Full wwPDB X-ray Structure Validation Report (i)

Jun 17, 2024 – 10:29 AM EDT

PDB ID	:	2Y3R
Title	:	Structure of the tirandamycin-bound FAD-dependent tirandamycin oxidase TamL in P21 space group
Authors	:	Carlson, J.C.; Li, S.; Gunatilleke, S.S.; Anzai, Y.; Burr, D.A.; Podust, L.M.; Sherman, D.H.
Deposited on	:	2010-12-22
Resolution	:	1.79 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 (i) 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 (i)) were used in the production of this report:

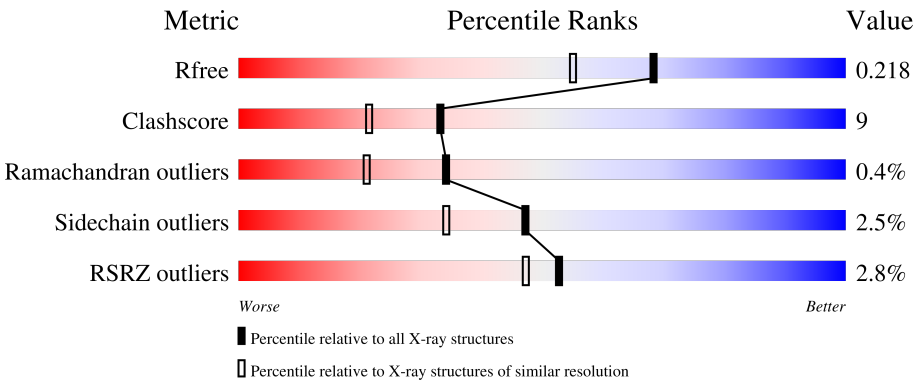
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.79 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 <=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	530	3% 78% 13% • 7%
1	B	530	2% 78% 12% • 8%
1	C	530	2% 81% 10% •• 7%
1	D	530	3% 78% 13% • 7%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	GOL	B	1501	-	X	X	-
5	GOL	C	1502	-	-	X	-

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 17755 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 TAML.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	492	3936	2473	716	739	8	0	20	0
			Total	C	N	O	S			
1	B	487	3768	2381	678	700	9	0	8	0
			Total	C	N	O	S			
1	C	492	3852	2424	704	716	8	0	10	0
			Total	C	N	O	S			
1	D	491	3819	2402	704	705	8	0	7	0

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-29	MET	-	expression tag	UNP D3Y1I2
A	-28	GLY	-	expression tag	UNP D3Y1I2
A	-27	SER	-	expression tag	UNP D3Y1I2
A	-26	HIS	-	expression tag	UNP D3Y1I2
A	-25	HIS	-	expression tag	UNP D3Y1I2
A	-24	HIS	-	expression tag	UNP D3Y1I2
A	-23	HIS	-	expression tag	UNP D3Y1I2
A	-22	HIS	-	expression tag	UNP D3Y1I2
A	-21	HIS	-	expression tag	UNP D3Y1I2
A	-20	HIS	-	expression tag	UNP D3Y1I2
A	-19	HIS	-	expression tag	UNP D3Y1I2
A	-18	GLY	-	expression tag	UNP D3Y1I2
A	-17	SER	-	expression tag	UNP D3Y1I2
A	-16	ASP	-	expression tag	UNP D3Y1I2
A	-15	TYR	-	expression tag	UNP D3Y1I2
A	-14	ASP	-	expression tag	UNP D3Y1I2
A	-13	ILE	-	expression tag	UNP D3Y1I2
A	-12	PRO	-	expression tag	UNP D3Y1I2
A	-11	THR	-	expression tag	UNP D3Y1I2
A	-10	THR	-	expression tag	UNP D3Y1I2
A	-9	GLU	-	expression tag	UNP D3Y1I2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	ASN	-	expression tag	UNP D3Y1I2
A	-7	LEU	-	expression tag	UNP D3Y1I2
A	-6	TYR	-	expression tag	UNP D3Y1I2
A	-5	PHE	-	expression tag	UNP D3Y1I2
A	-4	GLN	-	expression tag	UNP D3Y1I2
A	-3	GLY	-	expression tag	UNP D3Y1I2
A	-2	SER	-	expression tag	UNP D3Y1I2
A	-1	GLU	-	expression tag	UNP D3Y1I2
A	0	PHE	-	expression tag	UNP D3Y1I2
B	-29	MET	-	expression tag	UNP D3Y1I2
B	-28	GLY	-	expression tag	UNP D3Y1I2
B	-27	SER	-	expression tag	UNP D3Y1I2
B	-26	HIS	-	expression tag	UNP D3Y1I2
B	-25	HIS	-	expression tag	UNP D3Y1I2
B	-24	HIS	-	expression tag	UNP D3Y1I2
B	-23	HIS	-	expression tag	UNP D3Y1I2
B	-22	HIS	-	expression tag	UNP D3Y1I2
B	-21	HIS	-	expression tag	UNP D3Y1I2
B	-20	HIS	-	expression tag	UNP D3Y1I2
B	-19	HIS	-	expression tag	UNP D3Y1I2
B	-18	GLY	-	expression tag	UNP D3Y1I2
B	-17	SER	-	expression tag	UNP D3Y1I2
B	-16	ASP	-	expression tag	UNP D3Y1I2
B	-15	TYR	-	expression tag	UNP D3Y1I2
B	-14	ASP	-	expression tag	UNP D3Y1I2
B	-13	ILE	-	expression tag	UNP D3Y1I2
B	-12	PRO	-	expression tag	UNP D3Y1I2
B	-11	THR	-	expression tag	UNP D3Y1I2
B	-10	THR	-	expression tag	UNP D3Y1I2
B	-9	GLU	-	expression tag	UNP D3Y1I2
B	-8	ASN	-	expression tag	UNP D3Y1I2
B	-7	LEU	-	expression tag	UNP D3Y1I2
B	-6	TYR	-	expression tag	UNP D3Y1I2
B	-5	PHE	-	expression tag	UNP D3Y1I2
B	-4	GLN	-	expression tag	UNP D3Y1I2
B	-3	GLY	-	expression tag	UNP D3Y1I2
B	-2	SER	-	expression tag	UNP D3Y1I2
B	-1	GLU	-	expression tag	UNP D3Y1I2
B	0	PHE	-	expression tag	UNP D3Y1I2
C	-29	MET	-	expression tag	UNP D3Y1I2
C	-28	GLY	-	expression tag	UNP D3Y1I2
C	-27	SER	-	expression tag	UNP D3Y1I2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-26	HIS	-	expression tag	UNP D3Y1I2
C	-25	HIS	-	expression tag	UNP D3Y1I2
C	-24	HIS	-	expression tag	UNP D3Y1I2
C	-23	HIS	-	expression tag	UNP D3Y1I2
C	-22	HIS	-	expression tag	UNP D3Y1I2
C	-21	HIS	-	expression tag	UNP D3Y1I2
C	-20	HIS	-	expression tag	UNP D3Y1I2
C	-19	HIS	-	expression tag	UNP D3Y1I2
C	-18	GLY	-	expression tag	UNP D3Y1I2
C	-17	SER	-	expression tag	UNP D3Y1I2
C	-16	ASP	-	expression tag	UNP D3Y1I2
C	-15	TYR	-	expression tag	UNP D3Y1I2
C	-14	ASP	-	expression tag	UNP D3Y1I2
C	-13	ILE	-	expression tag	UNP D3Y1I2
C	-12	PRO	-	expression tag	UNP D3Y1I2
C	-11	THR	-	expression tag	UNP D3Y1I2
C	-10	THR	-	expression tag	UNP D3Y1I2
C	-9	GLU	-	expression tag	UNP D3Y1I2
C	-8	ASN	-	expression tag	UNP D3Y1I2
C	-7	LEU	-	expression tag	UNP D3Y1I2
C	-6	TYR	-	expression tag	UNP D3Y1I2
C	-5	PHE	-	expression tag	UNP D3Y1I2
C	-4	GLN	-	expression tag	UNP D3Y1I2
C	-3	GLY	-	expression tag	UNP D3Y1I2
C	-2	SER	-	expression tag	UNP D3Y1I2
C	-1	GLU	-	expression tag	UNP D3Y1I2
C	0	PHE	-	expression tag	UNP D3Y1I2
D	-29	MET	-	expression tag	UNP D3Y1I2
D	-28	GLY	-	expression tag	UNP D3Y1I2
D	-27	SER	-	expression tag	UNP D3Y1I2
D	-26	HIS	-	expression tag	UNP D3Y1I2
D	-25	HIS	-	expression tag	UNP D3Y1I2
D	-24	HIS	-	expression tag	UNP D3Y1I2
D	-23	HIS	-	expression tag	UNP D3Y1I2
D	-22	HIS	-	expression tag	UNP D3Y1I2
D	-21	HIS	-	expression tag	UNP D3Y1I2
D	-20	HIS	-	expression tag	UNP D3Y1I2
D	-19	HIS	-	expression tag	UNP D3Y1I2
D	-18	GLY	-	expression tag	UNP D3Y1I2
D	-17	SER	-	expression tag	UNP D3Y1I2
D	-16	ASP	-	expression tag	UNP D3Y1I2
D	-15	TYR	-	expression tag	UNP D3Y1I2

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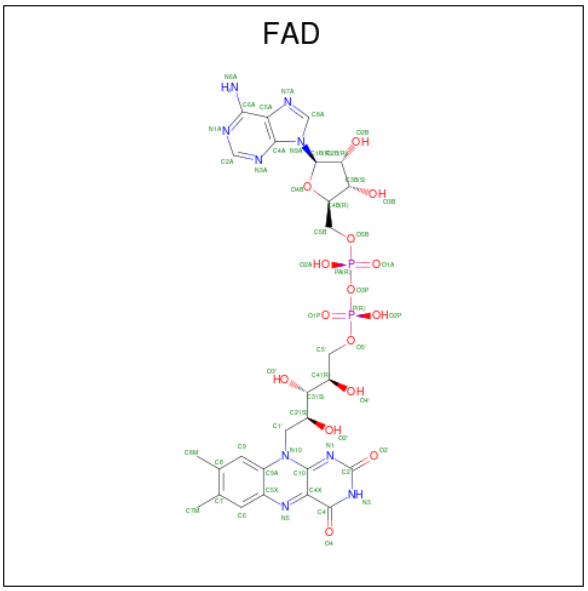
Chain	Residue	Modelled	Actual	Comment	Reference
D	-14	ASP	-	expression tag	UNP D3Y1I2
D	-13	ILE	-	expression tag	UNP D3Y1I2
D	-12	PRO	-	expression tag	UNP D3Y1I2
D	-11	THR	-	expression tag	UNP D3Y1I2
D	-10	THR	-	expression tag	UNP D3Y1I2
D	-9	GLU	-	expression tag	UNP D3Y1I2
D	-8	ASN	-	expression tag	UNP D3Y1I2
D	-7	LEU	-	expression tag	UNP D3Y1I2
D	-6	TYR	-	expression tag	UNP D3Y1I2
D	-5	PHE	-	expression tag	UNP D3Y1I2
D	-4	GLN	-	expression tag	UNP D3Y1I2
D	-3	GLY	-	expression tag	UNP D3Y1I2
D	-2	SER	-	expression tag	UNP D3Y1I2
D	-1	GLU	-	expression tag	UNP D3Y1I2
D	0	PHE	-	expression tag	UNP D3Y1I2

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	53	27	9	15	2	0	0
			Total	C	N	O	P		
2	B	1	53	27	9	15	2	0	0
			Total	C	N	O	P		
2	C	1	53	27	9	15	2	0	0

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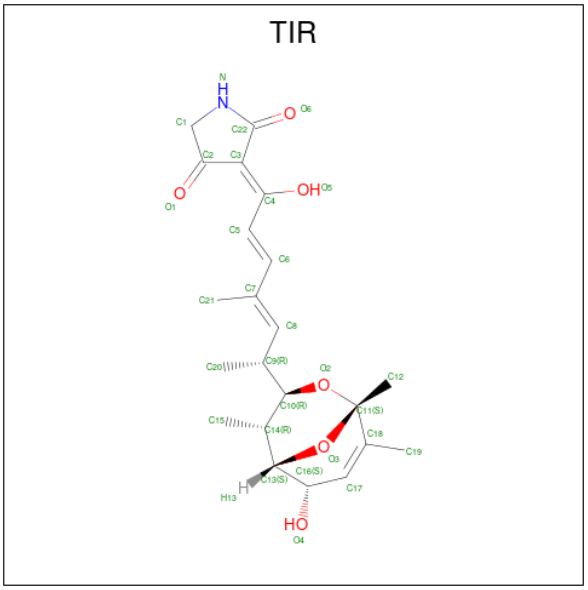
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	D	1	53	27	9	15	2	0	0

- Molecule 3 is TIRANDAMYCIN E (three-letter code: TIR) (formula: $C_{22}H_{29}NO_6$).

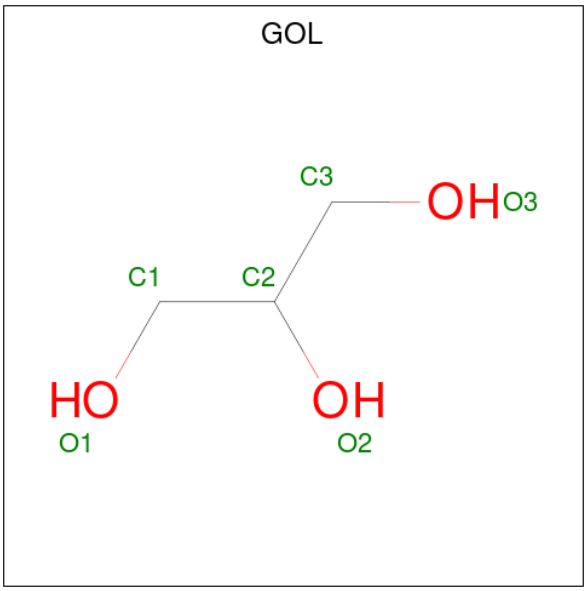


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	29	22	1	6	0	0
3	C	1	29	22	1	6	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
4	A	1	1	1	0	0
4	B	1	1	1	0	0
4	C	1	1	1	0	0
4	D	1	1	1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

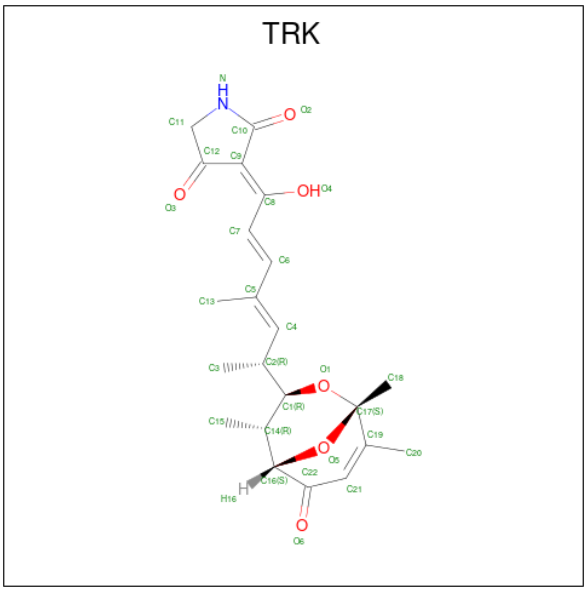


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	6	3	3	0	0
5	B	1	6	3	3	0	0
5	B	1	6	3	3	0	0
5	C	1	6	3	3	0	0
5	C	1	6	3	3	0	0
5	D	1	6	3	3	0	0
5	D	1	6	3	3	0	0

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
6	A	1	1	1	0	0
6	B	1	1	1	0	0
6	C	2	2	2	0	0
6	D	2	2	2	0	0

- Molecule 7 is TIRANDAMYCIN D (three-letter code: TRK) (formula: $C_{22}H_{27}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
7	B	1	29	22	1	6	0	0
			Total	C	N	O		
7	D	1	29	22	1	6	0	0

- Molecule 8 is water.

