



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 06:30 PM BST

PDB ID : 5YB7
Title : L-Amino acid oxidase/monooxygenase from *Pseudomonas* sp. AIU 813 - L-ornithine complex
Authors : Im, D.; Matsui, D.; Arakawa, T.; Isobe, K.; Asano, Y.; Fushinobu, S.
Deposited on : 2017-09-03
Resolution : 2.00 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 : 2.13.1
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) : 2.13.1

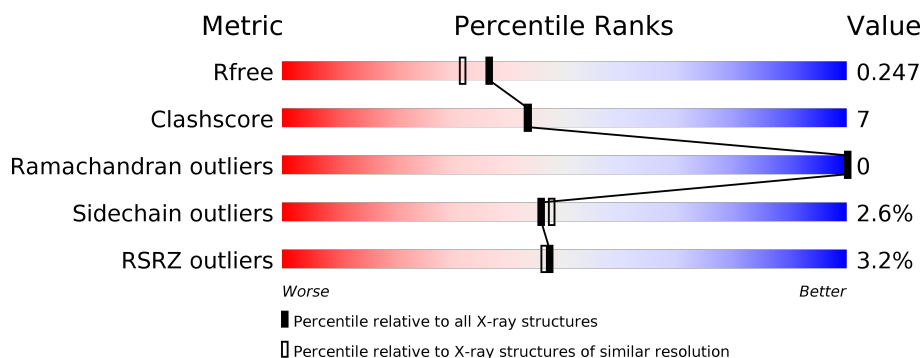
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	580	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 83% 11% • 5% </div> </div>
1	B	580	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 83% 10% • 5% </div> </div>
1	C	580	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 4% 84% 9% • 5% </div> </div>
1	D	580	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 6% 83% 11% • 5% </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18307 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 L-amino acid oxidase/monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	549	Total	C	N	O	S	0	0	0
			4305	2749	742	791	23			
1	B	549	Total	C	N	O	S	0	0	0
			4305	2749	742	791	23			
1	C	549	Total	C	N	O	S	0	0	0
			4305	2749	742	791	23			
1	D	549	Total	C	N	O	S	0	1	0
			4316	2758	743	792	23			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP W6JQJ6
A	-18	GLY	-	expression tag	UNP W6JQJ6
A	-17	SER	-	expression tag	UNP W6JQJ6
A	-16	SER	-	expression tag	UNP W6JQJ6
A	-15	HIS	-	expression tag	UNP W6JQJ6
A	-14	HIS	-	expression tag	UNP W6JQJ6
A	-13	HIS	-	expression tag	UNP W6JQJ6
A	-12	HIS	-	expression tag	UNP W6JQJ6
A	-11	HIS	-	expression tag	UNP W6JQJ6
A	-10	HIS	-	expression tag	UNP W6JQJ6
A	-9	SER	-	expression tag	UNP W6JQJ6
A	-8	SER	-	expression tag	UNP W6JQJ6
A	-7	GLY	-	expression tag	UNP W6JQJ6
A	-6	LEU	-	expression tag	UNP W6JQJ6
A	-5	VAL	-	expression tag	UNP W6JQJ6
A	-4	PRO	-	expression tag	UNP W6JQJ6
A	-3	ARG	-	expression tag	UNP W6JQJ6
A	-2	GLY	-	expression tag	UNP W6JQJ6
A	-1	SER	-	expression tag	UNP W6JQJ6
A	0	HIS	-	expression tag	UNP W6JQJ6
A	473	PHE	SER	conflict	UNP W6JQJ6

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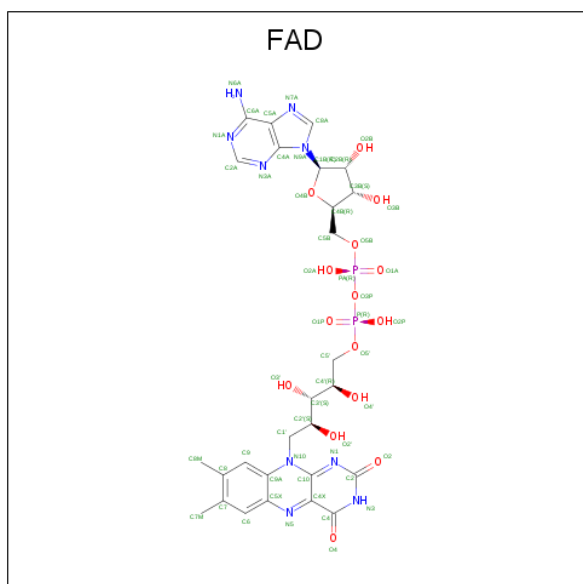
Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	expression tag	UNP W6JQJ6
B	-18	GLY	-	expression tag	UNP W6JQJ6
B	-17	SER	-	expression tag	UNP W6JQJ6
B	-16	SER	-	expression tag	UNP W6JQJ6
B	-15	HIS	-	expression tag	UNP W6JQJ6
B	-14	HIS	-	expression tag	UNP W6JQJ6
B	-13	HIS	-	expression tag	UNP W6JQJ6
B	-12	HIS	-	expression tag	UNP W6JQJ6
B	-11	HIS	-	expression tag	UNP W6JQJ6
B	-10	HIS	-	expression tag	UNP W6JQJ6
B	-9	SER	-	expression tag	UNP W6JQJ6
B	-8	SER	-	expression tag	UNP W6JQJ6
B	-7	GLY	-	expression tag	UNP W6JQJ6
B	-6	LEU	-	expression tag	UNP W6JQJ6
B	-5	VAL	-	expression tag	UNP W6JQJ6
B	-4	PRO	-	expression tag	UNP W6JQJ6
B	-3	ARG	-	expression tag	UNP W6JQJ6
B	-2	GLY	-	expression tag	UNP W6JQJ6
B	-1	SER	-	expression tag	UNP W6JQJ6
B	0	HIS	-	expression tag	UNP W6JQJ6
B	473	PHE	SER	conflict	UNP W6JQJ6
C	-19	MET	-	expression tag	UNP W6JQJ6
C	-18	GLY	-	expression tag	UNP W6JQJ6
C	-17	SER	-	expression tag	UNP W6JQJ6
C	-16	SER	-	expression tag	UNP W6JQJ6
C	-15	HIS	-	expression tag	UNP W6JQJ6
C	-14	HIS	-	expression tag	UNP W6JQJ6
C	-13	HIS	-	expression tag	UNP W6JQJ6
C	-12	HIS	-	expression tag	UNP W6JQJ6
C	-11	HIS	-	expression tag	UNP W6JQJ6
C	-10	HIS	-	expression tag	UNP W6JQJ6
C	-9	SER	-	expression tag	UNP W6JQJ6
C	-8	SER	-	expression tag	UNP W6JQJ6
C	-7	GLY	-	expression tag	UNP W6JQJ6
C	-6	LEU	-	expression tag	UNP W6JQJ6
C	-5	VAL	-	expression tag	UNP W6JQJ6
C	-4	PRO	-	expression tag	UNP W6JQJ6
C	-3	ARG	-	expression tag	UNP W6JQJ6
C	-2	GLY	-	expression tag	UNP W6JQJ6
C	-1	SER	-	expression tag	UNP W6JQJ6
C	0	HIS	-	expression tag	UNP W6JQJ6
C	473	PHE	SER	conflict	UNP W6JQJ6

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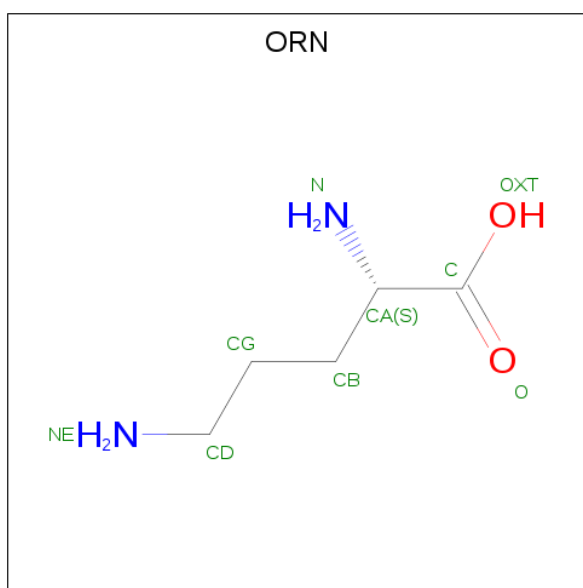
Chain	Residue	Modelled	Actual	Comment	Reference
D	-19	MET	-	expression tag	UNP W6JQJ6
D	-18	GLY	-	expression tag	UNP W6JQJ6
D	-17	SER	-	expression tag	UNP W6JQJ6
D	-16	SER	-	expression tag	UNP W6JQJ6
D	-15	HIS	-	expression tag	UNP W6JQJ6
D	-14	HIS	-	expression tag	UNP W6JQJ6
D	-13	HIS	-	expression tag	UNP W6JQJ6
D	-12	HIS	-	expression tag	UNP W6JQJ6
D	-11	HIS	-	expression tag	UNP W6JQJ6
D	-10	HIS	-	expression tag	UNP W6JQJ6
D	-9	SER	-	expression tag	UNP W6JQJ6
D	-8	SER	-	expression tag	UNP W6JQJ6
D	-7	GLY	-	expression tag	UNP W6JQJ6
D	-6	LEU	-	expression tag	UNP W6JQJ6
D	-5	VAL	-	expression tag	UNP W6JQJ6
D	-4	PRO	-	expression tag	UNP W6JQJ6
D	-3	ARG	-	expression tag	UNP W6JQJ6
D	-2	GLY	-	expression tag	UNP W6JQJ6
D	-1	SER	-	expression tag	UNP W6JQJ6
D	0	HIS	-	expression tag	UNP W6JQJ6
D	473	PHE	SER	conflict	UNP W6JQJ6

