD	2.61	0.41
1:G:108:GLN:OE1	1:H:108:GLN:NE2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:OD1	1:J:141:ASP:N	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

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	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
1	D	163/179 (91%)	155 (95%)	7 (4%)	1 (1%)	25	23
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%)	51	58

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%)	125 (98%)	3 (2%)	50	57
1	B	128/144 (89%)	123 (96%)	5 (4%)	32	35
1	C	128/144 (89%)	127 (99%)	1 (1%)	81	87
1	D	134/144 (93%)	129 (96%)	5 (4%)	34	38
1	E	128/144 (89%)	128 (100%)	0	100	100
1	F	128/144 (89%)	125 (98%)	3 (2%)	50	57
1	G	131/144 (91%)	126 (96%)	5 (4%)	33	36
1	H	129/144 (90%)	125 (97%)	4 (3%)	40	46
1	I	128/144 (89%)	126 (98%)	2 (2%)	62	70
1	J	128/144 (89%)	127 (99%)	1 (1%)	81	87
All	All	1290/1440 (90%)	1261 (98%)	29 (2%)	52	59

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	LYS
1	A	142	GLU
1	A	167	ASN
1	B	95	SER
1	B	116	LYS
1	B	125	LEU
1	B	149	HIS
1	B	167	ASN
1	C	11	GLN
1	D	3	VAL
1	D	4	LYS
1	D	6	LEU
1	D	9	LEU

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Mol	Chain	Res	Type
1	D	125	LEU
1	F	11	GLN
1	F	95	SER
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	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	J	125	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	149	HIS
1	C	108	GLN
1	E	108	GLN
1	I	108	GLN
1	I	149	HIS

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 carbohydrates 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	I	201	-	4,4,4	0.50	0	6,6,6	0.54	0
2	SO4	A	202	-	4,4,4	0.33	0	6,6,6	0.77	0
2	SO4	C	202	-	4,4,4	0.27	0	6,6,6	0.68	0
4	GOL	G	204	-	5,5,5	1.47	1 (20%)	5,5,5	1.53	2 (40%)
3	DLZ	C	204	-	21,24,24	3.02	6 (28%)	22,35,35	3.03	9 (40%)
2	SO4	G	203	-	4,4,4	0.48	0	6,6,6	0.87	0
3	DLZ	I	204	-	21,24,24	3.61	9 (42%)	22,35,35	3.38	10 (45%)
2	SO4	D	202	-	4,4,4	0.47	0	6,6,6	0.45	0
2	SO4	I	202	-	4,4,4	0.24	0	6,6,6	0.55	0
2	SO4	E	201	-	4,4,4	0.38	0	6,6,6	0.39	0
4	GOL	J	204	-	5,5,5	0.60	0	5,5,5	1.47	1 (20%)
2	SO4	E	202	-	4,4,4	0.57	0	6,6,6	0.81	0
2	SO4	J	201	-	4,4,4	0.50	0	6,6,6	0.45	0
2	SO4	F	202	-	4,4,4	0.49	0	6,6,6	0.41	0
2	SO4	J	202	-	4,4,4	0.39	0	6,6,6	0.86	0
4	GOL	D	203	-	5,5,5	0.56	0	5,5,5	0.77	0
4	GOL	E	203	-	5,5,5	0.84	0	5,5,5	0.81	0
2	SO4	B	202	-	4,4,4	0.62	0	6,6,6	0.48	0
2	SO4	J	203	-	4,4,4	0.50	0	6,6,6	0.24	0
2	SO4	A	201	-	4,4,4	0.48	0	6,6,6	0.66	0
2	SO4	F	203	-	4,4,4	0.21	0	6,6,6	0.48	0
2	SO4	H	201	-	4,4,4	0.36	0	6,6,6	1.04	0
4	GOL	I	203	-	5,5,5	0.57	0	5,5,5	0.94	0
2	SO4	B	201	-	4,4,4	0.57	0	6,6,6	0.42	0
2	SO4	C	201	-	4,4,4	0.35	0	6,6,6	0.58	0
2	SO4	F	201	-	4,4,4	0.49	0	6,6,6	0.97	0
2	SO4	D	201	-	4,4,4	0.60	0	6,6,6	0.86	0
2	SO4	G	201	-	4,4,4	0.53	0	6,6,6	0.81	0
2	SO4	G	202	-	4,4,4	0.42	0	6,6,6	0.35	0
2	SO4	H	202	-	4,4,4	0.37	0	6,6,6	0.44	0
2	SO4	C	203	-	4,4,4	0.51	0	6,6,6	0.33	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	E	203	-	-	0/4/4/4	-
4	GOL	G	204	-	-	2/4/4/4	-
3	DLZ	C	204	-	-	2/14/14/14	0/2/2/2
4	GOL	J	204	-	-	0/4/4/4	-
3	DLZ	I	204	-	-	4/14/14/14	0/2/2/2
4	GOL	I	203	-	-	1/4/4/4	-
4	GOL	D	203	-	-	2/4/4/4	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	204	DLZ	C4A-C8	12.54	1.51	1.38
3	C	204	DLZ	C4A-C8	11.01	1.49	1.38
3	I	204	DLZ	C1'-N8	-5.35	1.42	1.48
3	I	204	DLZ	C4-C4A	4.45	1.49	1.41
3	C	204	DLZ	C4-C4A	4.37	1.48	1.41
3	I	204	DLZ	C5A-N5	4.30	1.37	1.32
3	C	204	DLZ	C1'-N8	-3.50	1.44	1.48
3	I	204	DLZ	C7A-C5A	3.22	1.50	1.41
3	I	204	DLZ	C8-N1	3.00	1.37	1.33
3	C	204	DLZ	C7A-C5A	2.91	1.49	1.41
3	I	204	DLZ	C4A-N5	2.85	1.37	1.33
3	C	204	DLZ	C8-N1	2.82	1.36	1.33
3	I	204	DLZ	C2'-C3'	2.22	1.57	1.53
4	G	204	GOL	O3-C3	2.14	1.51	1.42
3	C	204	DLZ	C2'-C3'	2.06	1.57	1.53
3	I	204	DLZ	C4'-C3'	2.06	1.57	1.53