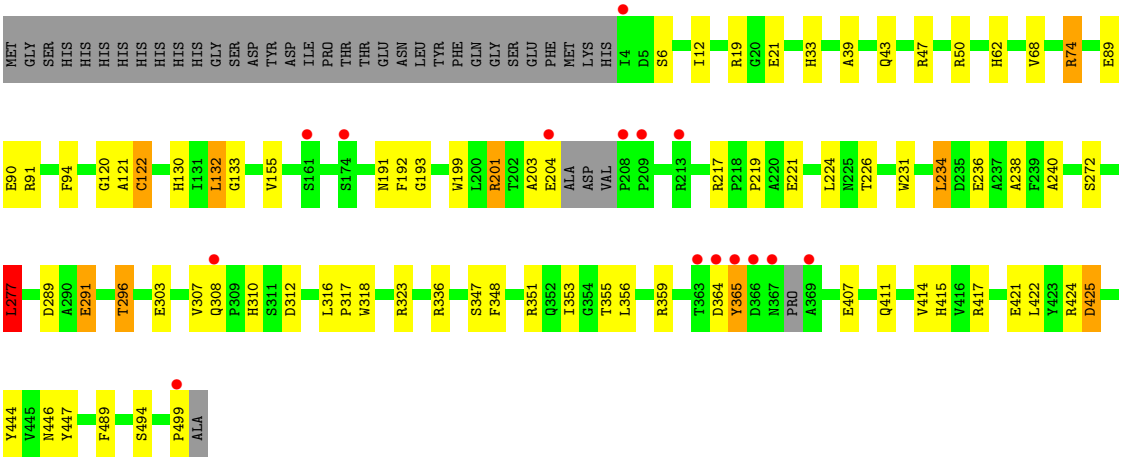
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	A	499	499	499	0	0
			Total	O		
8	B	495	495	495	0	0
			Total	O		
8	C	535	535	535	0	0
			Total	O		
8	D	471	471	471	0	0

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.

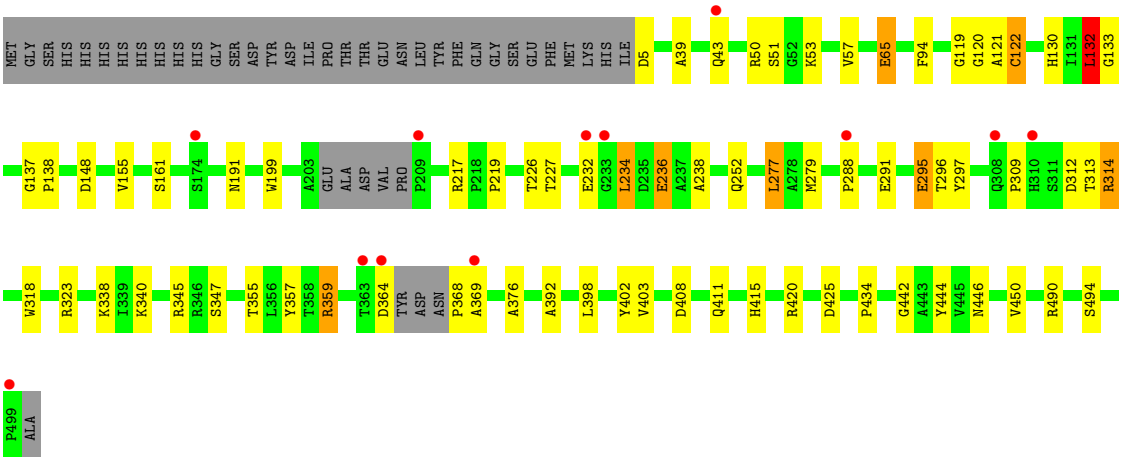
• Molecule 1: TAML

Chain A:



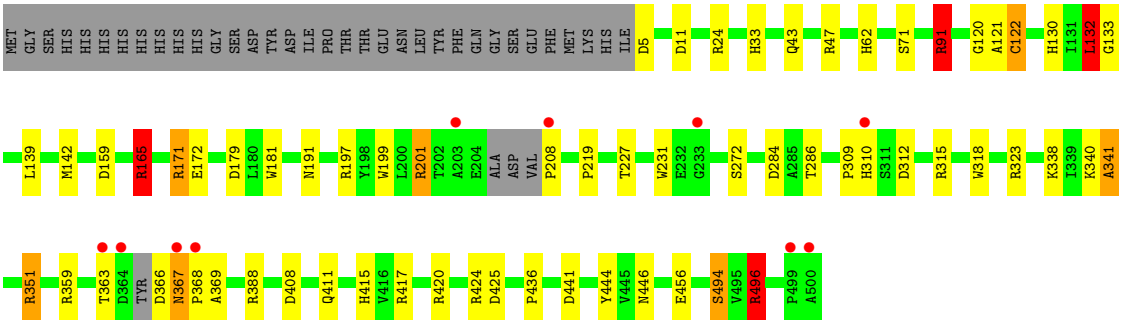
• Molecule 1: TAML

Chain B:



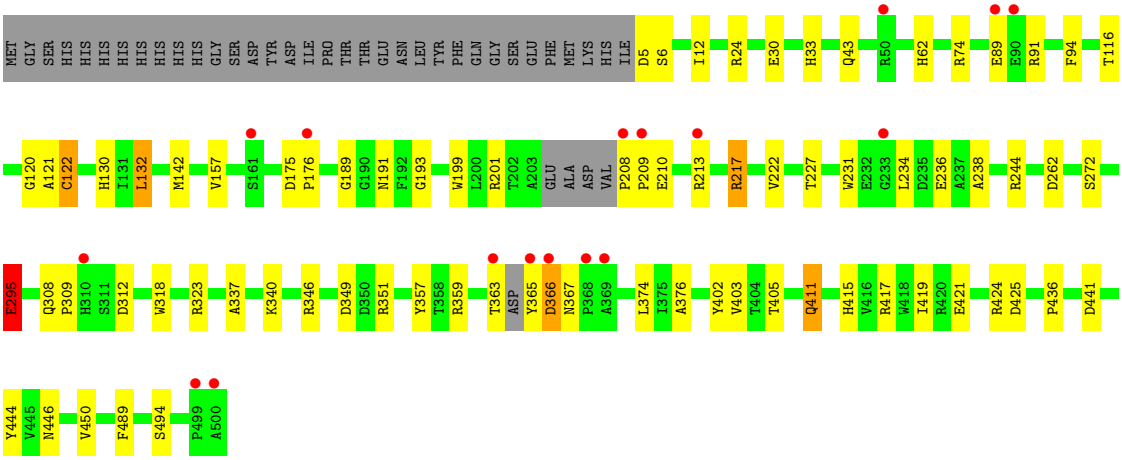
• Molecule 1: TAML

Chain C:



- Molecule 1: TAML

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	71.09Å 106.17Å 151.17Å 90.00° 94.63° 90.00°	Depositor
Resolution (Å)	150.68 – 1.79 19.82 – 1.79	Depositor EDS
% Data completeness (in resolution range)	76.5 (150.68-1.79) 76.5 (19.82-1.79)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.11 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.146 , 0.219 0.146 , 0.218	Depositor DCC
R_{free} test set	8138 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	16.6	Xtriage
Anisotropy	0.436	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 52.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	17755	wwPDB-VP
Average B, all atoms (Å^2)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.20 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3666e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CL, MG, TRK, GOL, FAD, TIR

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	1.14	6/4037 (0.1%)	0.99	12/5512 (0.2%)
1	B	1.15	7/3867 (0.2%)	0.99	12/5284 (0.2%)
1	C	1.14	5/3952 (0.1%)	1.00	19/5397 (0.4%)
1	D	1.13	6/3918 (0.2%)	0.97	6/5348 (0.1%)
All	All	1.14	24/15774 (0.2%)	0.99	49/21541 (0.2%)

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	201	ARG	CB-CG	-6.92	1.33	1.52
1	D	295	GLU	CB-CG	-6.48	1.39	1.52
1	A	21	GLU	CG-CD	6.40	1.61	1.51
1	B	347	SER	CB-OG	6.26	1.50	1.42
1	B	402	TYR	CD2-CE2	6.15	1.48	1.39
1	C	494	SER	CB-OG	6.14	1.50	1.42
1	A	348	PHE	CE2-CZ	6.11	1.49	1.37
1	B	65	GLU	CB-CG	5.96	1.63	1.52
1	C	341	ALA	CA-CB	5.91	1.64	1.52
1	B	392	ALA	CA-CB	5.86	1.64	1.52
1	B	295	GLU	CB-CG	-5.85	1.41	1.52
1	B	347	SER	CA-CB	5.77	1.61	1.52
1	C	179	ASP	CB-CG	5.72	1.63	1.51
1	D	157	VAL	CB-CG2	5.52	1.64	1.52
1	B	402	TYR	CD1-CE1	5.38	1.47	1.39
1	A	201	ARG	CB-CG	-5.26	1.38	1.52
1	A	407	GLU	CB-CG	-5.24	1.42	1.52
1	D	402	TYR	CD1-CE1	5.23	1.47	1.39
1	A	68	VAL	CB-CG2	5.16	1.63	1.52
1	A	192	PHE	CE2-CZ	5.13	1.47	1.37
1	D	89	GLU	CG-CD	5.12	1.59	1.51
1	D	222	VAL	CB-CG2	5.08	1.63	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	351	ARG	CZ-NH1	5.07	1.39	1.33
1	D	376	ALA	CA-CB	5.01	1.62	1.52

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	351	ARG	NE-CZ-NH2	-8.70	115.95	120.30
1	A	19	ARG	NE-CZ-NH2	-8.63	115.99	120.30
1	C	312	ASP	CB-CG-OD1	7.95	125.46	118.30
1	B	323	ARG	NE-CZ-NH2	-7.76	116.42	120.30
1	A	74	ARG	NE-CZ-NH1	7.69	124.14	120.30
1	B	277	LEU	CA-CB-CG	7.63	132.85	115.30
1	A	277	LEU	CA-CB-CG	7.09	131.60	115.30
1	C	315	ARG	NE-CZ-NH2	-6.93	116.84	120.30
1	A	336	ARG	NE-CZ-NH1	6.85	123.73	120.30
1	B	132[A]	LEU	CB-CG-CD2	-6.46	100.02	111.00
1	B	132[B]	LEU	CB-CG-CD2	-6.46	100.02	111.00
1	B	347	SER	N-CA-CB	6.43	120.15	110.50
1	D	208	PRO	N-CA-CB	6.36	110.94	103.30
1	B	277	LEU	CB-CG-CD2	-6.22	100.43	111.00
1	C	351	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	D	24	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	D	24	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	C	408	ASP	CB-CG-OD1	5.80	123.52	118.30
1	A	336	ARG	NE-CZ-NH2	-5.74	117.43	120.30
1	C	165[A]	ARG	NE-CZ-NH1	-5.73	117.44	120.30
1	C	165[B]	ARG	NE-CZ-NH1	-5.73	117.44	120.30
1	B	403	VAL	CB-CA-C	-5.72	100.53	111.40
1	C	179	ASP	CB-CG-OD1	5.72	123.45	118.30
1	A	201	ARG	CG-CD-NE	-5.70	99.82	111.80
1	B	345	ARG	NE-CZ-NH1	5.61	123.11	120.30
1	C	165[A]	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	C	165[B]	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	C	323	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	C	417	ARG	NE-CZ-NH1	5.53	123.06	120.30
1	B	408	ASP	CB-CG-OD2	-5.52	113.33	118.30
1	C	91	ARG	CG-CD-NE	-5.52	100.21	111.80
1	A	289	ASP	CB-CG-OD1	5.44	123.20	118.30
1	B	345	ARG	NE-CZ-NH2	-5.39	117.61	120.30
1	C	197	ARG	NE-CZ-NH1	5.37	122.98	120.30
1	D	312	ASP	CB-CG-OD1	5.34	123.11	118.30
1	D	323	ARG	NE-CZ-NH2	-5.33	117.64	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	11	ASP	CB-CG-OD1	5.27	123.05	118.30
1	A	91	ARG	CG-CD-NE	-5.23	100.81	111.80
1	B	408	ASP	CB-CG-OD1	5.23	123.00	118.30
1	C	132[A]	LEU	CB-CG-CD2	-5.21	102.14	111.00
1	C	132[B]	LEU	CB-CG-CD2	-5.21	102.14	111.00
1	C	91	ARG	NE-CZ-NH2	-5.19	117.71	120.30
1	A	291	GLU	OE1-CD-OE2	-5.18	117.09	123.30
1	C	496	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	A	91	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	A	323	ARG	NE-CZ-NH2	-5.09	117.76	120.30
1	A	19	ARG	NE-CZ-NH1	5.05	122.83	120.30
1	D	91	ARG	NE-CZ-NH2	-5.02	117.79	120.30
1	B	148	ASP	CB-CG-OD2	5.01	122.81	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3936	0	3761	65	0
1	B	3768	0	3615	56	0
1	C	3852	0	3705	65	0
1	D	3819	0	3671	60	0
2	A	53	0	30	9	0
2	B	53	0	29	1	0
2	C	53	0	29	9	0
2	D	53	0	29	6	0
3	A	29	0	28	5	0
3	C	29	0	28	5	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	6	0	8	1	0
5	B	12	0	15	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	12	0	15	11	0
5	D	12	0	15	2	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	2	0	0	0	0
6	D	2	0	0	0	0
7	B	29	0	26	3	0
7	D	29	0	26	2	0
8	A	499	0	0	17	0
8	B	495	0	0	13	1
8	C	535	0	0	17	0
8	D	471	0	0	16	0
All	All	17755	0	15030	261	1

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 9.