- 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
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is L-ornithine (three-letter code: ORN) (formula: $C_5H_{12}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			9	5	2	2		
3	A	1	Total	C	N	O	0	0
			9	5	2	2		
3	B	1	Total	C	N	O	0	0
			9	5	2	2		
3	B	1	Total	C	N	O	0	0
			9	5	2	2		
3	C	1	Total	C	N	O	0	0
			9	5	2	2		
3	C	1	Total	C	N	O	0	0
			9	5	2	2		
3	D	1	Total	C	N	O	0	0
			9	5	2	2		
3	D	1	Total	C	N	O	0	0
			9	5	2	2		

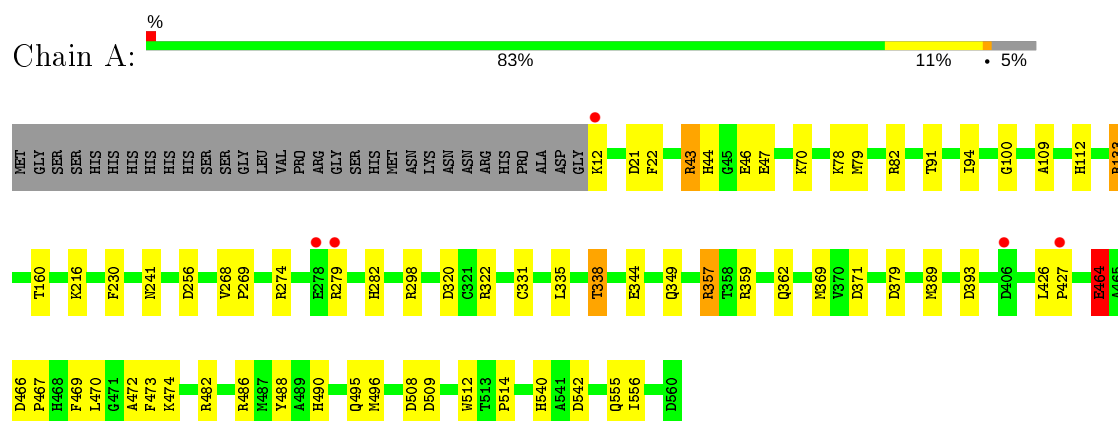
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	244	Total 244	O 244	0	0
4	B	266	Total 266	O 266	0	0
4	C	152	Total 152	O 152	0	0
4	D	130	Total 130	O 130	0	0

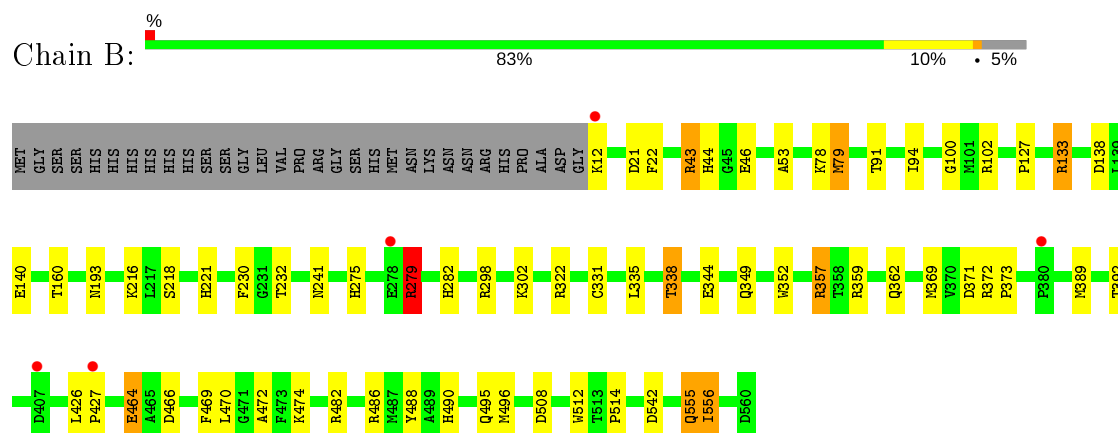
3 Residue-property plots [i](#)

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.

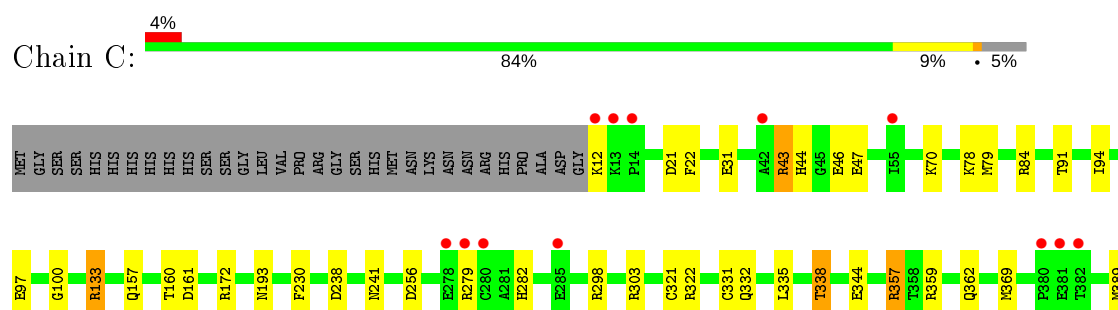
- Molecule 1: L-amino acid oxidase/monooxygenase



- Molecule 1: L-amino acid oxidase/monooxygenase

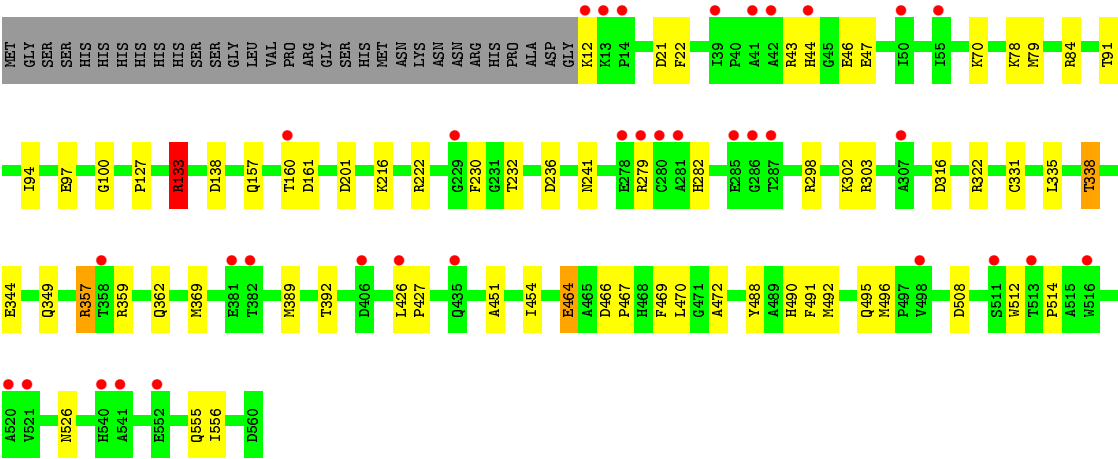
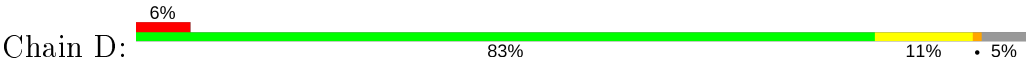


- Molecule 1: L-amino acid oxidase/monooxygenase





• Molecule 1: L-amino acid oxidase/monooxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.60Å 132.90Å 101.90Å 90.00° 111.50° 90.00°	Depositor
Resolution (Å)	94.81 – 2.00 36.69 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.0 (94.81-2.00) 93.1 (36.69-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.49 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.198 , 0.240 0.204 , 0.247	Depositor DCC
R_{free} test set	2015 reflections (1.37%)	wwPDB-VP
Wilson B-factor (Å ²)	29.5	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	18307	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.10% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.92	2/4428 (0.0%)	0.91	11/6019 (0.2%)
1	B	0.90	2/4428 (0.0%)	0.93	9/6019 (0.1%)
1	C	0.79	1/4428 (0.0%)	0.86	8/6019 (0.1%)
1	D	0.76	2/4440 (0.0%)	0.85	8/6035 (0.1%)
All	All	0.85	7/17724 (0.0%)	0.89	36/24092 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	344	GLU	CD-OE2	-7.29	1.17	1.25
1	B	344	GLU	CD-OE2	-6.88	1.18	1.25
1	C	344	GLU	CD-OE1	-6.80	1.18	1.25
1	B	344	GLU	CD-OE1	-6.59	1.18	1.25
1	A	344	GLU	CD-OE2	-5.97	1.19	1.25
1	D	344	GLU	CD-OE1	-5.55	1.19	1.25
1	A	344	GLU	CD-OE1	-5.05	1.20	1.25