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	204	DLZ	C4-N3-C2	10.95	124.39	115.14
3	C	204	DLZ	C4-N3-C2	9.12	122.84	115.14
3	I	204	DLZ	C4A-C4-N3	-5.47	115.95	123.43
3	C	204	DLZ	C4A-C4-N3	-5.00	116.59	123.43
3	I	204	DLZ	C4-C4A-C8	-4.49	116.98	119.95
3	I	204	DLZ	O2'-C2'-C3'	4.34	119.65	109.10
3	C	204	DLZ	O3'-C3'-C4'	-3.71	99.86	108.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	204	DLZ	C1'-N8-C8	3.55	121.59	118.41
3	C	204	DLZ	C4-C4A-C8	-3.19	117.84	119.95
3	I	204	DLZ	O2'-C2'-C1'	-3.13	102.05	109.59
3	C	204	DLZ	C4A-C8-N8	-2.96	117.26	120.30
3	I	204	DLZ	C6-C5A-N5	2.94	121.38	116.97
3	I	204	DLZ	O4'-C4'-C3'	2.89	116.13	109.10
3	I	204	DLZ	C7A-C5A-N5	-2.88	118.31	122.23
3	C	204	DLZ	C7A-C5A-N5	-2.75	118.49	122.23
3	C	204	DLZ	O2'-C2'-C3'	2.73	115.75	109.10
3	I	204	DLZ	C4A-C8-N8	-2.73	117.49	120.30
4	G	204	GOL	O3-C3-C2	2.58	122.59	110.20
3	C	204	DLZ	O4'-C4'-C3'	2.51	115.20	109.10
3	I	204	DLZ	C4-C4A-N5	2.38	121.32	118.60
4	G	204	GOL	O2-C2-C3	2.09	118.34	109.12
4	J	204	GOL	C3-C2-C1	-2.08	103.61	111.70