All (261) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:HIS:ND1	2:A:600:FAD:HM83	0.95	1.27
1:A:62:HIS:CE1	2:A:600:FAD:HM83	1.71	1.24
1:D:121:ALA:O	1:D:122:CYS:CB	1.86	1.20
1:C:121:ALA:O	1:C:122:CYS:CB	1.82	1.19
1:A:121:ALA:O	1:A:122:CYS:CB	1.87	1.19
1:B:121:ALA:O	1:B:122:CYS:CB	1.89	1.17
1:C:286:THR:HG22	8:C:2259:HOH:O	1.45	1.15
1:C:338:LYS:HE2	5:C:1502:GOL:H31	1.24	1.09
2:C:600:FAD:N5	3:C:700:TIR:H16	1.68	1.09
1:D:346[B]:ARG:HG3	1:D:346[B]:ARG:HH11	1.08	1.08
1:C:338:LYS:CE	5:C:1502:GOL:H31	1.82	1.08
1:C:24:ARG:HH22	5:C:1502:GOL:H12	1.13	1.06
1:C:208:PRO:HD3	8:C:2250:HOH:O	1.56	1.04
1:D:5:ASP:N	8:D:2002:HOH:O	1.89	1.04
1:D:210:GLU:CB	1:D:213:ARG:HH21	1.70	1.03
1:A:499:PRO:C	8:A:2492:HOH:O	1.95	1.03
1:C:71:SER:HB2	8:C:2043:HOH:O	1.60	1.01
2:A:600:FAD:N5	3:A:700:TIR:H16	1.76	0.99
1:B:368:PRO:CD	8:B:2358:HOH:O	2.09	0.99
1:C:165[A]:ARG:HG3	1:C:165[A]:ARG:HH11	1.24	0.98
1:A:121:ALA:O	1:A:122:CYS:HB2	1.16	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:121:ALA:O	1:D:122:CYS:HB2	1.12	0.96
1:B:121:ALA:O	1:B:122:CYS:HB2	1.14	0.95
1:C:165[A]:ARG:HH11	1:C:165[A]:ARG:CG	1.76	0.95
1:A:365:TYR:OH	1:A:421[A]:GLU:OE1	1.87	0.93
1:C:24:ARG:HH22	5:C:1502:GOL:C1	1.82	0.91
1:C:165[A]:ARG:HG3	1:C:165[A]:ARG:NH1	1.82	0.90
1:A:414[A]:VAL:HG23	8:A:2403:HOH:O	1.71	0.90
1:D:213:ARG:HG2	8:D:2205:HOH:O	1.74	0.88
1:C:121:ALA:O	1:C:122:CYS:HB2	1.04	0.88
1:C:24:ARG:NH2	5:C:1502:GOL:H12	1.89	0.87
1:A:74:ARG:HD3	8:A:2110:HOH:O	1.73	0.87
1:C:284:ASP:OD1	1:C:286:THR:HG23	1.75	0.87
1:D:346[B]:ARG:HG3	1:D:346[B]:ARG:NH1	1.82	0.84
1:A:89[B]:GLU:OE1	1:A:89[B]:GLU:N	2.09	0.84
1:A:89[B]:GLU:CD	1:A:89[B]:GLU:H	1.76	0.83
1:C:62:HIS:ND1	2:C:600:FAD:HM81	1.94	0.82
1:D:359:ARG:NH1	1:D:425:ASP:OD2	2.13	0.81
1:D:421[B]:GLU:OE1	1:D:424:ARG:NH2	2.13	0.80
1:C:62:HIS:CE1	2:C:600:FAD:C8M	2.64	0.80
1:D:244[B]:ARG:NH2	1:D:349:ASP:HA	1.96	0.80
1:C:338:LYS:HE2	5:C:1502:GOL:C3	2.09	0.79
1:C:159:ASP:OD1	1:C:165[A]:ARG:NH1	2.16	0.79
1:B:359:ARG:NH1	1:B:425:ASP:OD2	2.17	0.78
1:C:359:ARG:NH1	1:C:425:ASP:OD2	2.17	0.77
1:D:244[B]:ARG:HH21	1:D:349:ASP:HA	1.50	0.77
1:A:132[A]:LEU:HD23	1:A:132[A]:LEU:C	2.04	0.77
1:A:421[B]:GLU:OE1	1:A:424:ARG:NH2	2.17	0.77
1:D:62:HIS:CE1	2:D:600:FAD:C8M	2.67	0.76
1:C:366:ASP:O	1:C:368:PRO:HD3	1.86	0.75
1:B:279[B]:MET:CE	1:B:297:TYR:OH	2.34	0.74
1:B:39:ALA:O	1:B:43[A]:GLN:HG2	1.87	0.74
2:C:600:FAD:C4X	3:C:700:TIR:H16	2.17	0.73
2:A:600:FAD:C4X	3:A:700:TIR:H16	2.18	0.73
1:D:346[B]:ARG:NH1	1:D:346[B]:ARG:CG	2.50	0.72
1:A:62:HIS:ND1	2:A:600:FAD:HM81	1.97	0.72
1:C:369:ALA:O	8:C:2396:HOH:O	2.07	0.71
1:D:365:TYR:CB	8:D:2364:HOH:O	2.38	0.70
1:A:33:HIS:HE1	8:A:2043:HOH:O	1.73	0.70
1:C:5:ASP:CB	8:D:2131:HOH:O	2.40	0.69
1:B:279[B]:MET:HE2	1:B:297:TYR:OH	1.91	0.69
1:D:417[B]:ARG:HG2	8:D:2372:HOH:O	1.90	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:ALA:O	1:A:43[A]:GLN:HG2	1.91	0.68
1:B:368:PRO:CG	8:B:2358:HOH:O	2.40	0.68
1:B:217:ARG:NH2	8:B:2230:HOH:O	2.18	0.68
1:B:65:GLU:OE1	5:B:1501:GOL:H31	1.94	0.68
1:D:308:GLN:O	8:D:2274:HOH:O	2.11	0.68
1:D:62:HIS:ND1	2:D:600:FAD:HM81	1.98	0.67
1:B:227:THR:HG23	1:B:309:PRO:HB2	1.75	0.67
1:A:355:THR:HG21	1:A:425[A]:ASP:HB3	1.77	0.67
1:B:132[A]:LEU:C	1:B:132[A]:LEU:HD23	2.12	0.67
1:B:338:LYS:NZ	5:B:1501:GOL:H2	2.10	0.66
1:A:236:GLU:OE2	8:A:2259:HOH:O	2.12	0.66
1:A:234:LEU:HD23	1:A:307:VAL:HG21	1.77	0.65
1:A:310[B]:HIS:HD2	8:A:2250:HOH:O	1.80	0.64
1:B:369:ALA:HA	8:B:2164:HOH:O	1.98	0.64
1:D:374:LEU:HD12	1:D:374:LEU:N	2.13	0.63
1:A:132[A]:LEU:HD23	1:A:133:GLY:N	2.14	0.63
1:B:313:THR:C	1:B:314[B]:ARG:HG2	2.17	0.63
1:C:91:ARG:HD2	1:C:199:TRP:CE2	2.34	0.63
1:D:411:GLN:O	1:D:415:HIS:HD2	1.80	0.62
1:C:62:HIS:ND1	2:C:600:FAD:C8	2.59	0.62
1:B:279[B]:MET:HE1	1:B:297:TYR:OH	1.99	0.62
1:B:291:GLU:O	1:B:295:GLU:HG2	1.99	0.62
1:A:296:THR:HG21	8:A:2046:HOH:O	2.00	0.62
1:B:252:GLN:CG	8:B:2254:HOH:O	2.48	0.61
1:A:351[A]:ARG:NH1	8:A:2362:HOH:O	2.20	0.61
1:A:62:HIS:ND1	2:A:600:FAD:C8	2.59	0.61
1:C:310:HIS:HD2	8:C:2262:HOH:O	1.83	0.61
1:C:33:HIS:HE1	8:C:2054:HOH:O	1.83	0.61
1:C:227:THR:HG23	1:C:309:PRO:HB2	1.81	0.60
1:D:142:MET:HE2	1:D:262:ASP:HA	1.83	0.60
1:C:132[A]:LEU:HD22	1:C:132[A]:LEU:H	1.67	0.59
1:D:62:HIS:CG	2:D:600:FAD:C8M	2.76	0.59
1:D:130:HIS:HD2	2:D:600:FAD:O2'	1.85	0.59
1:A:130:HIS:HD2	2:A:600:FAD:O2'	1.85	0.59
1:D:244[B]:ARG:NH2	8:D:2225:HOH:O	2.11	0.59
5:C:1502:GOL:O1	8:C:2535:HOH:O	2.16	0.59
1:B:132[A]:LEU:N	1:B:132[A]:LEU:CD2	2.63	0.58
1:C:132[A]:LEU:CD2	1:C:132[A]:LEU:N	2.66	0.57
1:C:456:GLU:OE1	8:C:2469:HOH:O	2.17	0.57
1:C:91:ARG:NE	8:C:2150:HOH:O	2.26	0.57
1:C:366:ASP:O	1:C:368:PRO:CD	2.52	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:236:GLU:HG3	1:D:357:TYR:CG	2.38	0.57
1:D:210:GLU:CB	1:D:213:ARG:NH2	2.55	0.57
1:C:132[A]:LEU:C	1:C:132[A]:LEU:HD23	2.26	0.57
1:A:414[A]:VAL:CG2	8:A:2403:HOH:O	2.43	0.56
7:B:700:TRK:C7	5:B:1500:GOL:H31	2.34	0.56
1:C:338:LYS:HD3	5:C:1502:GOL:H11	1.87	0.56
1:A:411[A]:GLN:O	1:A:415:HIS:HD2	1.89	0.56
1:B:236:GLU:HG2	1:B:357:TYR:CG	2.41	0.55
1:A:234:LEU:HD22	1:A:238:ALA:HB3	1.89	0.55
7:D:700:TRK:H14	7:D:700:TRK:H132	1.89	0.55
1:C:341:ALA:HB3	8:C:2456:HOH:O	2.07	0.54
1:A:411[B]:GLN:O	1:A:415:HIS:HD2	1.91	0.54
1:A:308:GLN:O	8:A:2318:HOH:O	2.18	0.54
1:D:295:GLU:HG2	8:D:2261:HOH:O	2.08	0.54
1:B:368:PRO:HD3	8:B:2358:HOH:O	1.89	0.54
1:A:120:GLY:H	1:A:130:HIS:CE1	2.26	0.53
1:B:132[A]:LEU:HD23	1:B:132[A]:LEU:N	2.15	0.53
1:D:120:GLY:H	1:D:130:HIS:CE1	2.26	0.53
3:A:700:TIR:H121	3:A:700:TIR:H315	1.91	0.53
1:B:132[A]:LEU:CD2	1:B:132[A]:LEU:H	2.21	0.53
1:C:444:TYR:CZ	1:C:446:ASN:HB2	2.43	0.53
1:A:204:GLU:C	8:A:2237:HOH:O	2.46	0.53
1:A:444:TYR:CZ	1:A:446:ASN:HB2	2.44	0.53
1:B:234:LEU:HD22	1:B:238:ALA:HB3	1.91	0.53
1:B:368:PRO:HD2	8:B:2358:HOH:O	1.94	0.53
7:B:700:TRK:H132	7:B:700:TRK:H14	1.90	0.53
1:D:217:ARG:NH2	8:D:2209:HOH:O	2.21	0.53
1:D:244[B]:ARG:HH21	1:D:349:ASP:CA	2.22	0.53
3:C:700:TIR:H121	3:C:700:TIR:H315	1.91	0.53
1:B:338:LYS:HZ3	5:B:1501:GOL:H2	1.74	0.53
1:C:132[A]:LEU:H	1:C:132[A]:LEU:CD2	2.22	0.53
1:A:193:GLY:HA2	1:A:489:PHE:CE2	2.44	0.53
1:A:231:TRP:CE2	1:A:277:LEU:HD13	2.44	0.52
7:B:700:TRK:C6	5:B:1500:GOL:H31	2.40	0.52
1:D:444:TYR:CZ	1:D:446:ASN:HB2	2.44	0.52
1:B:450:VAL:CG2	5:B:1501:GOL:H32	2.39	0.52
1:D:227:THR:HG23	1:D:309:PRO:HB2	1.92	0.52
1:C:120:GLY:H	1:C:130:HIS:CE1	2.27	0.52
1:B:130:HIS:HD2	2:B:600:FAD:O2'	1.93	0.52
1:A:221:GLU:HG2	1:A:317:PRO:HA	1.91	0.52
1:A:47:ARG:HG2	1:A:50[B]:ARG:NH2	2.24	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:411:GLN:O	1:B:415:HIS:HD2	1.93	0.51
1:C:338:LYS:CD	5:C:1502:GOL:H31	2.38	0.51
1:B:191:ASN:HD21	1:B:494:SER:HA	1.75	0.51
1:B:444:TYR:CZ	1:B:446:ASN:HB2	2.45	0.51
1:B:288:PRO:HD2	8:B:2283:HOH:O	2.10	0.51
1:A:217:ARG:NH2	8:A:2243:HOH:O	2.35	0.51
1:C:191:ASN:HD21	1:C:494:SER:HA	1.75	0.51
1:A:231:TRP:CD1	1:A:272:SER:HA	2.46	0.51
1:A:296:THR:CG2	8:A:2046:HOH:O	2.59	0.51
1:B:132[A]:LEU:H	1:B:132[A]:LEU:HD22	1.75	0.51
1:B:132[A]:LEU:HD23	1:B:133:GLY:N	2.25	0.51
1:D:337:ALA:HB2	1:D:405:THR:HG22	1.91	0.51
1:C:231:TRP:CD1	1:C:272:SER:HA	2.47	0.50
1:A:226[A]:THR:HG23	1:A:226[A]:THR:O	2.11	0.50
1:B:5:ASP:HA	8:B:2004:HOH:O	2.10	0.50
1:B:119:GLY:O	1:B:137:GLY:HA3	2.12	0.49
1:A:365:TYR:HE1	1:A:417:ARG:HG2	1.76	0.49
1:C:132[A]:LEU:N	1:C:132[A]:LEU:HD23	2.27	0.49
1:D:62:HIS:ND1	2:D:600:FAD:C8	2.66	0.49
1:D:231:TRP:CD1	1:D:272:SER:HA	2.47	0.49
1:A:355:THR:HG21	1:A:425[B]:ASP:CG	2.33	0.49
1:D:417[B]:ARG:HG3	8:D:2367:HOH:O	2.12	0.49
1:A:191:ASN:HD21	1:A:494:SER:HA	1.78	0.49
1:B:120:GLY:H	1:B:130:HIS:CE1	2.30	0.48
1:A:356:LEU:HD23	1:A:422:LEU:HD11	1.95	0.48
1:C:132[A]:LEU:HD23	1:C:133:GLY:N	2.28	0.48
1:A:62:HIS:CG	2:A:600:FAD:C8M	2.82	0.48
1:B:57[B]:VAL:HG12	8:B:2486:HOH:O	2.13	0.48
1:C:436:PRO:HA	1:C:441:ASP:HA	1.96	0.48
1:D:33:HIS:HD2	8:D:2001:HOH:O	1.96	0.48
1:C:388:ARG:O	1:C:388:ARG:HG2	2.14	0.47
1:C:411:GLN:O	1:C:415:HIS:HD2	1.97	0.47
1:D:236:GLU:O	1:D:236:GLU:HG2	2.15	0.47
1:A:291:GLU:HG2	8:A:2144:HOH:O	2.14	0.47
1:B:234:LEU:HD22	1:B:238:ALA:CB	2.44	0.47
1:A:359:ARG:NH2	1:A:425[A]:ASP:OD2	2.40	0.47
1:B:338:LYS:HZ2	5:B:1501:GOL:H2	1.76	0.47
1:C:130:HIS:HD2	2:C:600:FAD:O2'	1.98	0.47
1:A:90:GLU:OE2	1:C:47:ARG:NH1	2.48	0.47
1:A:234:LEU:HD22	1:A:238:ALA:CB	2.45	0.47
1:D:193:GLY:HA2	1:D:489:PHE:CE2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:450:VAL:HG22	5:B:1501:GOL:H32	1.95	0.47
1:A:310[A]:HIS:CE1	8:A:2319:HOH:O	2.68	0.46
1:C:91:ARG:NH2	8:C:2150:HOH:O	2.46	0.46
1:B:313:THR:O	1:B:314[B]:ARG:HG2	2.15	0.46
1:D:234:LEU:HD12	1:D:238:ALA:HB3	1.96	0.46
1:A:121:ALA:O	1:A:122:CYS:SG	2.70	0.46
1:A:231:TRP:NE1	1:A:277:LEU:HD13	2.31	0.46
1:D:351[A]:ARG:NH2	8:D:2324:HOH:O	2.27	0.46
1:A:203:ALA:O	1:A:204:GLU:C	2.54	0.46
1:A:224:LEU:CD1	1:A:316:LEU:HD12	2.46	0.45
1:D:94:PHE:O	1:D:199:TRP:HA	2.16	0.45
1:A:347:SER:HB2	8:A:2360:HOH:O	2.15	0.45
1:D:30:GLU:OE2	1:D:74:ARG:HD2	2.16	0.45
7:D:700:TRK:H182	5:D:1501:GOL:H11	1.97	0.45
1:A:219:PRO:O	1:A:318:TRP:HB2	2.17	0.44
1:A:89[B]:GLU:OE1	1:C:43:GLN:OE1	2.35	0.44
1:B:450:VAL:HG22	5:B:1501:GOL:C3	2.47	0.44
1:A:364:ASP:O	1:A:365:TYR:HB2	2.18	0.44
3:A:700:TIR:C21	3:A:700:TIR:H14	2.48	0.44
1:D:351[B]:ARG:NH1	8:D:2324:HOH:O	2.50	0.44
1:D:121:ALA:O	1:D:122:CYS:SG	2.74	0.44
1:A:447:TYR:HH	3:A:700:TIR:HO4	1.57	0.43
1:B:368:PRO:HG3	8:B:2358:HOH:O	2.13	0.43
1:C:62:HIS:CG	2:C:600:FAD:C8M	2.84	0.43
8:C:2173:HOH:O	1:D:5:ASP:CB	2.65	0.43
5:C:1501:GOL:H31	8:C:2357:HOH:O	2.19	0.43
1:A:417:ARG:NH1	1:A:421[A]:GLU:OE2	2.48	0.43
1:A:94:PHE:O	1:A:199:TRP:HA	2.18	0.43
1:D:116:THR:HG1	1:D:318:TRP:HH2	1.65	0.43
1:C:424:ARG:NH1	8:C:2439:HOH:O	2.04	0.43
1:B:51:SER:HB3	1:B:53:LYS:HG3	2.00	0.43
1:D:436:PRO:HA	1:D:441:ASP:HA	2.00	0.43
1:C:171[B]:ARG:O	1:C:171[B]:ARG:HG2	2.18	0.43
1:D:417[A]:ARG:HD3	8:D:2373:HOH:O	2.19	0.43
1:A:226[A]:THR:HG22	1:A:312:ASP:HB3	2.01	0.43
1:C:338:LYS:NZ	5:C:1502:GOL:H31	2.31	0.42
2:C:600:FAD:C4X	3:C:700:TIR:C16	2.93	0.42
1:A:414[B]:VAL:HG13	8:A:2403:HOH:O	2.18	0.42
1:D:236:GLU:HG3	1:D:357:TYR:CD2	2.54	0.42
1:D:191:ASN:HD21	1:D:494:SER:HA	1.85	0.42
1:C:340:LYS:HE3	1:C:420:ARG:HA	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:351[A]:ARG:NE	8:D:2324:HOH:O	2.43	0.42
1:D:363:THR:HA	1:D:366:ASP:OD1	2.19	0.42
2:A:600:FAD:HM72	5:A:1500:GOL:O3	2.19	0.42
1:C:219:PRO:O	1:C:318:TRP:HB2	2.19	0.42
2:D:600:FAD:H9	2:D:600:FAD:H1'2	1.83	0.42
1:D:450:VAL:HG22	5:D:1502:GOL:H2	2.02	0.42
1:D:340:LYS:HD2	1:D:419:ILE:HG12	2.03	0.41
1:C:496:ARG:HD2	8:C:2522:HOH:O	2.20	0.41
1:A:240:ALA:HA	1:A:353:ILE:HG21	2.03	0.41
1:C:496:ARG:HD2	8:C:2523:HOH:O	2.20	0.41
1:D:351[A]:ARG:NH1	8:D:2321:HOH:O	2.53	0.41
1:C:351:ARG:NE	8:C:2384:HOH:O	2.53	0.41
2:C:600:FAD:C5X	3:C:700:TIR:H16	2.47	0.41
1:D:175:ASP:OD1	1:D:176:PRO:HD2	2.21	0.41
1:B:121:ALA:O	1:B:122:CYS:SG	2.77	0.41
1:B:364:ASP:C	8:B:2355:HOH:O	2.58	0.41
1:C:121:ALA:O	1:C:122:CYS:SG	2.77	0.41
1:B:490:ARG:HD2	8:B:2467:HOH:O	2.21	0.41
1:D:12:ILE:HD13	1:D:12:ILE:HG21	1.79	0.41
1:D:142:MET:HE2	1:D:142:MET:HB2	1.89	0.41
1:C:139:LEU:HD22	1:C:142:MET:HE3	2.03	0.41
1:D:132:LEU:HD13	1:D:189:GLY:HA2	2.02	0.41
1:A:12:ILE:HG21	1:A:12:ILE:HD13	1.72	0.40
1:B:161:SER:HB2	1:D:43[B]:GLN:HE22	1.86	0.40
1:B:434:PRO:O	1:B:442:GLY:HA2	2.20	0.40
1:B:376:ALA:HA	1:B:398:LEU:HD23	2.02	0.40
1:C:367:ASN:HD22	1:C:367:ASN:H	1.69	0.40
1:C:171[A]:ARG:HB3	1:C:181:TRP:CE2	2.56	0.40
1:B:219:PRO:O	1:B:318:TRP:HB2	2.22	0.40
1:B:94:PHE:O	1:B:199:TRP:HA	2.21	0.40
1:B:226[A]:THR:HG23	1:B:226[A]:THR:O	2.22	0.40
1:B:340:LYS:HE3	1:B:420:ARG:HA	2.03	0.40
1:B:355:THR:O	1:B:359:ARG:HG3	2.22	0.40