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	344	GLU	OE1-CD-OE2	-9.92	111.40	123.30
1	B	496	MET	CG-SD-CE	-9.13	85.59	100.20
1	D	344	GLU	OE1-CD-OE2	-8.80	112.74	123.30
1	C	344	GLU	OE1-CD-OE2	-8.74	112.81	123.30
1	A	344	GLU	OE1-CD-OE2	-8.74	112.81	123.30
1	B	322	ARG	NE-CZ-NH2	-8.65	115.97	120.30
1	A	322	ARG	NE-CZ-NH2	-8.48	116.06	120.30
1	A	496	MET	CG-SD-CE	-8.41	86.75	100.20
1	B	43	ARG	NE-CZ-NH1	-7.55	116.52	120.30
1	B	298	ARG	NE-CZ-NH2	-7.37	116.61	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	298	ARG	NE-CZ-NH2	-7.31	116.65	120.30
1	C	322	ARG	NE-CZ-NH2	-7.23	116.69	120.30
1	D	496	MET	CG-SD-CE	-7.03	88.95	100.20
1	C	298	ARG	NE-CZ-NH2	-6.99	116.81	120.30
1	A	393	ASP	CB-CG-OD1	6.71	124.33	118.30
1	D	322	ARG	NE-CZ-NH2	-6.63	116.99	120.30
1	C	496	MET	CG-SD-CE	-6.60	89.65	100.20
1	A	371	ASP	CB-CG-OD1	6.31	123.97	118.30
1	D	298	ARG	NE-CZ-NH2	-6.29	117.16	120.30
1	A	82	ARG	NE-CZ-NH1	6.03	123.31	120.30
1	C	464	GLU	CB-CA-C	-5.99	98.42	110.40
1	B	279	ARG	NE-CZ-NH1	5.97	123.28	120.30
1	A	43	ARG	NE-CZ-NH1	-5.83	117.38	120.30
1	D	464	GLU	CB-CA-C	-5.74	98.91	110.40
1	D	133	ARG	NE-CZ-NH1	5.70	123.15	120.30
1	B	371	ASP	CB-CG-OD1	5.64	123.37	118.30
1	A	464	GLU	CB-CA-C	-5.63	99.14	110.40
1	C	298	ARG	NE-CZ-NH1	5.38	122.99	120.30
1	B	102	ARG	NE-CZ-NH1	5.35	122.98	120.30
1	D	322	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	A	320	ASP	CB-CG-OD2	5.21	122.99	118.30
1	B	464	GLU	CB-CA-C	-5.21	99.99	110.40
1	C	322	ARG	NE-CZ-NH1	5.21	122.90	120.30
1	D	236	ASP	CB-CG-OD1	5.12	122.90	118.30
1	C	43	ARG	NE-CZ-NH1	-5.09	117.76	120.30
1	A	274	ARG	NE-CZ-NH1	5.04	122.82	120.30

There are no chirality outliers.

There are no planarity outliers.

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	4305	0	4166	54	1
1	B	4305	0	4166	64	1
1	C	4305	0	4166	55	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	4316	0	4174	59	0
2	A	53	0	31	4	0
2	B	53	0	31	4	0
2	C	53	0	31	4	0
2	D	53	0	31	4	0
3	A	18	0	22	1	0
3	B	18	0	22	2	0
3	C	18	0	22	2	0
3	D	18	0	22	1	0
4	A	244	0	0	2	0
4	B	266	0	0	10	0
4	C	152	0	0	8	0
4	D	130	0	0	8	0
All	All	18307	0	16884	229	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 7.