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	G	204	GOL	C1-C2-C3-O3
3	I	204	DLZ	C1'-C2'-C3'-O3'
3	I	204	DLZ	C1'-C2'-C3'-C4'
4	G	204	GOL	O2-C2-C3-O3
3	I	204	DLZ	O2'-C2'-C3'-C4'
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 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	204	DLZ	5	0
3	I	204	DLZ	5	0
4	J	204	GOL	3	0
4	D	203	GOL	3	0
4	E	203	GOL	2	0

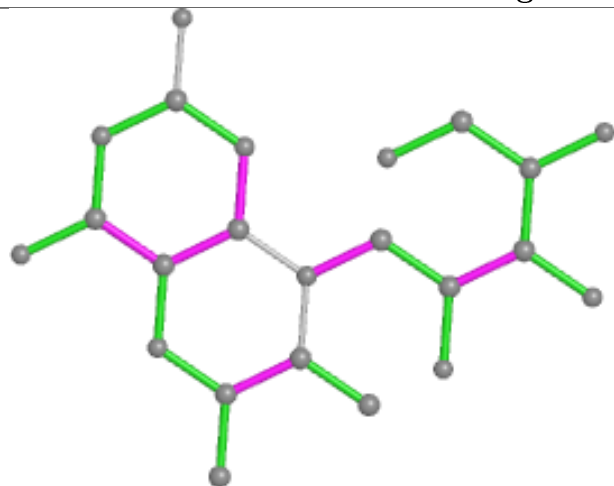
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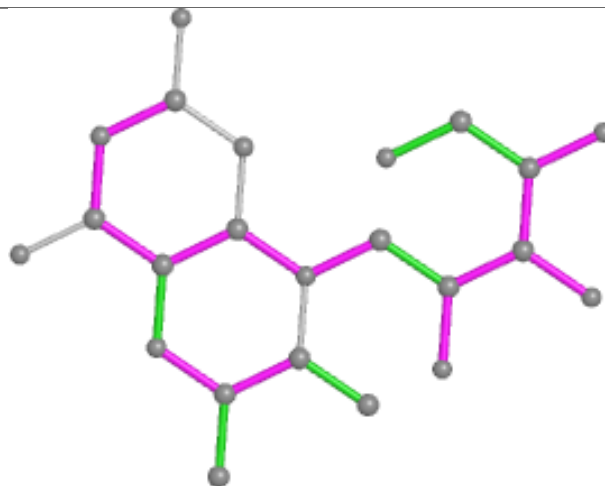
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201	SO4	1	0
4	I	203	GOL	1	0
2	G	201	SO4	1	0

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.