All (1) 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	Interatomic distance (Å)	Clash overlap (Å)
8:B:2285:HOH:O	8:B:2431:HOH:O[2_556]	2.13	0.07

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	506/530 (96%)	491 (97%)	13 (3%)	2 (0%)	34	21
1	B	489/530 (92%)	477 (98%)	11 (2%)	1 (0%)	47	33
1	C	496/530 (94%)	485 (98%)	9 (2%)	2 (0%)	34	21
1	D	492/530 (93%)	481 (98%)	8 (2%)	3 (1%)	25	12
All	All	1983/2120 (94%)	1934 (98%)	41 (2%)	8 (0%)	34	21

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	122	CYS
1	B	122	CYS
1	C	122	CYS
1	C	367	ASN
1	D	122	CYS
1	A	365	TYR
1	D	366	ASP
1	D	209	PRO

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	400/421 (95%)	388 (97%)	12 (3%)	41	27
1	B	378/421 (90%)	363 (96%)	15 (4%)	31	16

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	389/421 (92%)	377 (97%)	12 (3%)	40	25
1	D	382/421 (91%)	374 (98%)	8 (2%)	53	42
All	All	1549/1684 (92%)	1502 (97%)	47 (3%)	47	27

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	SER
1	A	132[A]	LEU
1	A	132[B]	LEU
1	A	155[A]	VAL
1	A	155[B]	VAL
1	A	201	ARG
1	A	234	LEU
1	A	277	LEU
1	A	296	THR
1	A	303	GLU
1	A	425[A]	ASP
1	A	425[B]	ASP
1	B	50	ARG
1	B	132[A]	LEU
1	B	132[B]	LEU
1	B	138	PRO
1	B	155[A]	VAL
1	B	155[B]	VAL
1	B	232	GLU
1	B	234	LEU
1	B	236	GLU
1	B	277	LEU
1	B	296	THR
1	B	312	ASP
1	B	314[A]	ARG
1	B	314[B]	ARG
1	B	359	ARG
1	C	91	ARG
1	C	132[A]	LEU
1	C	132[B]	LEU
1	C	165[A]	ARG
1	C	165[B]	ARG
1	C	171[A]	ARG
1	C	171[B]	ARG

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Mol	Chain	Res	Type
1	C	172[A]	GLU
1	C	172[B]	GLU
1	C	201	ARG
1	C	363	THR
1	C	496	ARG
1	D	6	SER
1	D	132	LEU
1	D	201	ARG
1	D	217	ARG
1	D	295	GLU
1	D	367	ASN
1	D	403	VAL
1	D	411	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	HIS
1	A	130	HIS
1	A	177	ASN
1	A	191	ASN
1	A	225	ASN
1	A	415	HIS
1	B	130	HIS
1	B	177	ASN
1	B	191	ASN
1	B	225	ASN
1	B	415	HIS
1	C	33	HIS
1	C	43	GLN
1	C	130	HIS
1	C	177	ASN
1	C	191	ASN
1	C	225	ASN
1	C	252	GLN
1	C	310	HIS
1	C	415	HIS
1	D	130	HIS
1	D	151	HIS
1	D	177	ASN
1	D	191	ASN
1	D	225	ASN

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Mol	Chain	Res	Type
1	D	415	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 25 ligands modelled in this entry, 10 are monoatomic - leaving 15 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	TIR	A	700	4	31,31,31	1.49	5 (16%)	30,47,47	3.82	9 (30%)
2	FAD	D	600	1	54,58,58	1.18	6 (11%)	71,89,89	1.70	15 (21%)
5	GOL	A	1500	-	5,5,5	0.92	0	5,5,5	1.48	1 (20%)
5	GOL	D	1501	-	5,5,5	0.95	0	5,5,5	1.08	0
5	GOL	C	1502	-	5,5,5	1.07	0	5,5,5	1.98	2 (40%)
5	GOL	B	1500	-	5,5,5	0.78	0	5,5,5	1.00	0
3	TIR	C	700	4	31,31,31	1.63	9 (29%)	30,47,47	3.58	7 (23%)
5	GOL	C	1501	-	5,5,5	0.71	0	5,5,5	0.36	0
5	GOL	D	1502	-	5,5,5	1.52	1 (20%)	5,5,5	2.02	2 (40%)
2	FAD	A	600	1	54,58,58	1.60	10 (18%)	71,89,89	1.83	21 (29%)
5	GOL	B	1501	-	5,5,5	1.31	1 (20%)	5,5,5	1.44	1 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
7	TRK	D	700	4	31,31,31	1.66	7 (22%)	27,47,47	3.64	10 (37%)
2	FAD	C	600	1	54,58,58	1.21	4 (7%)	71,89,89	1.78	16 (22%)
7	TRK	B	700	4	31,31,31	1.42	4 (12%)	27,47,47	4.12	10 (37%)
2	FAD	B	600	1	54,58,58	1.30	7 (12%)	71,89,89	1.84	14 (19%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	TIR	A	700	4	-	1/17/62/62	0/3/3/3
2	FAD	D	600	1	-	6/30/50/50	0/6/6/6
5	GOL	A	1500	-	-	2/4/4/4	-
5	GOL	D	1501	-	-	4/4/4/4	-
5	GOL	C	1502	-	-	2/4/4/4	-
5	GOL	B	1500	-	-	2/4/4/4	-
3	TIR	C	700	4	-	2/17/62/62	0/3/3/3
5	GOL	C	1501	-	-	2/4/4/4	-
5	GOL	D	1502	-	-	2/4/4/4	-
2	FAD	A	600	1	-	6/30/50/50	0/6/6/6
5	GOL	B	1501	-	-	4/4/4/4	-
7	TRK	D	700	4	-	2/17/62/62	0/2/3/3
2	FAD	C	600	1	-	5/30/50/50	0/6/6/6
7	TRK	B	700	4	-	2/17/62/62	0/2/3/3
2	FAD	B	600	1	-	6/30/50/50	0/6/6/6

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	FAD	P-O3P	5.16	1.65	1.59
2	A	600	FAD	O4-C4	4.40	1.31	1.23
7	D	700	TRK	O3-C12	4.01	1.31	1.23
2	A	600	FAD	C4X-N5	3.81	1.39	1.30
7	D	700	TRK	C21-C19	3.69	1.38	1.33
7	B	700	TRK	O3-C12	3.61	1.30	1.23
2	C	600	FAD	C2A-N3A	3.52	1.37	1.32
3	A	700	TIR	C17-C18	3.52	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	700	TIR	C1-C2	-3.52	1.48	1.52
2	C	600	FAD	C9A-N10	-3.50	1.35	1.41
2	B	600	FAD	C4X-N5	3.41	1.38	1.30
3	C	700	TIR	C1-N	3.41	1.50	1.44
2	B	600	FAD	O3'-C3'	3.38	1.51	1.43
2	D	600	FAD	C4-N3	-3.35	1.32	1.38
2	A	600	FAD	O4B-C1B	3.29	1.45	1.40
3	A	700	TIR	C16-C17	3.19	1.54	1.50
2	C	600	FAD	C4X-N5	3.15	1.37	1.30
7	B	700	TRK	C9-C8	3.07	1.49	1.40
2	A	600	FAD	C9A-N10	-3.07	1.35	1.41
5	D	1502	GOL	O2-C2	-3.05	1.34	1.43
3	A	700	TIR	O1-C2	2.99	1.29	1.23
2	D	600	FAD	C9A-N10	-2.95	1.36	1.41
2	D	600	FAD	C4X-N5	2.89	1.37	1.30
3	C	700	TIR	O6-C22	2.87	1.29	1.23
3	C	700	TIR	C16-C17	2.75	1.53	1.50
3	C	700	TIR	O1-C2	2.72	1.28	1.23
3	A	700	TIR	O6-C22	2.59	1.28	1.23
2	A	600	FAD	C9-C9A	2.56	1.43	1.39
7	D	700	TRK	C7-C8	2.55	1.48	1.45
2	B	600	FAD	C4-N3	-2.52	1.34	1.38
5	B	1501	GOL	O2-C2	-2.49	1.36	1.43
7	D	700	TRK	C13-C5	2.48	1.55	1.50
3	C	700	TIR	C3-C4	2.48	1.47	1.40
2	B	600	FAD	C9A-N10	-2.41	1.37	1.41
7	D	700	TRK	C9-C8	2.40	1.47	1.40
2	B	600	FAD	C5'-C4'	2.39	1.55	1.51
2	A	600	FAD	C6-C7	-2.38	1.36	1.39
3	C	700	TIR	C17-C18	2.31	1.36	1.33
3	C	700	TIR	C13-C16	2.29	1.55	1.51
2	D	600	FAD	O4-C4	2.28	1.27	1.23
2	A	600	FAD	C8M-C8	2.20	1.55	1.51
7	D	700	TRK	C21-C22	2.20	1.50	1.46
2	D	600	FAD	C5'-C4'	2.19	1.54	1.51
7	D	700	TRK	O1-C1	2.17	1.47	1.44
3	A	700	TIR	C1-N	2.17	1.48	1.44
2	A	600	FAD	C2A-N1A	2.12	1.37	1.33
2	B	600	FAD	O4-C4	2.07	1.27	1.23
7	B	700	TRK	O5-C16	2.07	1.46	1.43
3	C	700	TIR	O2-C11	2.07	1.45	1.42
2	C	600	FAD	C2-N3	-2.07	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	FAD	O3'-C3'	2.06	1.48	1.43
2	D	600	FAD	C9-C8	2.04	1.42	1.39
7	B	700	TRK	O2-C10	2.01	1.27	1.23
2	B	600	FAD	C1'-C2'	-2.01	1.49	1.52