All (229) 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:B:555:GLN:OE1	1:B:555:GLN:N	1.61	1.31
1:C:555:GLN:N	1:C:555:GLN:OE1	1.63	1.26
1:D:555:GLN:OE1	1:D:555:GLN:N	1.69	1.25
1:A:555:GLN:N	1:A:555:GLN:OE1	1.69	1.25
1:A:555:GLN:H	1:A:555:GLN:CD	1.58	1.04
1:C:555:GLN:CD	1:C:555:GLN:H	1.58	1.03
1:B:555:GLN:CD	1:B:555:GLN:H	1.62	1.02
1:C:22:PHE:HB2	1:C:555:GLN:NE2	1.76	1.01
1:D:555:GLN:CD	1:D:555:GLN:H	1.63	0.99
1:C:466:ASP:O	1:C:469:PHE:O	1.81	0.98
1:A:22:PHE:HB2	1:A:555:GLN:NE2	1.80	0.97
1:D:22:PHE:HB2	1:D:555:GLN:NE2	1.81	0.95
1:A:466:ASP:O	1:A:469:PHE:O	1.85	0.95
1:A:94:ILE:HG21	1:A:369:MET:CE	1.97	0.94
1:D:466:ASP:O	1:D:469:PHE:O	1.86	0.94
1:B:466:ASP:O	1:B:469:PHE:O	1.85	0.94
1:B:94:ILE:HG21	1:B:369:MET:CE	1.99	0.91
1:B:279:ARG:NH2	4:B:701:HOH:O	1.97	0.89
1:B:22:PHE:HB2	1:B:555:GLN:NE2	1.89	0.87
1:A:94:ILE:HG21	1:A:369:MET:HE2	1.57	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:508:ASP:OD2	1:B:514:PRO:O	1.93	0.85
1:A:508:ASP:OD2	1:A:514:PRO:O	1.94	0.85
1:B:302:LYS:HE3	4:B:705:HOH:O	1.80	0.82
1:D:302:LYS:HE3	4:D:1514:HOH:O	1.79	0.81
1:B:555:GLN:CD	1:B:555:GLN:N	2.25	0.81
1:D:508:ASP:OD2	1:D:514:PRO:O	1.99	0.81
1:C:94:ILE:HG21	1:C:369:MET:CE	2.11	0.80
1:C:472:ALA:HB1	2:C:601:FAD:HM83	1.65	0.79
1:C:508:ASP:OD2	1:C:514:PRO:O	2.02	0.78
1:D:94:ILE:HG21	1:D:369:MET:CE	2.14	0.77
1:B:94:ILE:HG21	1:B:369:MET:HE2	1.67	0.74
1:C:321:CYS:SG	4:C:809:HOH:O	2.45	0.74
1:D:472:ALA:HB1	2:D:601:FAD:HM83	1.69	0.73
1:C:555:GLN:N	1:C:555:GLN:CD	2.24	0.72
1:A:241:ASN:OD1	1:A:359:ARG:NH2	2.24	0.71
1:B:241:ASN:OD1	1:B:359:ARG:NH2	2.25	0.70
1:D:94:ILE:HG21	1:D:369:MET:HE2	1.73	0.70
1:B:362:GLN:HG3	1:B:464:GLU:CG	2.22	0.70
1:A:362:GLN:HG3	1:A:464:GLU:CG	2.22	0.69
1:B:362:GLN:CG	1:B:464:GLU:HG2	2.23	0.69
1:D:241:ASN:OD1	1:D:359:ARG:NH2	2.25	0.69
1:D:555:GLN:CD	1:D:555:GLN:N	2.27	0.69
1:C:241:ASN:OD1	1:C:359:ARG:NH2	2.28	0.67
1:B:94:ILE:CG2	1:B:369:MET:CE	2.73	0.67
1:C:94:ILE:HG21	1:C:369:MET:HE2	1.75	0.67
1:B:472:ALA:HB1	2:B:601:FAD:HM83	1.77	0.66
1:A:94:ILE:CG2	1:A:369:MET:CE	2.72	0.66
1:B:362:GLN:HG3	1:B:464:GLU:HG2	1.77	0.65
1:A:362:GLN:CG	1:A:464:GLU:HG2	2.27	0.65
1:C:362:GLN:HG3	1:C:464:GLU:CG	2.28	0.64
1:A:362:GLN:HG3	1:A:464:GLU:HG3	1.80	0.63
1:C:362:GLN:CG	1:C:464:GLU:HG2	2.28	0.63
1:B:302:LYS:CE	4:B:705:HOH:O	2.43	0.63
1:A:555:GLN:N	1:A:555:GLN:CD	2.26	0.63
1:D:362:GLN:HG3	1:D:464:GLU:CG	2.30	0.62
1:A:472:ALA:HB1	2:A:601:FAD:HM83	1.82	0.61
1:A:349:GLN:HE22	3:A:603:ORN:HE2	1.49	0.61
1:D:490:HIS:HD2	4:D:1613:HOH:O	1.84	0.59
1:D:362:GLN:CG	1:D:464:GLU:HG2	2.32	0.59
1:D:84:ARG:HD3	4:D:1626:HOH:O	2.03	0.59
1:A:362:GLN:HG3	1:A:464:GLU:HG2	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:362:GLN:HG3	1:D:464:GLU:HG3	1.85	0.58
1:A:512:TRP:CZ2	1:A:555:GLN:HG2	2.38	0.58
1:D:22:PHE:HB2	1:D:555:GLN:HE21	1.68	0.58
1:B:302:LYS:NZ	4:B:705:HOH:O	2.37	0.58
1:A:426:LEU:HD21	1:C:426:LEU:HD21	1.86	0.57
1:A:512:TRP:HZ2	1:A:555:GLN:HG2	1.69	0.57
1:D:512:TRP:HZ2	1:D:555:GLN:HG2	1.70	0.57
1:D:491[B]:PHE:CD1	1:D:526:ASN:O	2.58	0.57
1:B:94:ILE:HG21	1:B:369:MET:HE1	1.84	0.57
1:B:362:GLN:HG3	1:B:464:GLU:HG3	1.87	0.56
1:D:302:LYS:CE	4:D:1514:HOH:O	2.47	0.56
1:A:362:GLN:CG	1:A:464:GLU:CG	2.83	0.55
1:D:512:TRP:CZ2	1:D:555:GLN:HG2	2.41	0.55
1:C:362:GLN:HG3	1:C:464:GLU:HG2	1.86	0.55
1:B:362:GLN:CG	1:B:464:GLU:CG	2.84	0.55
1:C:362:GLN:HG3	1:C:464:GLU:HG3	1.88	0.54
1:D:94:ILE:CG2	1:D:369:MET:CE	2.86	0.54
1:D:43:ARG:HD2	1:D:46:GLU:OE2	2.07	0.54
1:C:472:ALA:CB	2:C:601:FAD:HM83	2.37	0.54
1:A:22:PHE:HB2	1:A:555:GLN:HE22	1.71	0.54
1:C:43:ARG:HD2	1:C:46:GLU:OE2	2.08	0.54
1:A:335:LEU:HA	1:A:338:THR:HG22	1.91	0.53
1:C:22:PHE:HB2	1:C:555:GLN:HE22	1.68	0.53
1:C:94:ILE:CG2	1:C:369:MET:CE	2.84	0.53
1:C:426:LEU:HB3	1:C:427:PRO:HD3	1.90	0.53
1:C:512:TRP:HZ2	1:C:555:GLN:HG2	1.73	0.53
1:D:491[B]:PHE:CE2	1:D:526:ASN:HB3	2.44	0.53
1:C:512:TRP:CZ2	1:C:555:GLN:HG2	2.43	0.53
1:D:491[A]:PHE:CE1	1:D:492:MET:SD	3.01	0.52
1:D:349:GLN:HE22	3:D:603:ORN:HE2	1.57	0.52
1:D:426:LEU:HB3	1:D:427:PRO:HD3	1.91	0.52
1:D:362:GLN:CG	1:D:464:GLU:CG	2.87	0.52
1:C:488:TYR:CE2	1:C:555:GLN:NE2	2.78	0.52
1:B:488:TYR:CE2	1:B:555:GLN:NE2	2.78	0.52
1:D:22:PHE:HB2	1:D:555:GLN:HE22	1.73	0.52
1:D:303:ARG:CZ	4:D:1590:HOH:O	2.58	0.52
1:B:349:GLN:HE22	3:B:603:ORN:HE2	1.58	0.51
1:A:338:THR:HG21	1:A:470:LEU:HD11	1.92	0.51
1:B:426:LEU:HB3	1:B:427:PRO:HD3	1.93	0.50
1:C:21:ASP:HB3	1:C:555:GLN:HG3	1.94	0.50
1:B:335:LEU:HA	1:B:338:THR:HG22	1.91	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:362:GLN:CG	1:C:464:GLU:CG	2.87	0.50
1:D:222:ARG:NH2	4:D:1507:HOH:O	2.43	0.50
1:C:84:ARG:HD3	4:C:848:HOH:O	2.11	0.50
1:B:100:GLY:HA2	2:B:601:FAD:N5	2.26	0.50
1:B:512:TRP:CZ2	1:B:555:GLN:HG2	2.46	0.50
1:C:157:GLN:NE2	1:C:161:ASP:OD1	2.42	0.50
1:A:488:TYR:CE2	1:A:555:GLN:NE2	2.80	0.50
1:C:22:PHE:HB2	1:C:555:GLN:HE21	1.69	0.50
1:A:426:LEU:HB3	1:A:427:PRO:HD3	1.94	0.49
1:B:160:THR:OG1	1:B:230:PHE:HB2	2.12	0.49
1:C:490:HIS:HD2	4:C:825:HOH:O	1.94	0.49
1:C:172:ARG:HD3	4:C:769:HOH:O	2.11	0.49
1:A:490:HIS:HE1	4:A:705:HOH:O	1.94	0.49
1:A:22:PHE:HB2	1:A:555:GLN:HE21	1.71	0.49
1:C:418:TRP:CZ2	3:C:602:ORN:HD2	2.47	0.49
1:D:100:GLY:HA2	2:D:601:FAD:N5	2.28	0.49
1:D:335:LEU:HA	1:D:338:THR:HG22	1.93	0.49
1:D:338:THR:HG21	1:D:470:LEU:HD11	1.95	0.48
1:D:362:GLN:HG3	1:D:464:GLU:HG2	1.93	0.48
1:A:21:ASP:HB3	1:A:555:GLN:HG3	1.95	0.48
1:A:357:ARG:HH11	1:A:357:ARG:HA	1.78	0.48
1:A:490:HIS:HD2	4:A:853:HOH:O	1.97	0.48
1:A:542:ASP:N	1:A:542:ASP:OD1	2.46	0.48
1:B:490:HIS:HD2	4:B:870:HOH:O	1.96	0.48
1:C:160:THR:OG1	1:C:230:PHE:HB2	2.13	0.48
1:B:426:LEU:HD21	1:D:426:LEU:HD21	1.94	0.48
1:D:100:GLY:HA2	2:D:601:FAD:C5X	2.44	0.48
1:C:335:LEU:HA	1:C:338:THR:HG22	1.94	0.48
1:A:482:ARG:O	1:A:486:ARG:HG3	2.14	0.48
1:B:357:ARG:HH11	1:B:357:ARG:HA	1.77	0.48
1:A:160:THR:OG1	1:A:230:PHE:HB2	2.14	0.47
1:D:157:GLN:NE2	1:D:161:ASP:OD1	2.46	0.47
1:C:357:ARG:HA	1:C:357:ARG:HH11	1.80	0.47
1:B:43:ARG:HD2	1:B:46:GLU:OE2	2.14	0.47
1:A:94:ILE:HG21	1:A:369:MET:HE1	1.94	0.47
1:D:160:THR:OG1	1:D:230:PHE:HB2	2.14	0.47
1:B:21:ASP:HB3	1:B:555:GLN:HG3	1.96	0.47
1:B:279:ARG:NH1	4:B:709:HOH:O	2.42	0.47
1:B:512:TRP:HZ2	1:B:555:GLN:HG2	1.79	0.47
1:D:472:ALA:CB	2:D:601:FAD:HM83	2.41	0.47
1:B:22:PHE:HB2	1:B:555:GLN:HE21	1.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:21:ASP:HB3	1:D:555:GLN:HG3	1.97	0.47
1:C:466:ASP:OD1	1:C:467:PRO:HD2	2.15	0.46
1:D:331:CYS:HB2	1:D:335:LEU:HD12	1.96	0.46
1:B:140:GLU:HA	1:D:349:GLN:HE21	1.80	0.46
1:C:100:GLY:HA2	2:C:601:FAD:N5	2.29	0.46
1:B:338:THR:HG21	1:B:470:LEU:HD11	1.97	0.46
1:B:474:LYS:HB3	1:B:514:PRO:HB3	1.97	0.46
1:C:338:THR:HG21	1:C:470:LEU:HD11	1.98	0.46
1:D:488:TYR:CE2	1:D:555:GLN:NE2	2.84	0.46
1:B:275:HIS:HD2	4:B:954:HOH:O	1.98	0.46
1:B:218:SER:OG	1:B:221:HIS:ND1	2.46	0.46
1:B:472:ALA:CB	2:B:601:FAD:HM83	2.45	0.46
1:D:91:THR:HG21	1:D:369:MET:CE	2.46	0.46
1:B:133:ARG:HH11	1:B:133:ARG:HB3	1.81	0.46
1:B:53:ALA:HB1	1:B:79:MET:HE2	1.98	0.45
1:A:335:LEU:HD23	1:A:338:THR:HG22	1.97	0.45
1:A:43:ARG:HD2	1:A:46:GLU:OE2	2.15	0.45
1:D:44:HIS:HB3	1:D:282:HIS:CG	2.52	0.45
1:C:91:THR:HG21	1:C:369:MET:CE	2.47	0.45
1:B:335:LEU:HD23	1:B:338:THR:HG22	1.98	0.45
1:C:94:ILE:HG21	1:C:369:MET:HE1	1.96	0.45
1:B:94:ILE:CG2	1:B:369:MET:HE2	2.42	0.45
1:A:100:GLY:HA2	2:A:601:FAD:N5	2.32	0.45
1:B:193:ASN:ND2	1:B:556:ILE:CD1	2.80	0.45
1:C:44:HIS:HB3	1:C:282:HIS:CG	2.52	0.45
1:A:91:THR:HG21	1:A:369:MET:HE1	1.99	0.45
1:C:238:ASP:OD2	3:C:602:ORN:NE	2.45	0.45
1:D:47:GLU:HA	1:D:70:LYS:O	2.17	0.45
1:B:490:HIS:HE1	4:B:710:HOH:O	1.99	0.44
1:C:490:HIS:HE1	4:C:714:HOH:O	1.99	0.44
1:D:466:ASP:OD1	1:D:467:PRO:HD2	2.17	0.44
1:A:47:GLU:HA	1:A:70:LYS:O	2.18	0.44
1:B:482:ARG:O	1:B:486:ARG:HG3	2.17	0.44
1:A:133:ARG:HB3	1:A:133:ARG:HH11	1.83	0.44
1:A:474:LYS:HB3	1:A:514:PRO:HB3	1.98	0.44
1:B:91:THR:HG21	1:B:369:MET:CE	2.48	0.44
1:D:451:ALA:HA	1:D:454:ILE:HD12	1.99	0.44
1:C:100:GLY:HA2	2:C:601:FAD:C5X	2.48	0.44
1:A:509:ASP:N	1:A:509:ASP:OD1	2.51	0.43
1:A:91:THR:HG21	1:A:369:MET:CE	2.48	0.43
1:D:357:ARG:HA	1:D:357:ARG:HH11	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:LEU:HD23	1:A:338:THR:CG2	2.48	0.43
1:B:44:HIS:HB3	1:B:282:HIS:CG	2.52	0.43
1:B:279:ARG:CZ	4:B:701:HOH:O	2.52	0.43
1:C:97:GLU:OE1	1:C:97:GLU:HA	2.18	0.43
1:D:133:ARG:HH11	1:D:133:ARG:HB3	1.84	0.43
1:A:466:ASP:OD1	1:A:467:PRO:HD2	2.19	0.43
1:B:362:GLN:HG2	1:B:464:GLU:HG2	2.00	0.43
1:C:47:GLU:HA	1:C:70:LYS:O	2.19	0.43
1:B:372:ARG:HB2	1:B:373:PRO:HD2	2.01	0.43
1:C:303:ARG:CZ	4:C:812:HOH:O	2.67	0.42
1:D:127:PRO:HG2	1:D:232:THR:HA	2.01	0.42
1:B:335:LEU:HD23	1:B:338:THR:CG2	2.50	0.42
1:A:331:CYS:HB2	1:A:335:LEU:HD12	2.01	0.42
1:C:332:GLN:NE2	4:C:702:HOH:O	2.38	0.42
1:B:193:ASN:CG	1:B:556:ILE:HD13	2.40	0.42
1:D:138:ASP:O	1:D:392:THR:HA	2.19	0.42
1:C:256:ASP:OD1	1:C:256:ASP:N	2.53	0.42
1:A:109:ALA:O	1:A:112:HIS:HB3	2.19	0.42
1:C:331:CYS:HB2	1:C:335:LEU:HD12	2.01	0.42
1:A:44:HIS:HB3	1:A:282:HIS:CG	2.55	0.42
1:B:127:PRO:HG2	1:B:232:THR:HA	2.02	0.42
1:A:362:GLN:HG2	1:A:464:GLU:HG2	2.01	0.41
1:A:100:GLY:HA2	2:A:601:FAD:C5X	2.50	0.41
1:D:302:LYS:NZ	4:D:1514:HOH:O	2.51	0.41
1:D:362:GLN:HG2	1:D:464:GLU:HG2	2.02	0.41
1:B:542:ASP:N	1:B:542:ASP:OD1	2.46	0.41
1:B:279:ARG:NH1	4:B:701:HOH:O	2.54	0.41
1:B:22:PHE:HB2	1:B:555:GLN:HE22	1.79	0.41
1:D:201:ASP:HB2	4:D:1543:HOH:O	2.20	0.41
1:A:256:ASP:OD1	1:A:256:ASP:N	2.54	0.41
1:C:31:GLU:CG	4:C:708:HOH:O	2.68	0.41
1:C:362:GLN:HG2	1:C:464:GLU:HG2	2.03	0.41
1:D:426:LEU:HB3	1:D:427:PRO:CD	2.51	0.41
1:B:100:GLY:HA2	2:B:601:FAD:C4X	2.50	0.41
1:A:379:ASP:C	1:A:379:ASP:OD1	2.59	0.41
1:D:316:ASP:C	1:D:316:ASP:OD1	2.60	0.41
1:B:352:TRP:HB3	3:B:603:ORN:HD2	2.03	0.40
1:B:331:CYS:HB2	1:B:335:LEU:HD12	2.02	0.40
1:C:193:ASN:CG	1:C:556:ILE:HD13	2.42	0.40
1:A:268:VAL:HB	1:A:269:PRO:CD	2.52	0.40
1:D:97:GLU:HA	1:D:97:GLU:OE1	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:472:ALA:CB	2:A:601:FAD:HM83	2.50	0.40
1:B:138:ASP:O	1:B:392:THR:HA	2.22	0.40
1:C:133:ARG:HH11	1:C:133:ARG:HB3	1.86	0.40
1:C:425:MET:O	1:C:426:LEU:C	2.60	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)
1:A:540:HIS:CD2	1:B:279:ARG:NH2[2_556]	1.90	0.30