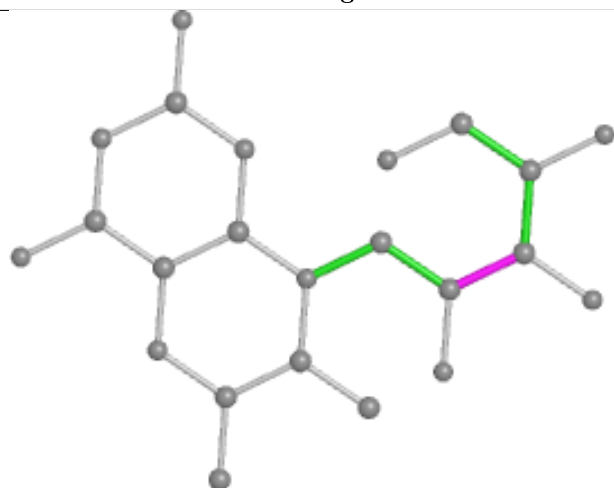
Ligand DLZ C 204



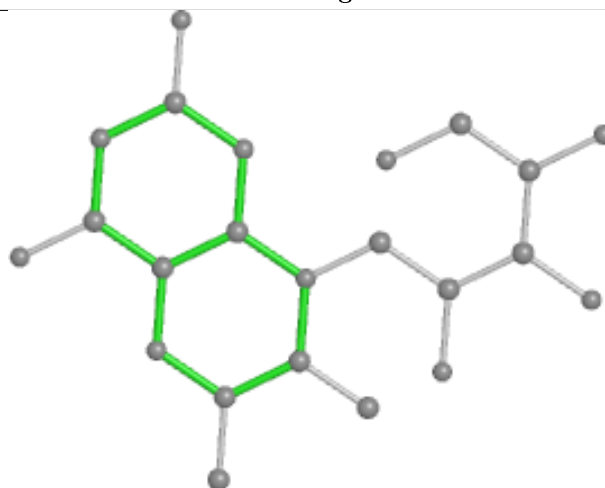
Bond lengths



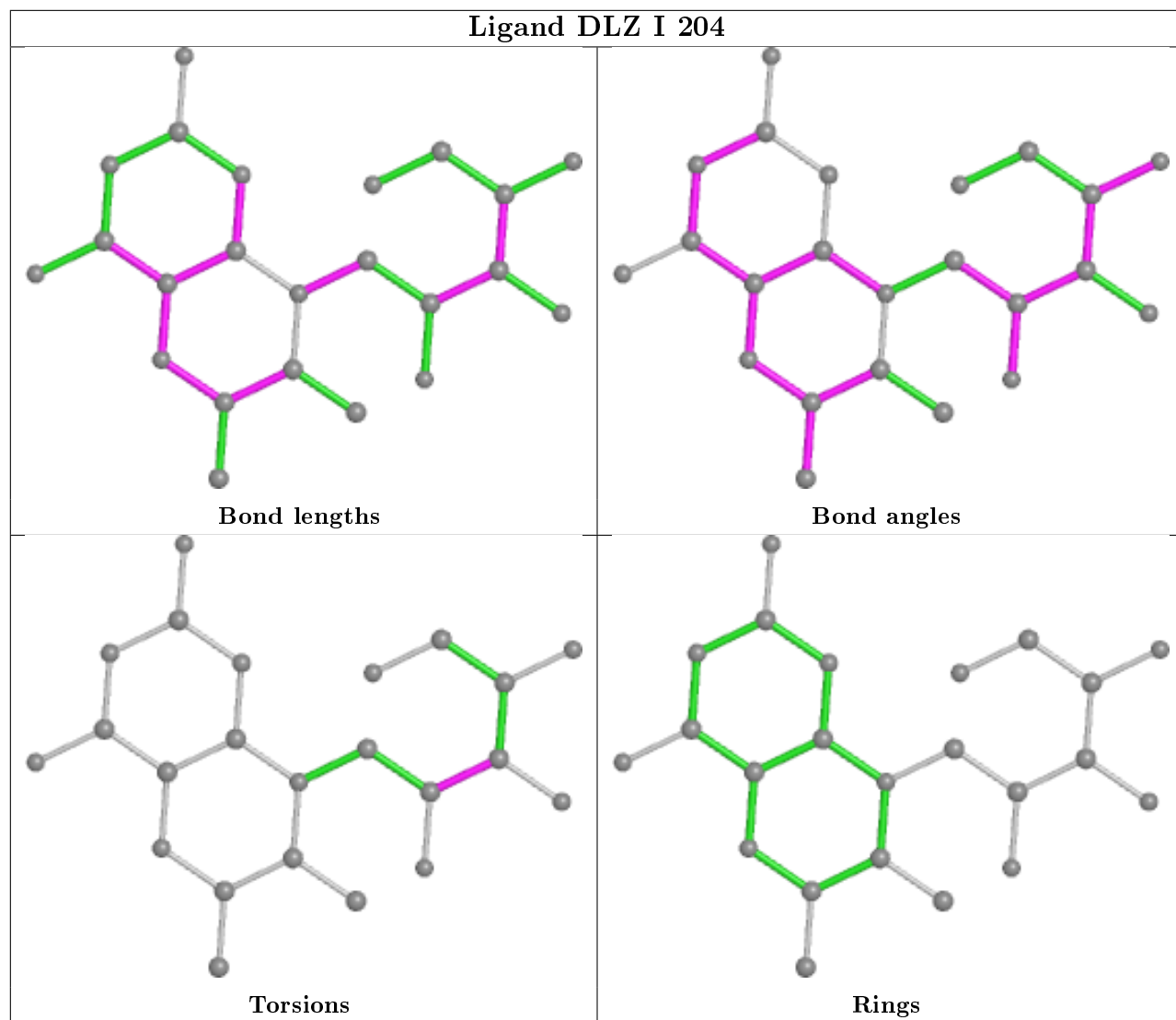
Bond angles



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%)	-0.60	1 (0%) 89 89	8, 16, 40, 77	2 (1%)
1	B	158/179 (88%)	-0.62	3 (1%) 66 68	9, 15, 37, 60	6 (3%)
1	C	158/179 (88%)	-0.52	2 (1%) 77 78	9, 15, 38, 71	4 (2%)
1	D	165/179 (92%)	-0.47	4 (2%) 59 60	8, 15, 32, 59	10 (6%)
1	E	157/179 (87%)	-0.57	2 (1%) 77 78	8, 13, 33, 71	2 (1%)
1	F	157/179 (87%)	-0.57	2 (1%) 77 78	9, 14, 32, 61	4 (2%)
1	G	160/179 (89%)	-0.58	0 100 100	10, 19, 40, 56	6 (3%)
1	H	158/179 (88%)	-0.36	2 (1%) 77 78	12, 22, 43, 70	7 (4%)
1	I	158/179 (88%)	-0.44	3 (1%) 66 68	11, 20, 45, 72	3 (1%)
1	J	157/179 (87%)	-0.54	0 100 100	8, 17, 38, 73	3 (1%)
All	All	1586/1790 (88%)	-0.53	19 (1%) 79 80	8, 17, 39, 77	47 (2%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	9	LEU	5.0
1	C	12	ASN	4.2
1	D	3	VAL	4.0
1	F	12	ASN	3.8
1	D	4	LYS	3.8
1	E	11	GLN	3.3