All (108) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	B	700	TRK	C11-C12-C9	17.70	115.85	106.86
3	A	700	TIR	C1-C2-C3	16.65	115.32	106.86
3	C	700	TIR	C1-C2-C3	15.46	114.72	106.86
7	D	700	TRK	C11-C12-C9	13.94	113.94	106.86
7	B	700	TRK	C9-C10-N	7.48	115.96	107.25
3	A	700	TIR	C3-C22-N	7.34	115.80	107.25
7	D	700	TRK	C9-C10-N	7.14	115.56	107.25
2	D	600	FAD	N3A-C2A-N1A	-6.81	119.42	128.67
3	C	700	TIR	C3-C22-N	6.37	114.66	107.25
7	D	700	TRK	O4-C8-C7	5.89	123.74	115.70
2	B	600	FAD	N3A-C2A-N1A	-5.15	121.68	128.67
2	A	600	FAD	N3A-C2A-N1A	-5.06	121.80	128.67
2	C	600	FAD	N3A-C2A-N1A	-4.99	121.90	128.67
2	A	600	FAD	C10-C4X-N5	-4.86	114.88	124.81
2	B	600	FAD	C4'-C3'-C2'	-4.81	105.56	113.57
3	A	700	TIR	O1-C2-C3	-4.68	120.98	128.60
2	B	600	FAD	C4X-C10-N10	4.65	123.14	116.48
7	B	700	TRK	C7-C6-C5	-4.63	119.39	126.23
3	C	700	TIR	C5-C4-C3	-4.61	119.47	124.39
3	A	700	TIR	C5-C6-C7	-4.55	119.51	126.23
3	C	700	TIR	C5-C6-C7	-4.43	119.68	126.23
2	A	600	FAD	C4-C4X-N5	4.39	124.27	118.21
7	D	700	TRK	C7-C6-C5	-4.36	119.78	126.23
2	D	600	FAD	C4'-C3'-C2'	-4.27	106.46	113.57
2	B	600	FAD	O2-C2-N1	-4.26	114.72	121.80
2	C	600	FAD	C4X-C10-N10	4.12	122.38	116.48
3	C	700	TIR	O5-C4-C5	4.02	121.19	115.70
3	A	700	TIR	O5-C4-C5	4.02	121.19	115.70
2	C	600	FAD	C4-N3-C2	-4.01	118.52	125.64
2	C	600	FAD	O2A-PA-O3P	3.98	118.04	107.27
7	D	700	TRK	O3-C12-C9	-3.96	122.16	128.60
2	A	600	FAD	C4X-C10-N10	3.88	122.04	116.48
2	D	600	FAD	C4X-C10-N10	3.80	121.92	116.48
2	B	600	FAD	C10-C4X-N5	-3.65	117.35	124.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	600	FAD	C10-C4X-N5	-3.57	117.52	124.81
2	B	600	FAD	C5'-C4'-C3'	-3.52	105.58	112.22
3	C	700	TIR	O1-C2-C3	-3.48	122.93	128.60
2	C	600	FAD	O2P-P-O3P	3.45	116.61	107.27
5	C	1502	GOL	O3-C3-C2	-3.45	94.85	110.38
2	C	600	FAD	C4-C4X-N5	3.44	122.96	118.21
2	B	600	FAD	C4-C4X-N5	3.43	122.94	118.21
7	D	700	TRK	C7-C8-C9	-3.28	120.88	124.39
3	A	700	TIR	C5-C4-C3	-3.27	120.89	124.39
2	A	600	FAD	O2'-C2'-C3'	3.18	116.68	109.25
2	D	600	FAD	O3P-PA-O1A	-3.10	101.37	110.70
2	D	600	FAD	O2P-P-O3P	3.08	115.61	107.27
2	C	600	FAD	O3P-PA-O1A	-3.08	101.43	110.70
2	B	600	FAD	C4-N3-C2	-3.07	120.19	125.64
7	B	700	TRK	O3-C12-C9	-3.04	123.65	128.60
2	A	600	FAD	C9-C9A-N10	3.02	125.91	121.85
2	A	600	FAD	C4'-C3'-C2'	-2.93	108.69	113.57
7	B	700	TRK	C15-C14-C16	2.92	115.59	111.20
5	D	1502	GOL	O2-C2-C3	-2.91	97.12	109.18
2	A	600	FAD	C5X-N5-C4X	2.87	122.73	118.09
2	D	600	FAD	C10-C4X-N5	-2.87	118.95	124.81
7	B	700	TRK	C7-C8-C9	-2.84	121.35	124.39
3	C	700	TIR	C6-C5-C4	-2.83	117.96	122.49
2	A	600	FAD	O3P-PA-O1A	-2.76	102.39	110.70
5	C	1502	GOL	O1-C1-C2	-2.71	98.17	110.38
2	A	600	FAD	O4B-C1B-N9A	2.69	112.31	108.75
2	A	600	FAD	C7M-C7-C8	2.66	126.20	120.76
7	D	700	TRK	O2-C10-C9	-2.66	119.83	129.43
2	C	600	FAD	C4X-C10-N1	-2.64	118.11	124.59
2	C	600	FAD	C4A-C5A-N7A	-2.64	106.55	109.34
2	A	600	FAD	C4A-C5A-N7A	-2.63	106.56	109.34
2	D	600	FAD	C4-N3-C2	-2.62	120.98	125.64
2	C	600	FAD	C4X-C4-N3	2.61	119.89	113.25
5	A	1500	GOL	O2-C2-C3	-2.60	98.42	109.18
2	B	600	FAD	O4'-C4'-C3'	2.60	115.33	109.25
7	B	700	TRK	O4-C8-C7	2.55	119.18	115.70
7	B	700	TRK	C3-C2-C4	-2.55	105.76	110.08
7	B	700	TRK	O3-C12-C11	-2.52	120.47	123.01
3	A	700	TIR	O4-C16-C17	2.52	114.96	109.27
5	D	1502	GOL	C3-C2-C1	2.52	121.03	111.80
5	B	1501	GOL	O3-C3-C2	2.50	121.64	110.38
2	A	600	FAD	C7M-C7-C6	-2.50	115.17	119.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	600	FAD	O2P-P-O3P	2.48	113.97	107.27
2	C	600	FAD	O2-C2-N1	-2.47	117.69	121.80
2	B	600	FAD	C6-C7-C8	-2.44	116.11	119.69
3	A	700	TIR	C20-C9-C10	-2.43	106.84	111.11
2	C	600	FAD	C1B-N9A-C4A	-2.43	122.37	126.64
2	B	600	FAD	O2A-PA-O3P	2.43	113.83	107.27
3	A	700	TIR	O6-C22-C3	-2.41	120.73	129.43
2	A	600	FAD	C5'-C4'-C3'	-2.40	107.68	112.22
2	D	600	FAD	O3B-C3B-C2B	-2.38	104.19	111.82
2	D	600	FAD	O2A-PA-O3P	2.38	113.70	107.27
2	A	600	FAD	O2-C2-N1	-2.28	118.01	121.80
2	A	600	FAD	C4-C4X-C10	2.26	120.81	116.93
2	B	600	FAD	C1B-N9A-C4A	-2.26	122.67	126.64
2	D	600	FAD	C7M-C7-C8	2.26	125.37	120.76
2	D	600	FAD	C4X-C10-N1	-2.25	119.07	124.59
7	D	700	TRK	O4-C8-C9	-2.24	115.24	119.12
7	B	700	TRK	O2-C10-C9	-2.22	121.42	129.43
2	A	600	FAD	O2'-C2'-C1'	-2.21	101.12	110.20
2	B	600	FAD	O3P-PA-O1A	-2.21	104.06	110.70
2	D	600	FAD	O2'-C2'-C3'	2.16	114.29	109.25
2	A	600	FAD	O2A-PA-O3P	2.15	113.09	107.27
2	C	600	FAD	C4'-C3'-C2'	-2.13	110.02	113.57
2	A	600	FAD	O3P-P-O1P	-2.11	104.36	110.70
2	B	600	FAD	C5X-C6-C7	2.09	124.70	120.83
2	D	600	FAD	C4A-C5A-N7A	-2.07	107.15	109.34
2	C	600	FAD	C6-C7-C8	-2.06	116.67	119.69
2	D	600	FAD	C4-C4X-N5	2.06	121.05	118.21
7	D	700	TRK	O6-C22-C16	-2.06	118.43	121.05
2	C	600	FAD	O2'-C2'-C1'	-2.04	101.83	110.20
7	D	700	TRK	C3-C2-C4	-2.04	106.63	110.08
2	A	600	FAD	C9A-C5X-N5	-2.03	120.31	122.45
2	D	600	FAD	O2A-PA-O1A	2.01	121.78	112.44

There are no chirality outliers.

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1500	GOL	C1-C2-C3-O3
5	B	1500	GOL	C1-C2-C3-O3
5	B	1500	GOL	O2-C2-C3-O3
5	B	1501	GOL	O1-C1-C2-O2
5	B	1501	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
5	C	1502	GOL	O1-C1-C2-C3
5	D	1502	GOL	C1-C2-C3-O3
5	D	1502	GOL	O2-C2-C3-O3
3	C	700	TIR	C5-C6-C7-C21
7	B	700	TRK	C13-C5-C6-C7
5	A	1500	GOL	O2-C2-C3-O3
5	C	1501	GOL	O1-C1-C2-O2
2	D	600	FAD	O3'-C3'-C4'-C5'
2	B	600	FAD	C2'-C3'-C4'-C5'
7	D	700	TRK	C13-C5-C6-C7
3	C	700	TIR	C5-C6-C7-C8
7	B	700	TRK	C4-C5-C6-C7
2	D	600	FAD	O3'-C3'-C4'-O4'
5	B	1501	GOL	C1-C2-C3-O3
5	C	1501	GOL	O1-C1-C2-C3
5	D	1501	GOL	O1-C1-C2-C3
5	D	1501	GOL	C1-C2-C3-O3
5	D	1501	GOL	O1-C1-C2-O2
2	B	600	FAD	O3'-C3'-C4'-C5'
2	C	600	FAD	O3'-C3'-C4'-O4'
5	D	1501	GOL	O2-C2-C3-O3
2	B	600	FAD	O3'-C3'-C4'-O4'
2	A	600	FAD	C2'-C3'-C4'-O4'
2	B	600	FAD	C2'-C3'-C4'-O4'
2	C	600	FAD	C2'-C3'-C4'-O4'
2	A	600	FAD	C2'-C3'-C4'-C5'
2	C	600	FAD	C2'-C3'-C4'-C5'
2	D	600	FAD	C2'-C3'-C4'-C5'
2	D	600	FAD	C2'-C3'-C4'-O4'
2	A	600	FAD	O3'-C3'-C4'-C5'
2	A	600	FAD	O3'-C3'-C4'-O4'
5	C	1502	GOL	O1-C1-C2-O2
5	B	1501	GOL	O2-C2-C3-O3
2	C	600	FAD	O3'-C3'-C4'-C5'
2	A	600	FAD	PA-O3P-P-O2P
2	D	600	FAD	PA-O3P-P-O2P
2	B	600	FAD	PA-O3P-P-O2P
2	C	600	FAD	PA-O3P-P-O2P
2	A	600	FAD	PA-O3P-P-O1P
2	D	600	FAD	PA-O3P-P-O1P
3	A	700	TIR	C7-C8-C9-C20
7	D	700	TRK	C4-C5-C6-C7

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Mol	Chain	Res	Type	Atoms
2	B	600	FAD	PA-O3P-P-O1P

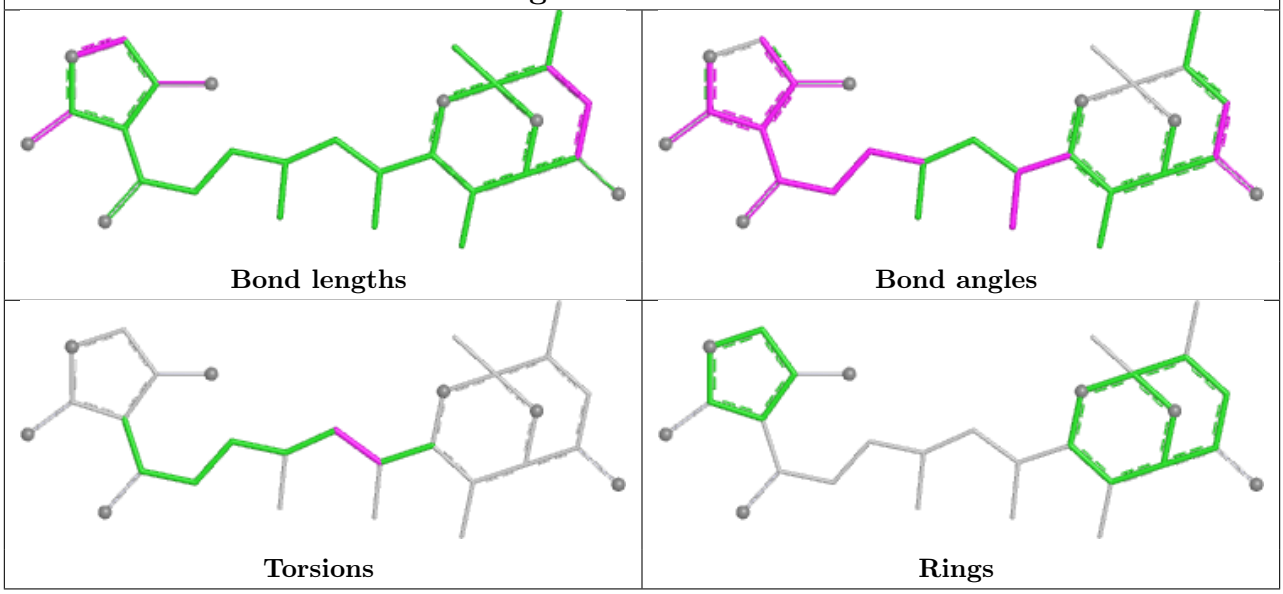
There are no ring outliers.