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	547/580 (94%)	535 (98%)	12 (2%)	0	100	100
1	B	547/580 (94%)	536 (98%)	11 (2%)	0	100	100
1	C	547/580 (94%)	533 (97%)	14 (3%)	0	100	100
1	D	548/580 (94%)	534 (97%)	14 (3%)	0	100	100
All	All	2189/2320 (94%)	2138 (98%)	51 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	448/474 (94%)	435 (97%)	13 (3%)	42	43
1	B	448/474 (94%)	436 (97%)	12 (3%)	44	46
1	C	448/474 (94%)	438 (98%)	10 (2%)	52	55
1	D	449/474 (95%)	438 (98%)	11 (2%)	49	51
All	All	1793/1896 (95%)	1747 (97%)	46 (3%)	46	48

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LYS
1	A	78	LYS
1	A	79	MET
1	A	133	ARG
1	A	216	LYS
1	A	279	ARG
1	A	338	THR
1	A	357	ARG
1	A	389	MET
1	A	464	GLU
1	A	473	PHE
1	A	495	GLN
1	A	556	ILE
1	B	12	LYS
1	B	78	LYS
1	B	79	MET
1	B	133	ARG
1	B	216	LYS
1	B	279	ARG
1	B	338	THR
1	B	357	ARG
1	B	389	MET
1	B	495	GLN
1	B	555	GLN
1	B	556	ILE
1	C	12	LYS
1	C	78	LYS
1	C	79	MET
1	C	133	ARG
1	C	279	ARG
1	C	338	THR
1	C	357	ARG
1	C	389	MET

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Mol	Chain	Res	Type
1	C	495	GLN
1	C	556	ILE
1	D	12	LYS
1	D	78	LYS
1	D	79	MET
1	D	133	ARG
1	D	216	LYS
1	D	279	ARG
1	D	338	THR
1	D	357	ARG
1	D	389	MET
1	D	495	GLN
1	D	556	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	332	GLN
1	A	349	GLN
1	A	484	ASN
1	A	490	HIS
1	A	533	ASN
1	A	540	HIS
1	B	275	HIS
1	B	332	GLN
1	B	349	GLN
1	B	484	ASN
1	B	490	HIS
1	B	540	HIS
1	C	349	GLN
1	C	484	ASN
1	C	490	HIS
1	D	332	GLN
1	D	349	GLN
1	D	484	ASN
1	D	490	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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