1	C	168	ALA	3.3
1	E	12	ASN	3.1
1	I	168	ALA	3.0
1	A	168	ALA	2.9
1	B	168	ALA	2.9
1	F	167	ASN	2.6
1	D	10	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	H	168	ALA	2.6
1	H	12	ASN	2.6
1	I	167	ASN	2.4
1	I	143	GLY	2.2
1	B	11	GLN	2.2
1	B	12	ASN	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	DLZ	I	204	23/23	0.70	0.38	16,28,34,39	17
4	GOL	G	204	6/6	0.73	0.19	32,33,36,37	0
3	DLZ	C	204	23/23	0.78	0.36	9,19,29,35	17
2	SO4	J	203	5/5	0.80	0.28	80,89,94,97	0
4	GOL	E	203	6/6	0.88	0.17	28,29,33,33	0
2	SO4	I	201	5/5	0.90	0.23	66,71,78,81	0
4	GOL	D	203	6/6	0.90	0.15	29,30,33,35	0
4	GOL	I	203	6/6	0.91	0.14	32,33,36,37	0
2	SO4	G	201	5/5	0.91	0.23	56,59,69,70	0
2	SO4	C	201	5/5	0.92	0.18	68,68,73,74	0
2	SO4	B	201	5/5	0.92	0.16	54,57,62,68	0
2	SO4	G	202	5/5	0.93	0.24	67,74,77,81	0
4	GOL	J	204	6/6	0.94	0.13	22,25,26,27	0
2	SO4	E	202	5/5	0.95	0.17	31,37,47,49	0
2	SO4	H	202	5/5	0.95	0.14	50,54,58,66	0
2	SO4	D	201	5/5	0.96	0.13	41,47,54,56	0
2	SO4	A	201	5/5	0.96	0.17	39,44,47,49	0