15 monomers are involved in 53 short contacts:

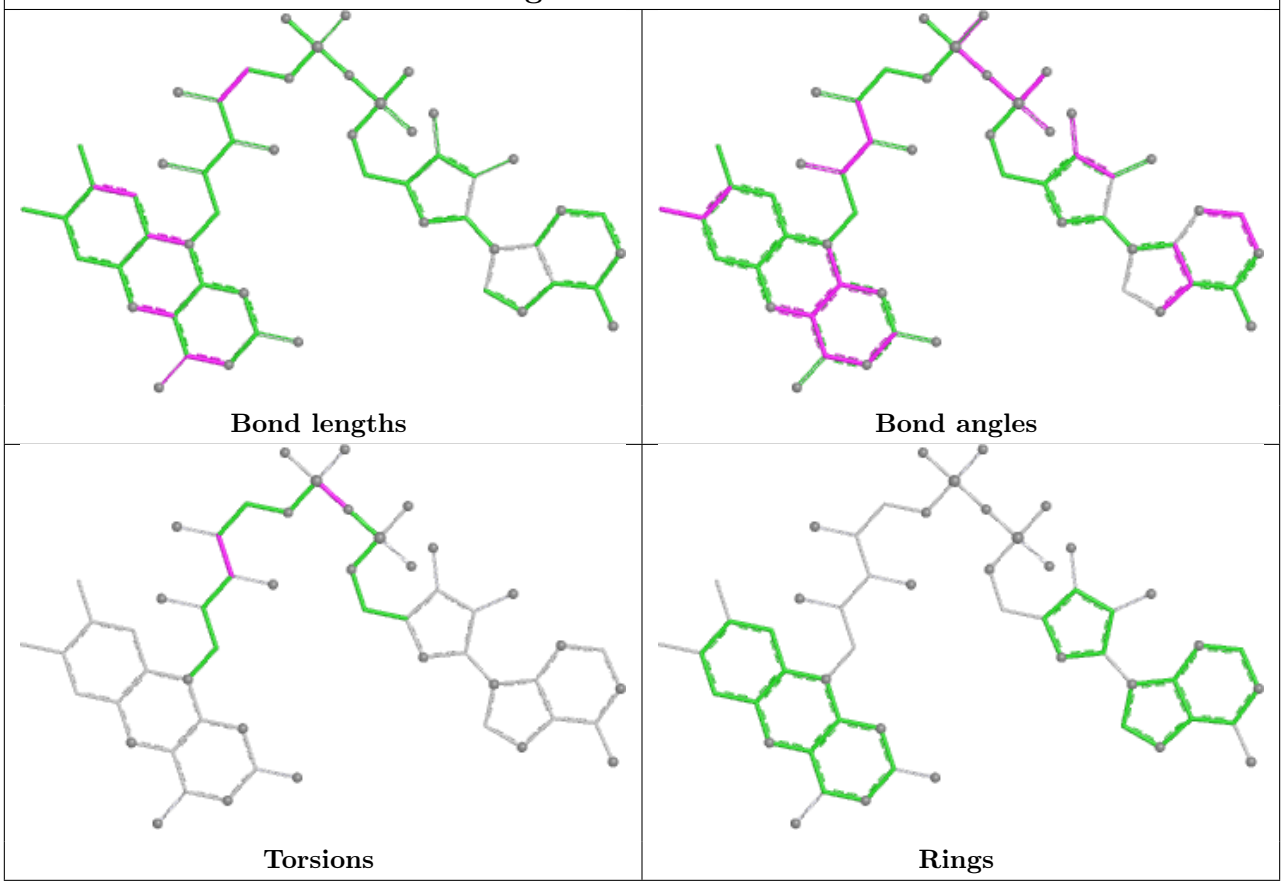
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	700	TIR	5	0
2	D	600	FAD	6	0
5	A	1500	GOL	1	0
5	D	1501	GOL	1	0
5	C	1502	GOL	10	0
5	B	1500	GOL	2	0
3	C	700	TIR	5	0
5	C	1501	GOL	1	0
5	D	1502	GOL	1	0
2	A	600	FAD	9	0
5	B	1501	GOL	7	0
7	D	700	TRK	2	0
2	C	600	FAD	9	0
7	B	700	TRK	3	0
2	B	600	FAD	1	0

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 then 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.