12 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	$\# Z > 2$
3	ORN	C	603	-	4,8,8	0.63	0	3,9,9	1.30	1 (33%)
3	ORN	D	602	-	4,8,8	0.53	0	3,9,9	0.89	0
3	ORN	D	603	-	4,8,8	0.68	0	3,9,9	0.25	0
2	FAD	B	601	-	51,58,58	1.95	10 (19%)	60,89,89	2.89	21 (35%)
3	ORN	B	603	-	4,8,8	0.70	0	3,9,9	0.37	0
2	FAD	D	601	-	51,58,58	2.11	11 (21%)	60,89,89	2.38	19 (31%)
3	ORN	A	602	-	4,8,8	0.93	0	3,9,9	0.66	0
3	ORN	B	602	-	4,8,8	0.95	0	3,9,9	0.30	0
3	ORN	C	602	-	4,8,8	0.64	0	3,9,9	0.87	0
2	FAD	A	601	-	51,58,58	1.71	9 (17%)	60,89,89	2.53	18 (30%)
2	FAD	C	601	-	51,58,58	2.05	10 (19%)	60,89,89	2.48	19 (31%)
3	ORN	A	603	-	4,8,8	0.85	0	3,9,9	0.35	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
3	ORN	C	603	-	-	0/4/8/8	-
3	ORN	D	602	-	-	2/4/8/8	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ORN	D	603	-	-	0/4/8/8	-
2	FAD	B	601	-	-	2/30/50/50	0/6/6/6
3	ORN	B	603	-	-	0/4/8/8	-
2	FAD	D	601	-	-	8/30/50/50	0/6/6/6
3	ORN	A	602	-	-	0/4/8/8	-
3	ORN	B	602	-	-	3/4/8/8	-
3	ORN	C	602	-	-	3/4/8/8	-
2	FAD	A	601	-	-	2/30/50/50	0/6/6/6
2	FAD	C	601	-	-	3/30/50/50	0/6/6/6
3	ORN	A	603	-	-	0/4/8/8	-

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	FAD	C4X-C10	10.38	1.49	1.38
2	C	601	FAD	C4X-C10	9.73	1.48	1.38
2	B	601	FAD	C4X-C10	9.15	1.48	1.38
2	A	601	FAD	C4X-C10	5.95	1.44	1.38
2	D	601	FAD	C9A-C5X	4.29	1.51	1.42
2	C	601	FAD	C9A-N10	4.24	1.44	1.38
2	C	601	FAD	C9A-C5X	3.98	1.50	1.42
2	D	601	FAD	C4-C4X	3.84	1.48	1.41
2	C	601	FAD	C4-C4X	3.83	1.48	1.41
2	A	601	FAD	C10-N1	3.73	1.38	1.33
2	C	601	FAD	C8-C7	3.57	1.49	1.40
2	D	601	FAD	C8-C7	3.53	1.49	1.40
2	B	601	FAD	C10-N1	3.30	1.37	1.33
2	B	601	FAD	C4-C4X	3.26	1.47	1.41
2	D	601	FAD	C9A-N10	3.21	1.42	1.38
2	D	601	FAD	C5A-C4A	3.08	1.49	1.40
2	A	601	FAD	C2A-N3A	3.08	1.37	1.32
2	A	601	FAD	C2-N1	-2.99	1.32	1.38
2	B	601	FAD	C1'-N10	-2.96	1.45	1.48
2	A	601	FAD	C4-C4X	2.95	1.46	1.41
2	A	601	FAD	C2A-N1A	2.92	1.39	1.33
2	A	601	FAD	C5A-C4A	2.91	1.48	1.40
2	A	601	FAD	C1'-N10	-2.90	1.45	1.48
2	A	601	FAD	C9A-C5X	2.89	1.48	1.42
2	B	601	FAD	C4X-N5	2.88	1.37	1.33
2	B	601	FAD	C9A-N10	2.69	1.42	1.38
2	C	601	FAD	C5A-C4A	2.63	1.47	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	FAD	C9A-C5X	2.51	1.47	1.42
2	D	601	FAD	C2A-N1A	2.44	1.38	1.33
2	C	601	FAD	C1'-N10	-2.39	1.45	1.48
2	D	601	FAD	C2A-N3A	2.33	1.35	1.32
2	D	601	FAD	C1'-N10	-2.28	1.45	1.48
2	C	601	FAD	C4-N3	2.25	1.37	1.33
2	D	601	FAD	C10-N1	2.24	1.36	1.33
2	B	601	FAD	C8-C7	2.19	1.46	1.40
2	C	601	FAD	C2A-N3A	2.12	1.35	1.32
2	D	601	FAD	C4-N3	2.10	1.36	1.33
2	C	601	FAD	O2'-C2'	-2.09	1.38	1.43
2	B	601	FAD	O2'-C2'	-2.01	1.39	1.43
2	B	601	FAD	C5X-N5	2.00	1.38	1.35

All (78) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	FAD	C1'-N10-C9A	11.45	127.30	118.29
2	B	601	FAD	C4-N3-C2	9.31	123.00	115.14
2	A	601	FAD	C4-N3-C2	9.30	122.99	115.14
2	C	601	FAD	C4-N3-C2	9.23	122.93	115.14
2	A	601	FAD	C1'-N10-C9A	8.09	124.66	118.29
2	D	601	FAD	C4-N3-C2	7.97	121.87	115.14
2	A	601	FAD	C4-C4X-C10	-6.73	115.50	119.95
2	B	601	FAD	C4-C4X-C10	-6.63	115.56	119.95
2	C	601	FAD	C1'-N10-C9A	6.27	123.22	118.29
2	B	601	FAD	N3A-C2A-N1A	-5.85	119.53	128.68
2	D	601	FAD	C4-C4X-C10	-5.79	116.11	119.95
2	D	601	FAD	O4B-C1B-C2B	-5.68	98.62	106.93
2	C	601	FAD	C4-C4X-C10	-5.64	116.22	119.95
2	B	601	FAD	C1'-N10-C10	-5.17	113.78	118.41
2	D	601	FAD	C1'-N10-C9A	5.06	122.27	118.29
2	B	601	FAD	C10-C4X-N5	4.96	124.69	121.26
2	D	601	FAD	O4B-C4B-C5B	4.85	125.34	109.37
2	B	601	FAD	C5X-C9A-N10	4.84	121.22	117.72
2	C	601	FAD	C4X-C4-N3	-4.54	117.22	123.43
2	A	601	FAD	N3A-C2A-N1A	-4.50	121.65	128.68
2	A	601	FAD	C5'-C4'-C3'	-4.36	103.78	112.20
2	C	601	FAD	C4X-N5-C5X	3.97	120.73	116.77
2	C	601	FAD	C5X-C9A-N10	3.89	120.53	117.72
2	C	601	FAD	O4B-C1B-C2B	-3.87	101.27	106.93
2	A	601	FAD	C10-C4X-N5	3.85	123.92	121.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	FAD	N3A-C2A-N1A	-3.82	122.71	128.68
2	B	601	FAD	C5'-C4'-C3'	-3.76	104.95	112.20
2	D	601	FAD	C4X-N5-C5X	3.74	120.50	116.77
2	C	601	FAD	C10-C4X-N5	3.70	123.82	121.26
2	D	601	FAD	O2P-P-O1P	3.66	130.31	112.24
2	B	601	FAD	C4X-C4-N3	-3.52	118.62	123.43
2	A	601	FAD	C4X-C4-N3	-3.47	118.69	123.43
2	D	601	FAD	C4X-C4-N3	-3.44	118.72	123.43
2	B	601	FAD	C4X-C10-N10	-3.34	116.87	120.30
2	A	601	FAD	C8M-C8-C7	-3.30	113.98	120.74
2	D	601	FAD	O5B-C5B-C4B	3.29	120.32	108.99
2	A	601	FAD	C2A-N1A-C6A	3.11	124.07	118.75
2	D	601	FAD	C4A-C5A-N7A	-3.02	106.25	109.40
2	D	601	FAD	O2A-PA-O1A	3.02	127.17	112.24
2	D	601	FAD	N3A-C2A-N1A	-2.90	124.14	128.68
2	C	601	FAD	C4A-C5A-N7A	-2.90	106.38	109.40
2	C	601	FAD	O2A-PA-O1A	2.81	126.15	112.24
2	C	601	FAD	C9A-N10-C10	-2.81	118.23	121.91
2	D	601	FAD	C5'-C4'-C3'	-2.78	106.83	112.20
2	B	601	FAD	O4B-C1B-C2B	-2.78	102.86	106.93
2	C	601	FAD	O2P-P-O1P	2.68	125.49	112.24
2	D	601	FAD	C10-C4X-N5	2.68	123.11	121.26
2	C	601	FAD	C4X-C10-N10	-2.66	117.57	120.30
2	B	601	FAD	C9-C8-C7	2.57	124.24	119.91
2	A	601	FAD	O2P-P-O1P	2.56	124.90	112.24
2	B	601	FAD	O4'-C4'-C3'	2.51	115.20	109.10
2	A	601	FAD	O4'-C4'-C3'	2.46	115.08	109.10
2	A	601	FAD	N6A-C6A-N1A	2.45	123.66	118.57
2	D	601	FAD	O2B-C2B-C1B	2.44	119.85	110.85
2	C	601	FAD	C2A-N1A-C6A	2.43	122.91	118.75
2	B	601	FAD	C8M-C8-C7	-2.39	115.83	120.74
2	B	601	FAD	C1B-N9A-C4A	-2.37	122.48	126.64
2	A	601	FAD	O5B-PA-O1A	-2.33	99.95	109.07
2	D	601	FAD	C9A-N10-C10	-2.32	118.87	121.91
2	B	601	FAD	C5A-C6A-N6A	-2.32	116.83	120.35
2	A	601	FAD	C9-C8-C7	2.31	123.81	119.91
2	A	601	FAD	C1'-C2'-C3'	-2.31	103.32	109.79
2	B	601	FAD	C9A-N10-C10	-2.30	118.90	121.91
2	C	601	FAD	C8M-C8-C9	-2.29	114.87	120.34
2	A	601	FAD	C1'-N10-C10	-2.28	116.37	118.41
2	A	601	FAD	C9A-N10-C10	-2.26	118.94	121.91
2	A	601	FAD	C4X-C10-N10	-2.26	117.98	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	FAD	O2P-P-O1P	2.26	123.39	112.24
2	B	601	FAD	O2A-PA-O1A	2.25	123.39	112.24
2	C	601	FAD	C6-C5X-C9A	2.19	121.92	119.05
2	C	601	FAD	C1'-C2'-C3'	-2.19	103.67	109.79
2	D	601	FAD	C5X-C9A-N10	2.17	119.29	117.72
2	D	601	FAD	C4X-C10-N10	-2.15	118.09	120.30
2	B	601	FAD	N6A-C6A-N1A	2.13	122.98	118.57
2	B	601	FAD	C2A-N1A-C6A	2.11	122.37	118.75
2	C	601	FAD	C9A-C5X-N5	-2.08	119.10	122.36
2	D	601	FAD	O5'-P-O1P	-2.02	101.16	109.07
3	C	603	ORN	CG-CB-CA	2.00	119.66	113.35

There are no chirality outliers.