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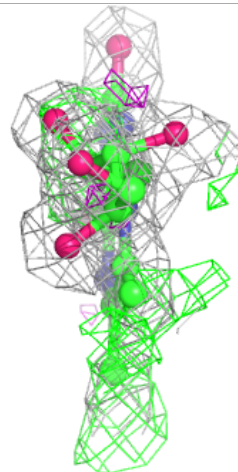
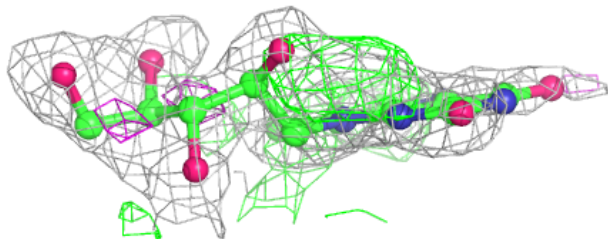
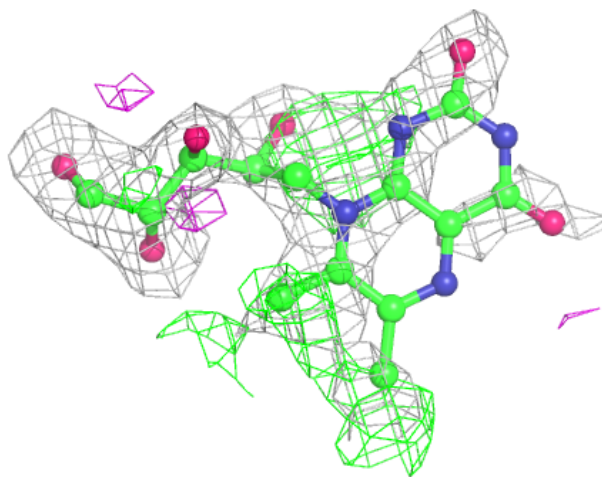
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	F	202	5/5	0.96	0.13	41,47,53,58	0
2	SO4	F	201	5/5	0.96	0.19	36,43,49,56	0
2	SO4	J	201	5/5	0.97	0.16	56,56,64,69	0
2	SO4	E	201	5/5	0.98	0.10	30,32,34,38	0
2	SO4	J	202	5/5	0.98	0.09	24,28,31,33	0
2	SO4	G	203	5/5	0.98	0.11	34,37,40,45	0
2	SO4	H	201	5/5	0.98	0.12	35,40,42,50	0
2	SO4	A	202	5/5	0.99	0.07	28,29,34,35	0
2	SO4	C	202	5/5	0.99	0.09	27,27,30,32	0
2	SO4	F	203	5/5	0.99	0.10	32,34,36,40	0
2	SO4	D	202	5/5	0.99	0.09	25,26,29,30	0
2	SO4	I	202	5/5	0.99	0.11	30,30,32,34	0
2	SO4	B	202	5/5	0.99	0.08	32,33,35,39	0
2	SO4	C	203	5/5	0.99	0.10	42,43,55,57	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.

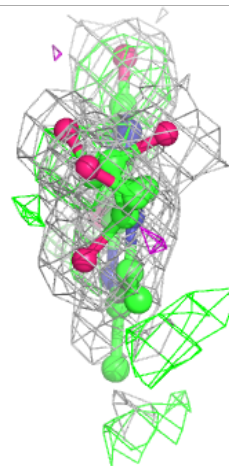
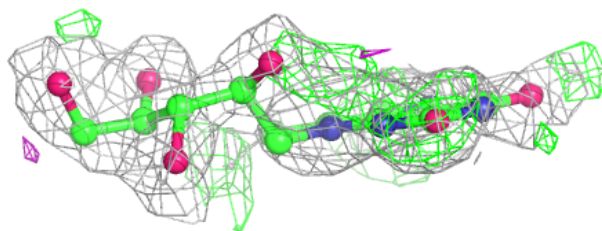
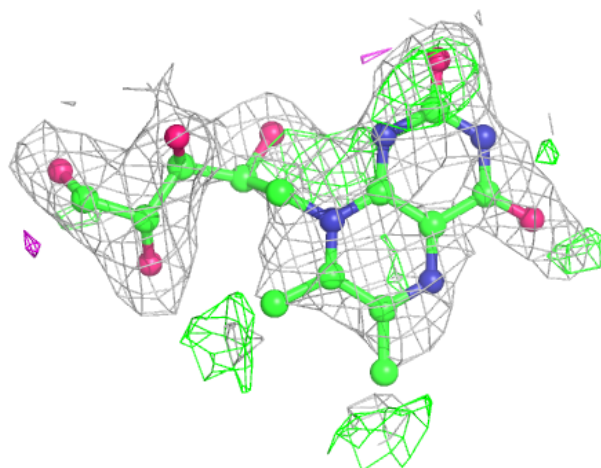
Electron density around DLZ I 204:

$2mF_o-DF_c$ (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.