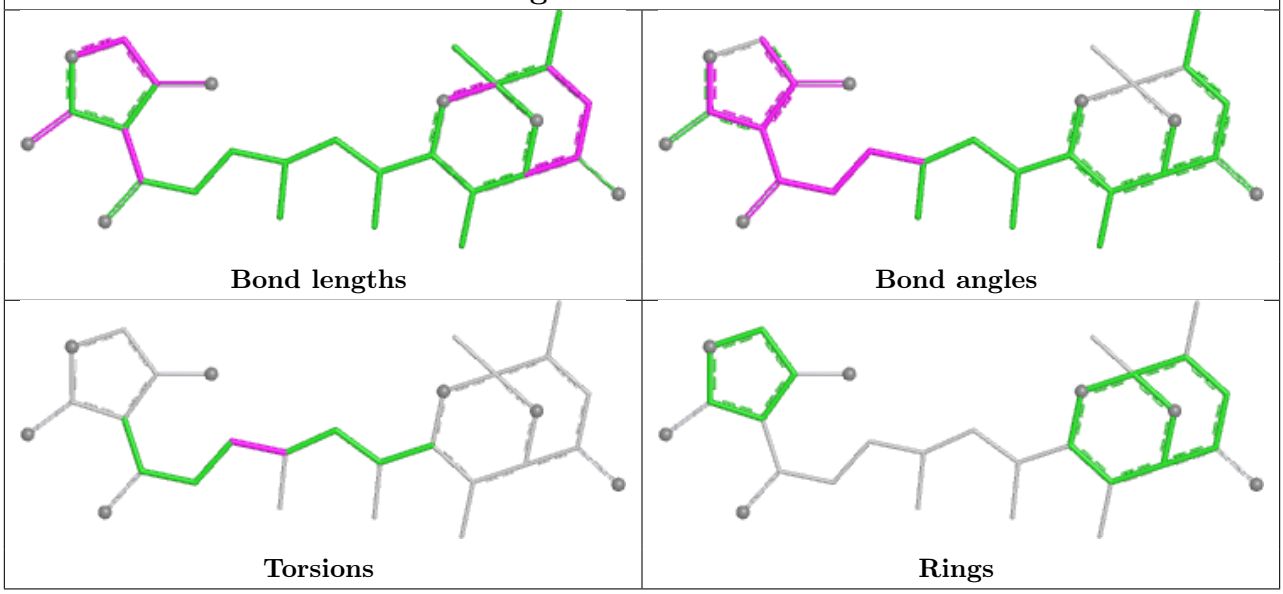
Ligand TIR A 700



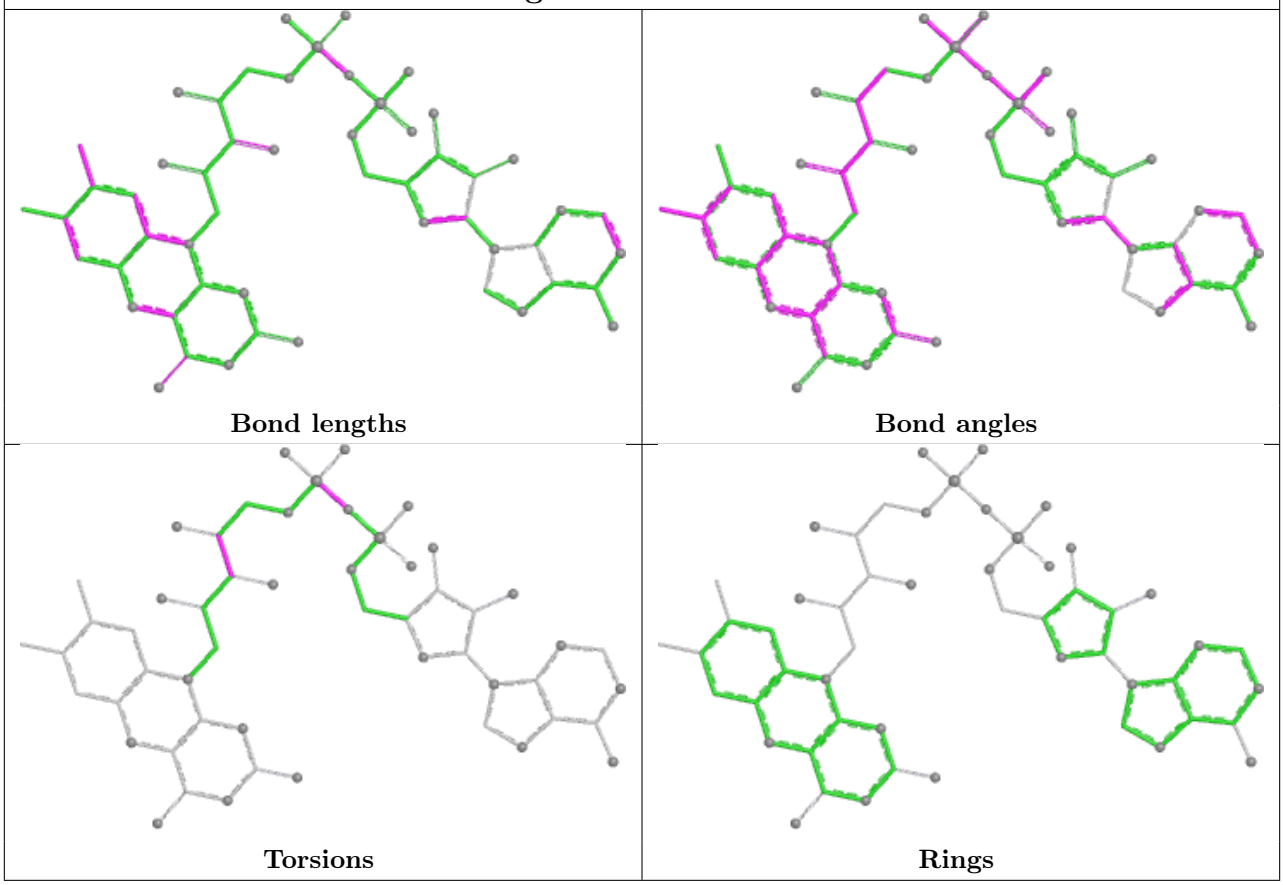
Ligand FAD D 600



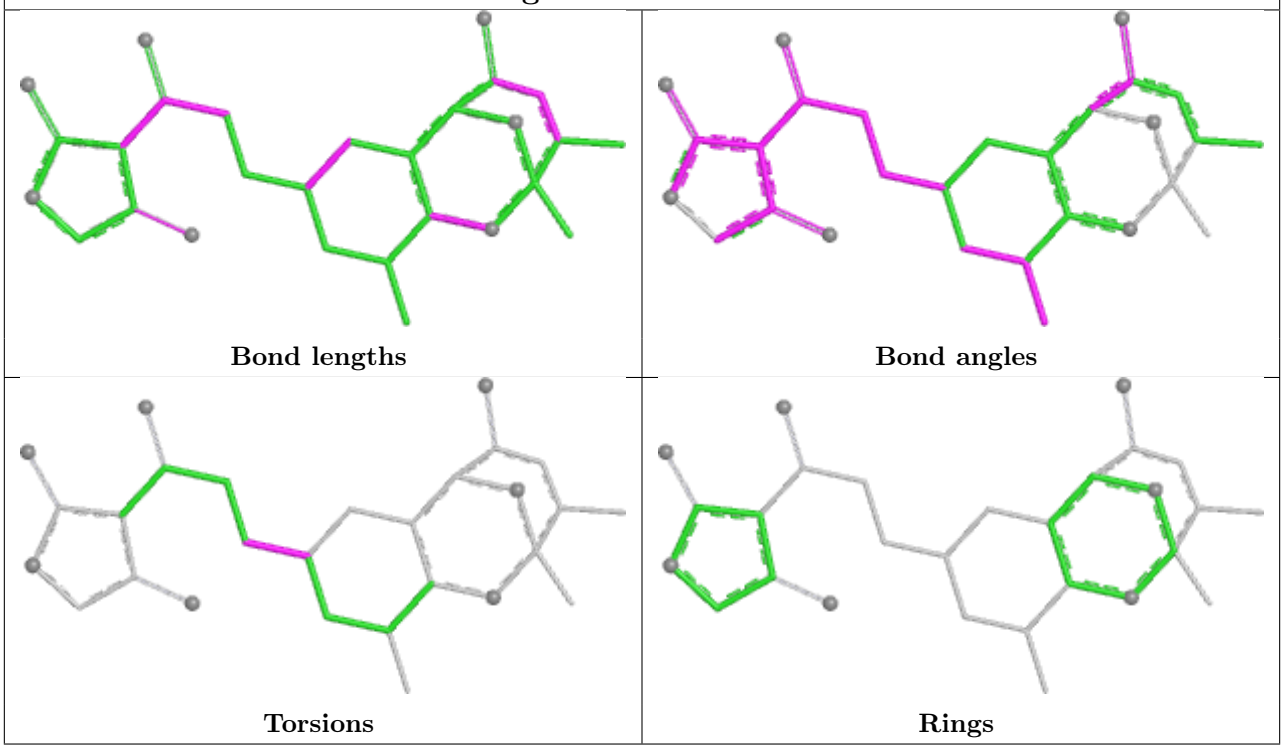
Ligand TIR C 700



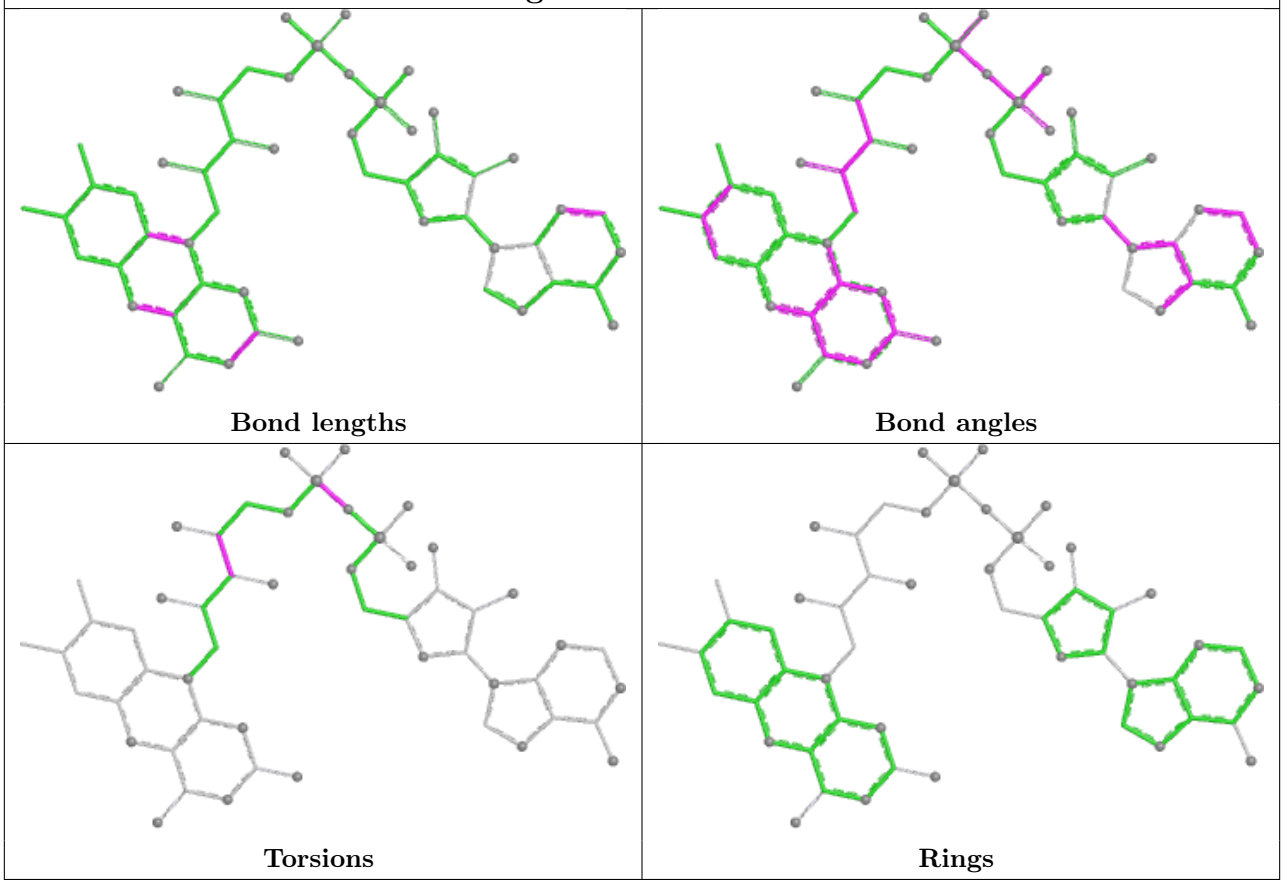
Ligand FAD A 600



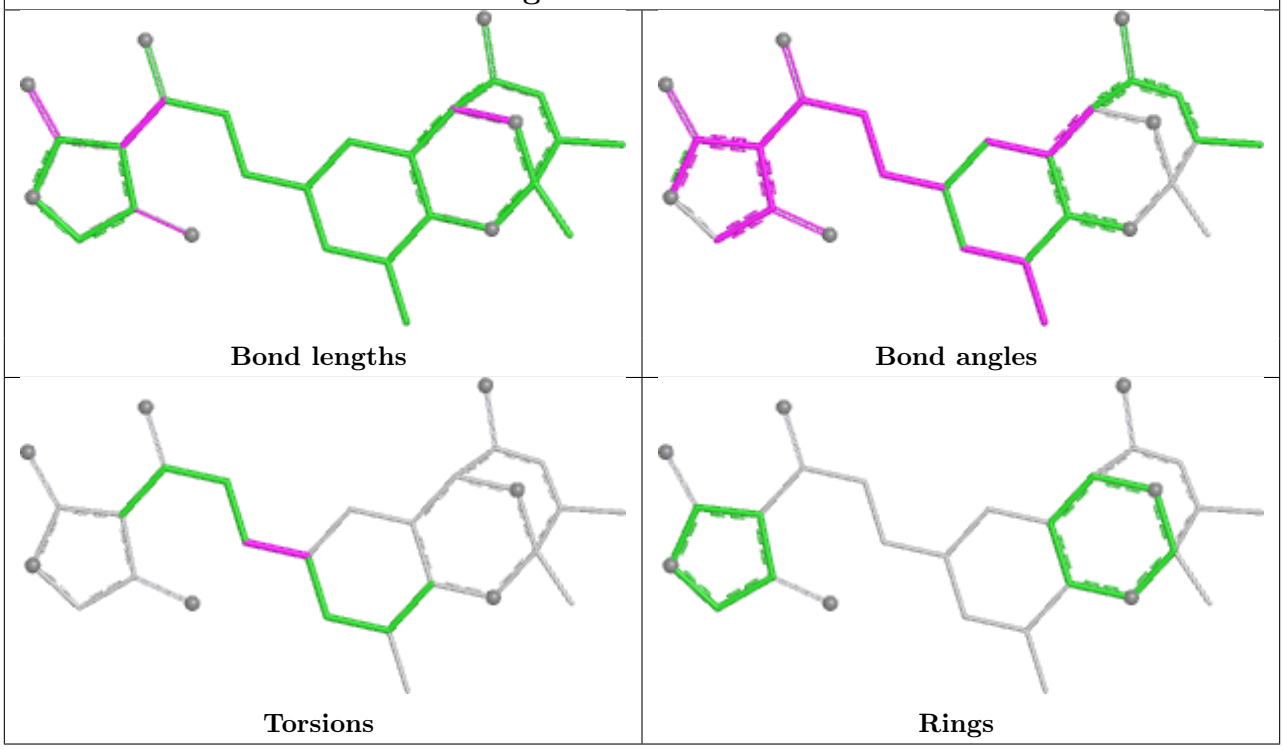
Ligand TRK D 700



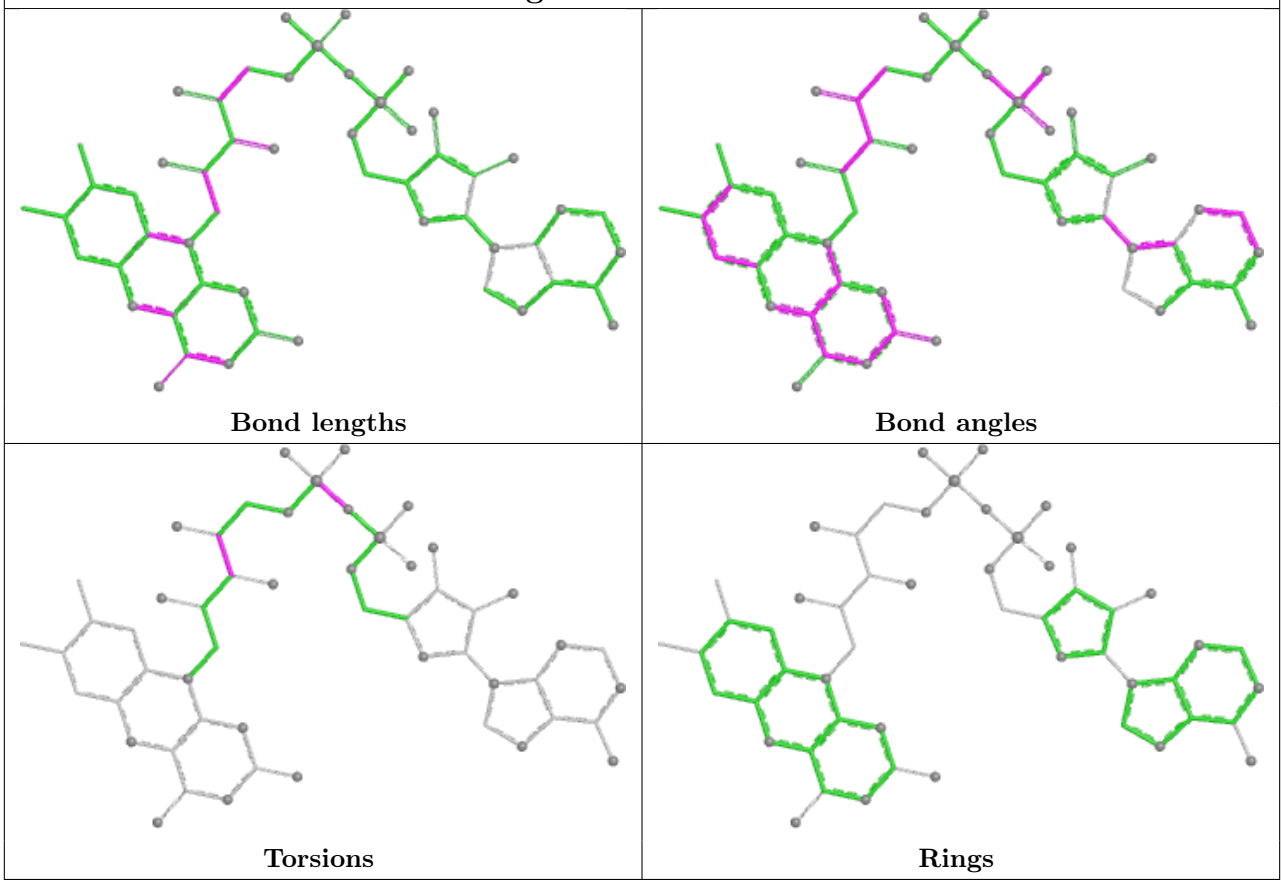
Ligand FAD C 600



Ligand TRK B 700



Ligand FAD B 600



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 (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	492/530 (92%)	-0.24	15 (3%)	50	44	8, 15, 31, 45	0
1	B	487/530 (91%)	-0.26	12 (2%)	57	52	7, 15, 29, 50	0
1	C	492/530 (92%)	-0.23	10 (2%)	65	61	7, 14, 31, 42	0
1	D	491/530 (92%)	-0.20	17 (3%)	44	38	7, 15, 32, 45	0
All	All	1962/2120 (92%)	-0.24	54 (2%)	53	47	7, 15, 31, 50	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	365	TYR	9.8
1	D	500	ALA	8.2
1	C	500	ALA	7.7
1	D	208	PRO	7.1
1	C	368	PRO	7.1
1	A	208	PRO	6.5
1	B	363	THR	5.2
1	A	364	ASP	4.9
1	D	209	PRO	4.9
1	B	369	ALA	4.7
1	D	368	PRO	4.4
1	A	209	PRO	4.1
1	A	204	GLU	3.9
1	B	364	ASP	3.9
1	C	499	PRO	3.8
1	C	363	THR	3.8
1	C	208	PRO	3.8
1	D	50	ARG	3.6
1	A	363	THR	3.5
1	D	90	GLU	3.5
1	C	203	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	89	GLU	3.4
1	D	365	TYR	3.4
1	D	233	GLY	3.4
1	C	364	ASP	3.4
1	B	233	GLY	3.3
1	D	369	ALA	3.2
1	D	366	ASP	3.2
1	D	161	SER	3.1
1	A	499	PRO	3.1
1	B	310	HIS	3.0
1	B	232	GLU	3.0
1	A	4	ILE	2.9
1	A	369	ALA	2.9
1	A	366	ASP	2.9
1	C	233	GLY	2.9
1	A	367	ASN	2.9
1	D	499	PRO	2.7
1	B	499	PRO	2.7
1	B	209	PRO	2.6
1	A	161[A]	SER	2.6
1	D	213	ARG	2.6
1	B	288	PRO	2.5
1	D	363	THR	2.5
1	D	310	HIS	2.4
1	A	213	ARG	2.4
1	C	310	HIS	2.4
1	A	308	GLN	2.2
1	B	43[A]	GLN	2.2
1	C	367	ASN	2.2
1	A	174	SER	2.1
1	B	174	SER	2.1
1	D	176	PRO	2.0
1	B	308	GLN	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

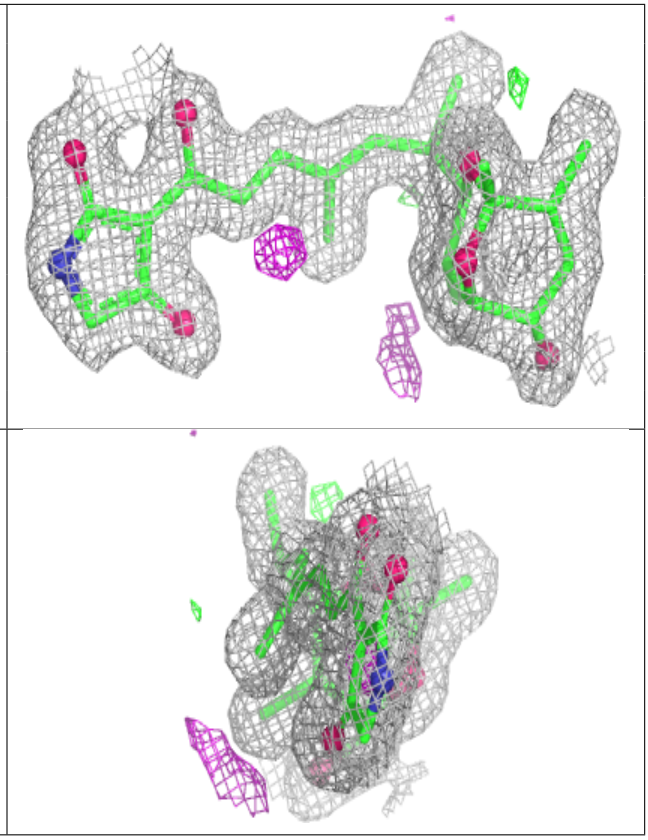
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	GOL	D	1502	6/6	0.87	0.15	27,29,30,30	0
5	GOL	B	1501	6/6	0.90	0.11	28,30,31,33	0
5	GOL	B	1500	6/6	0.90	0.12	31,33,35,35	0
5	GOL	D	1501	6/6	0.91	0.13	29,33,36,38	0
5	GOL	C	1501	6/6	0.93	0.11	31,35,37,40	0
5	GOL	A	1500	6/6	0.95	0.10	30,33,34,38	0
6	CL	D	1504	1/1	0.95	0.07	54,54,54,54	0
3	TIR	C	700	29/29	0.96	0.08	12,15,18,21	0
7	TRK	B	700	29/29	0.96	0.08	14,18,20,20	0
7	TRK	D	700	29/29	0.96	0.07	12,16,20,21	0
5	GOL	C	1502	6/6	0.97	0.16	23,28,31,35	0
3	TIR	A	700	29/29	0.97	0.07	12,15,20,21	0
2	FAD	C	600	53/53	0.98	0.06	5,8,11,13	0
2	FAD	D	600	53/53	0.98	0.06	7,9,11,11	0
2	FAD	A	600	53/53	0.98	0.06	6,9,11,16	0
2	FAD	B	600	53/53	0.98	0.07	6,9,11,14	0
4	MG	D	800	1/1	0.98	0.05	23,23,23,23	0
4	MG	B	800	1/1	0.99	0.09	21,21,21,21	0
4	MG	C	800	1/1	0.99	0.05	19,19,19,19	0
6	CL	A	1501	1/1	0.99	0.04	19,19,19,19	0
6	CL	C	1504	1/1	1.00	0.04	16,16,16,16	0
6	CL	D	1503	1/1	1.00	0.04	19,19,19,19	0
4	MG	A	800	1/1	1.00	0.03	19,19,19,19	0
6	CL	B	1502	1/1	1.00	0.03	18,18,18,18	0
6	CL	C	1503	1/1	1.00	0.04	24,24,24,24	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.