All (23) torsion outliers are listed below:

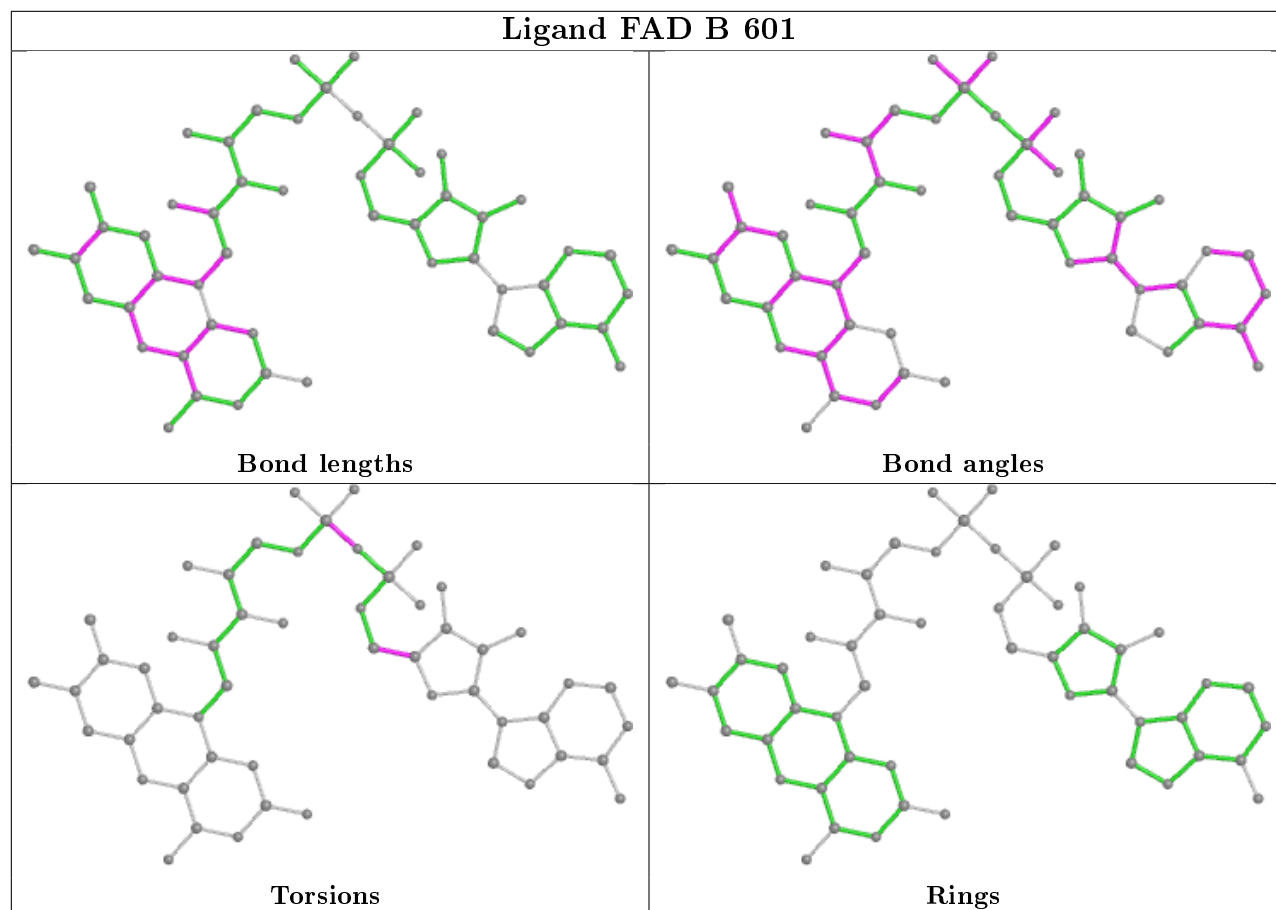
Mol	Chain	Res	Type	Atoms
3	D	602	ORN	N-CA-CB-CG
3	D	602	ORN	C-CA-CB-CG
2	D	601	FAD	C5B-O5B-PA-O1A
2	D	601	FAD	C5B-O5B-PA-O2A
2	D	601	FAD	C5B-O5B-PA-O3P
3	B	602	ORN	C-CA-CB-CG
2	B	601	FAD	PA-O3P-P-O5'
2	A	601	FAD	PA-O3P-P-O5'
2	C	601	FAD	C5B-O5B-PA-O1A
2	D	601	FAD	C2'-C3'-C4'-O4'
2	C	601	FAD	PA-O3P-P-O5'
3	C	602	ORN	CA-CB-CG-CD
3	B	602	ORN	N-CA-CB-CG
3	C	602	ORN	NE-CD-CG-CB
2	A	601	FAD	O4B-C4B-C5B-O5B
2	D	601	FAD	O3'-C3'-C4'-O4'
2	D	601	FAD	C3B-C4B-C5B-O5B
3	B	602	ORN	NE-CD-CG-CB
2	D	601	FAD	C2'-C3'-C4'-C5'
2	C	601	FAD	O4B-C4B-C5B-O5B
2	B	601	FAD	O4B-C4B-C5B-O5B
3	C	602	ORN	N-CA-CB-CG
2	D	601	FAD	C4B-C5B-O5B-PA

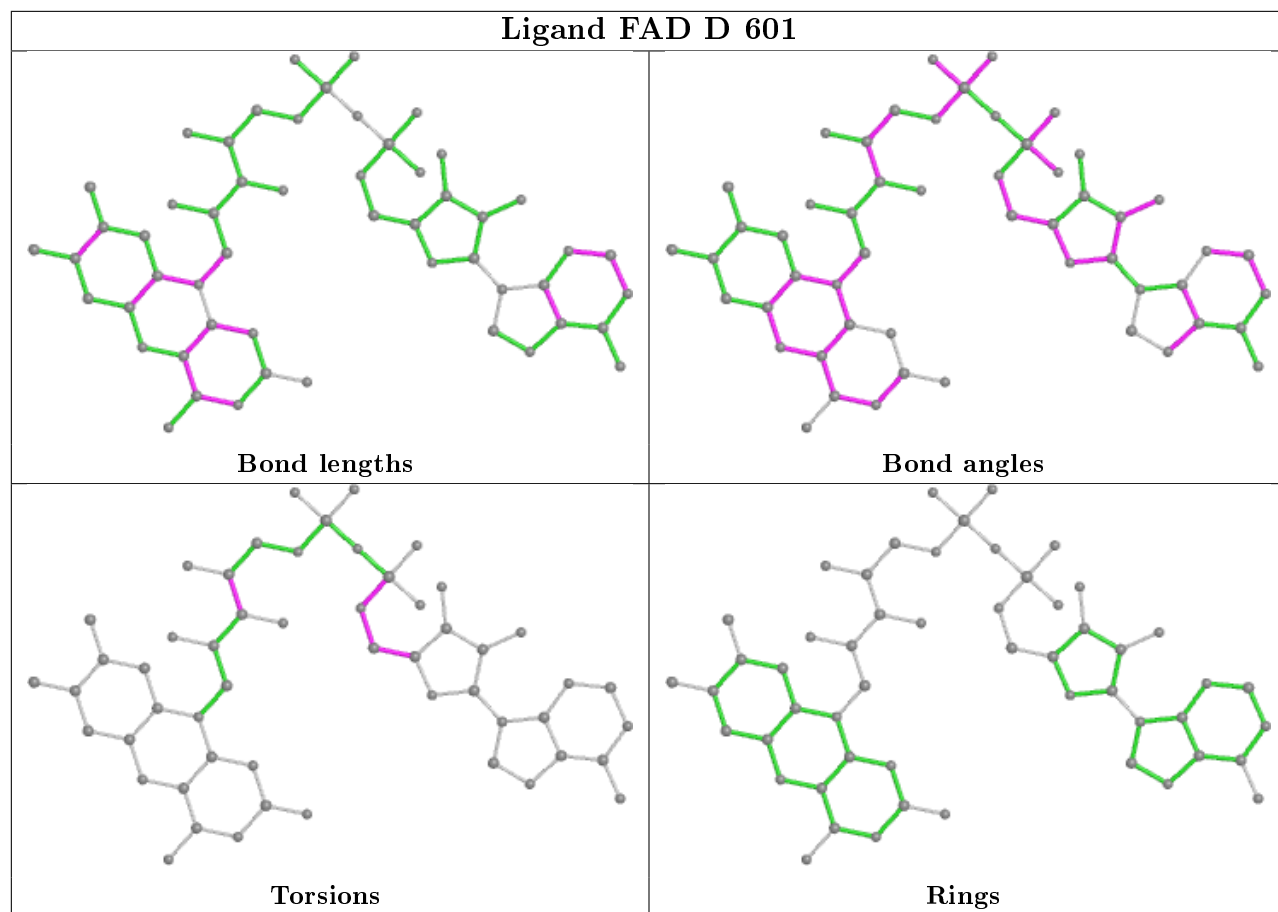
There are no ring outliers.