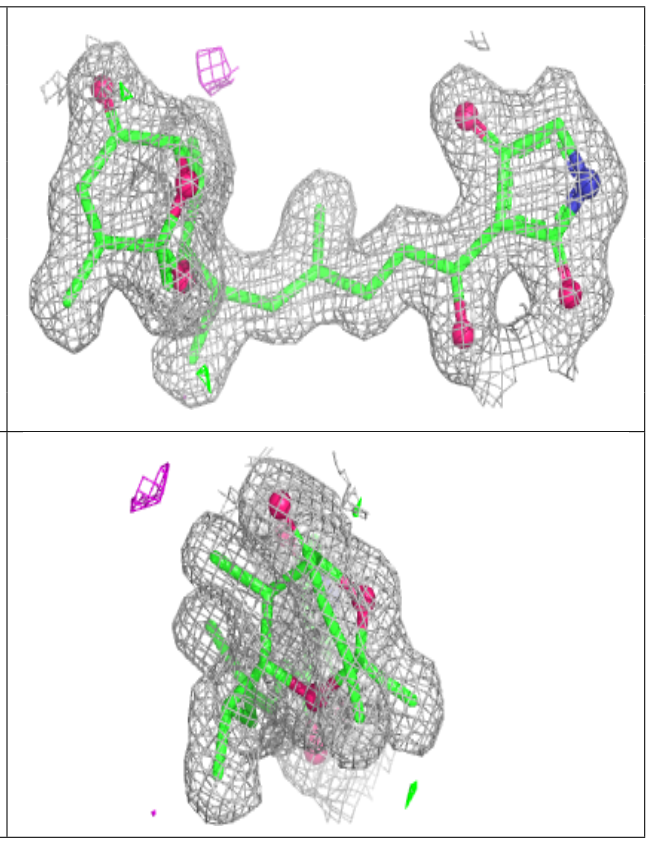
Electron density around TIR C 700:

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



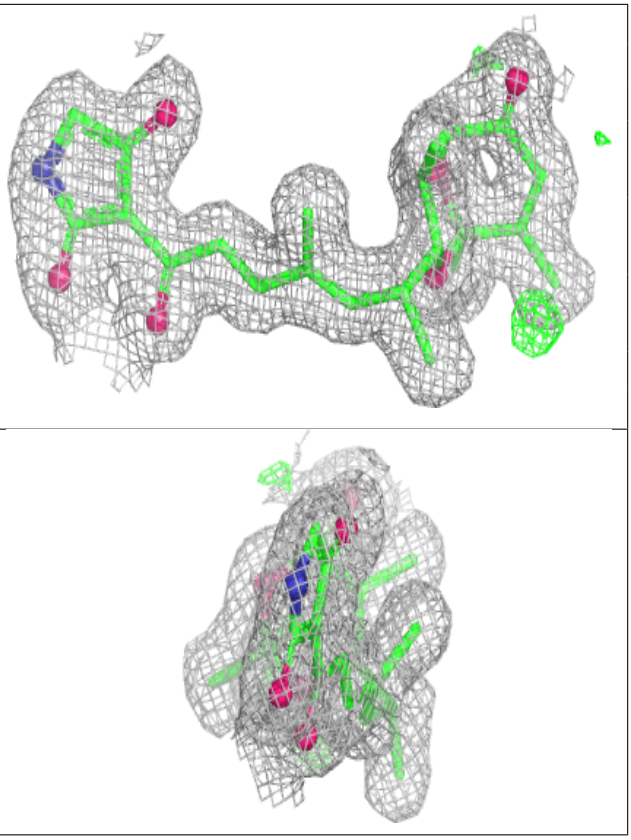
Electron density around TRK B 700:

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



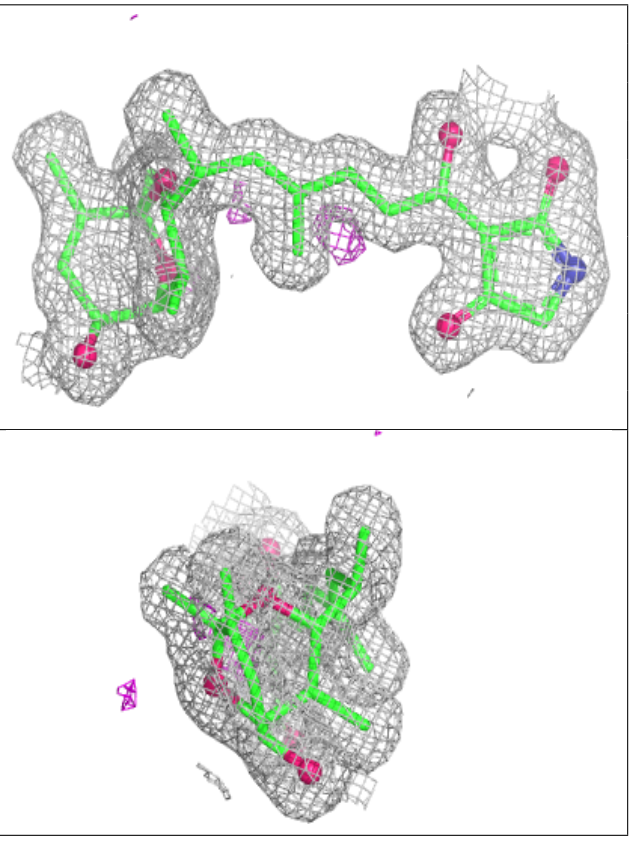
Electron density around TRK D 700:

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



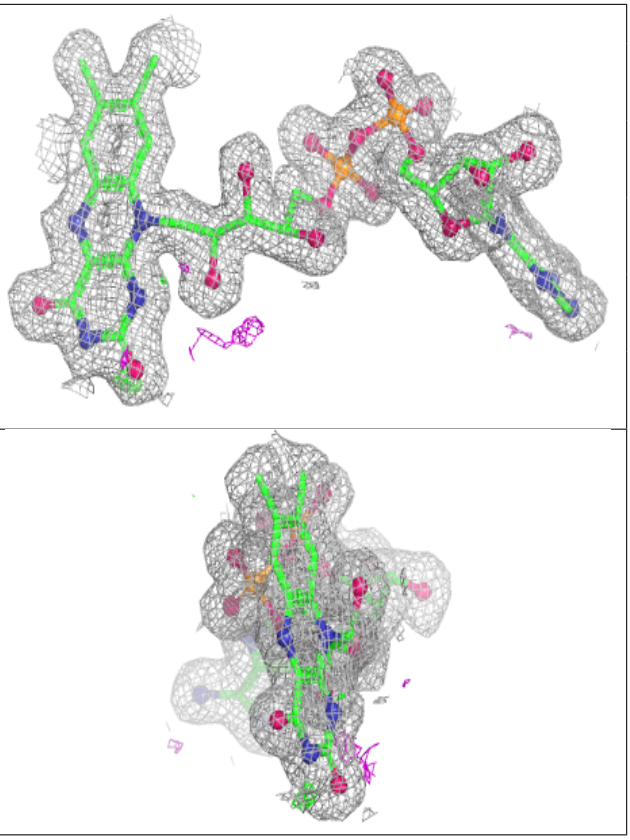
Electron density around TIR A 700:

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



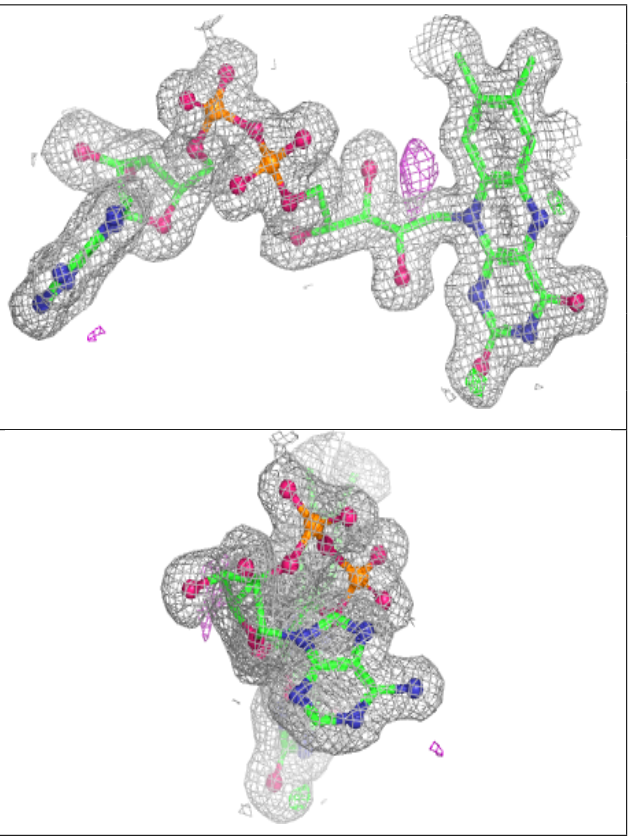
Electron density around FAD C 600:

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



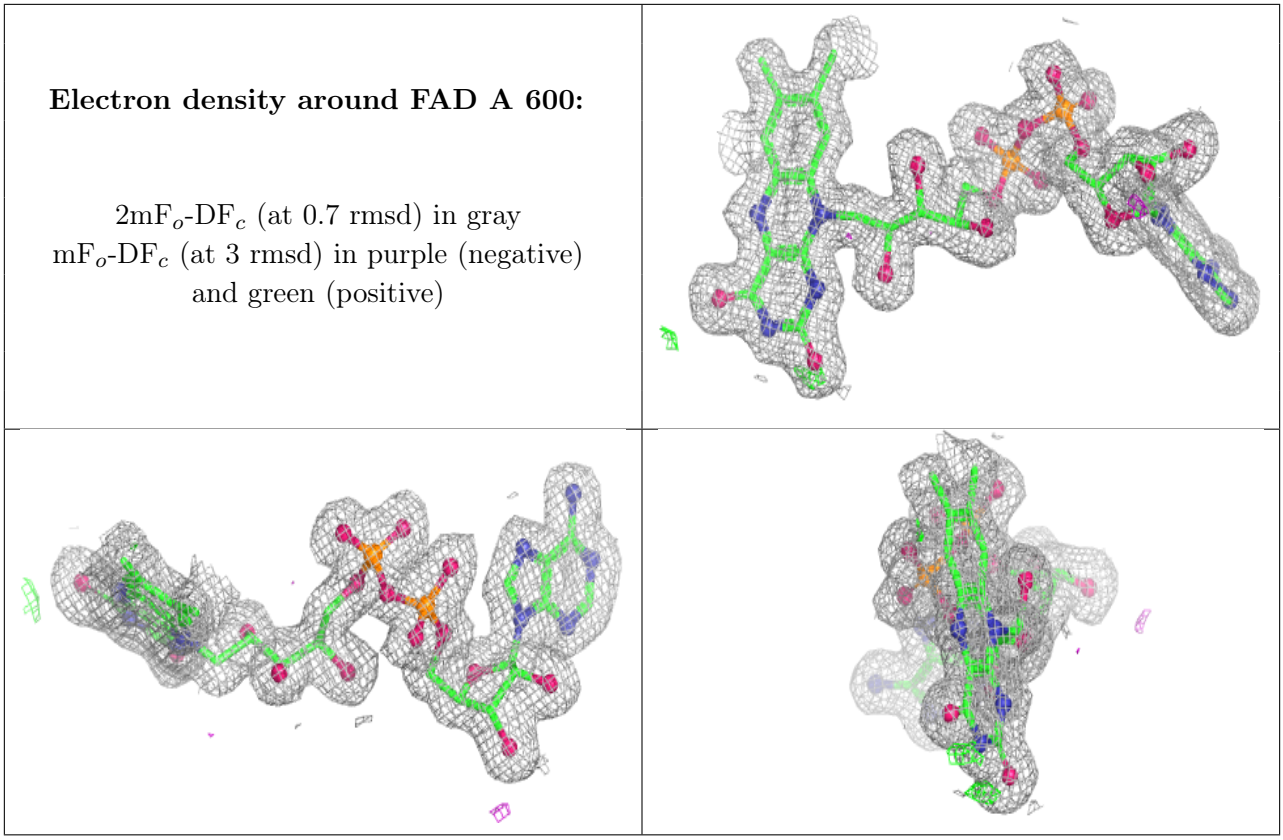
Electron density around FAD D 600:

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



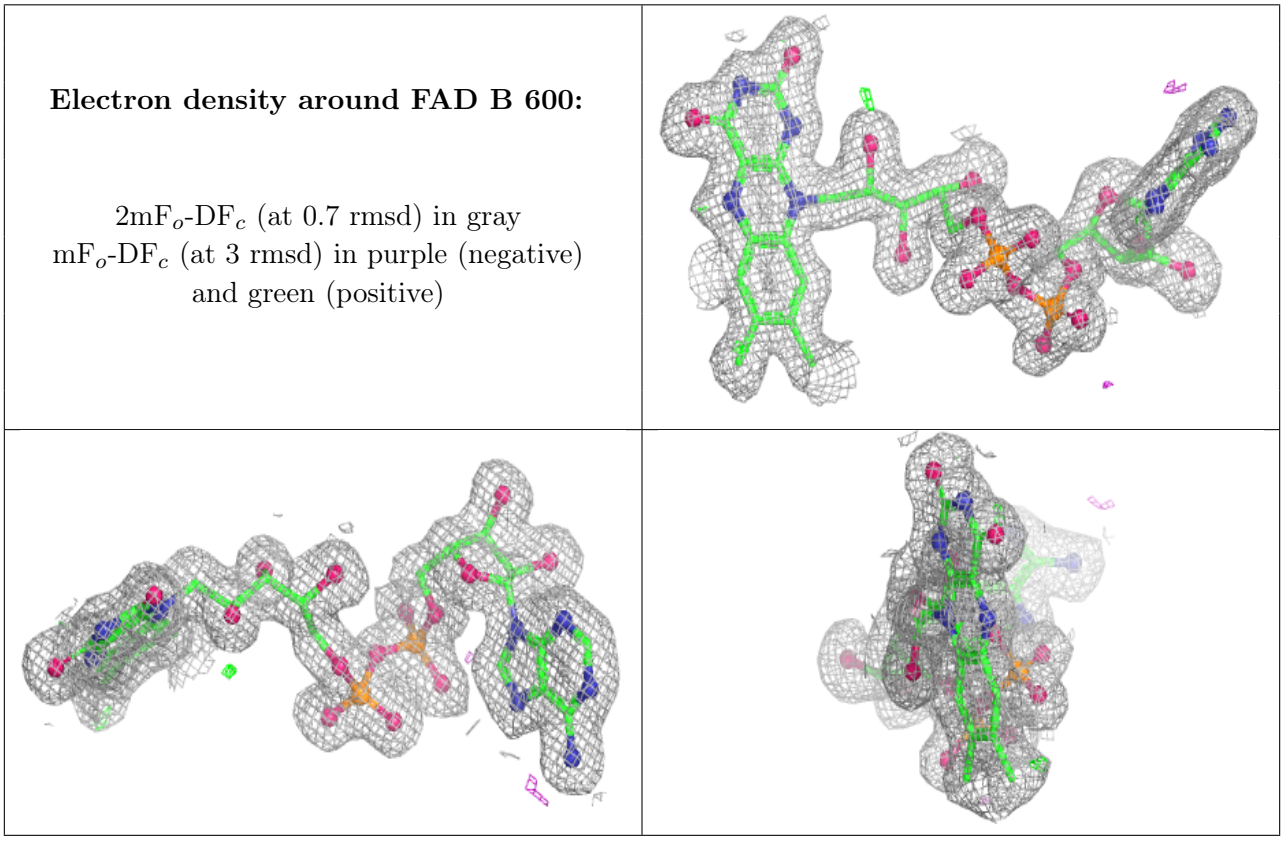
Electron density around FAD A 600:

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



Electron density around FAD B 600:

$2mF_o$-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.