8 monomers are involved in 22 short contacts:

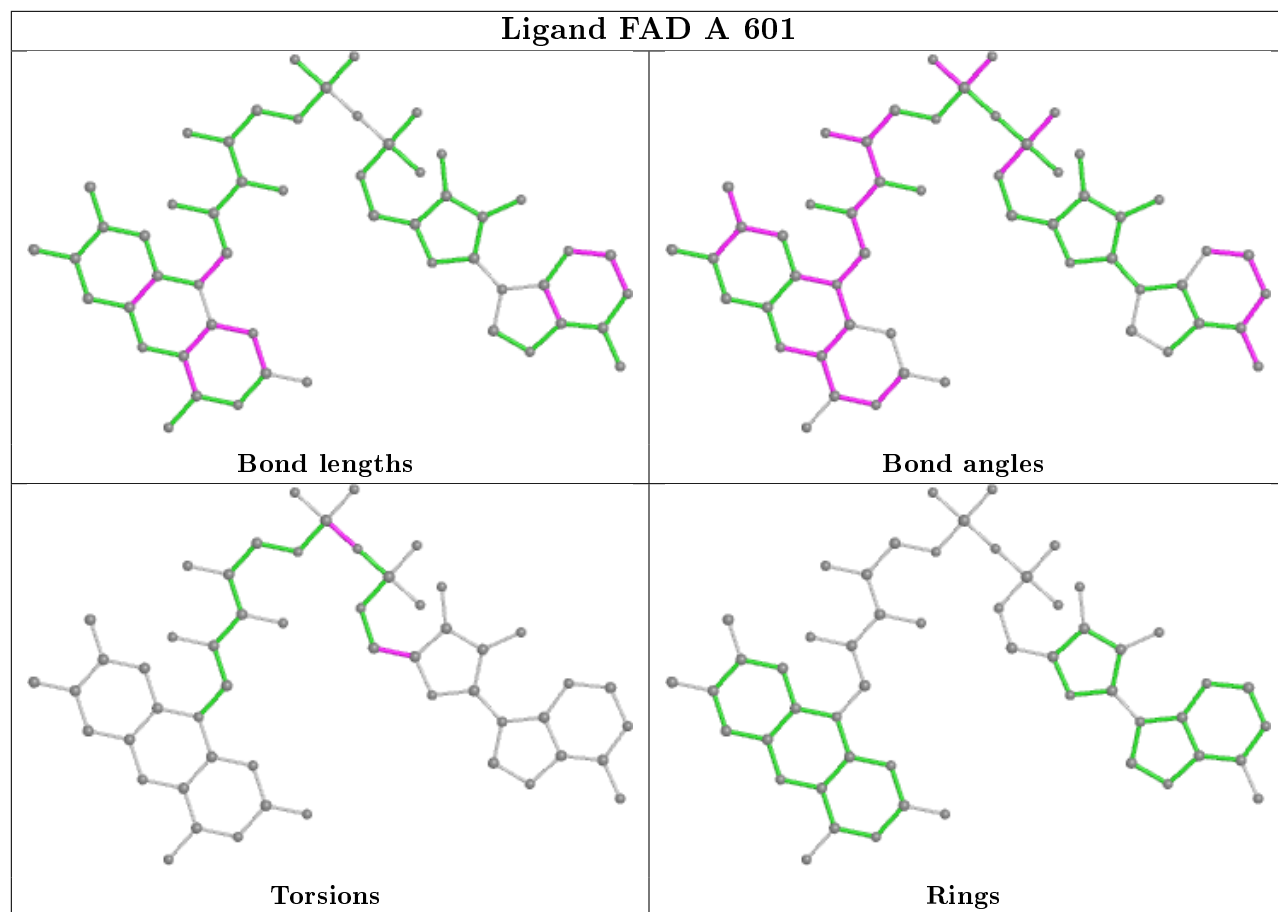
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	603	ORN	1	0
2	B	601	FAD	4	0
3	B	603	ORN	2	0
2	D	601	FAD	4	0
3	C	602	ORN	2	0
2	A	601	FAD	4	0
2	C	601	FAD	4	0
3	A	603	ORN	1	0

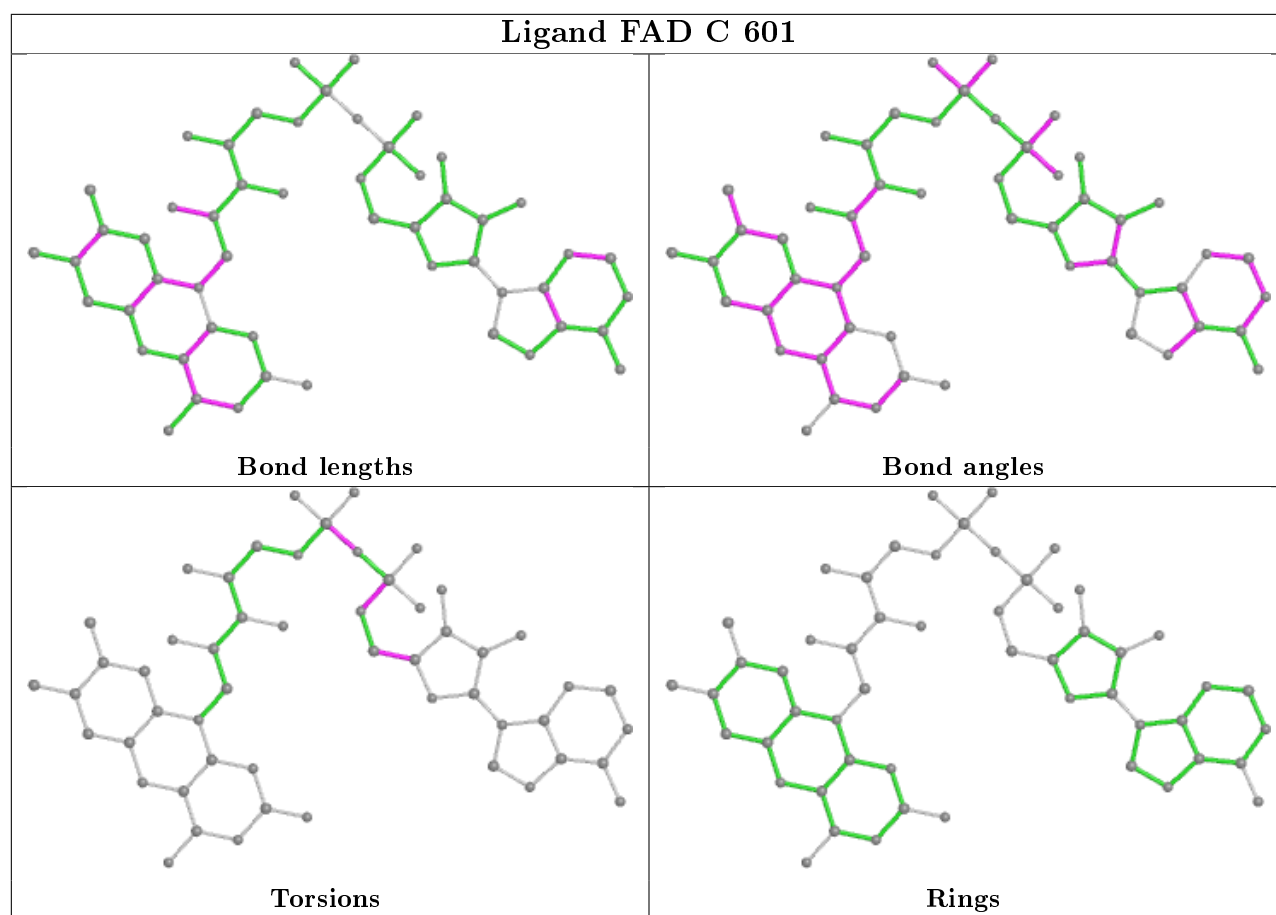
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 FAD A 601





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	549/580 (94%)	-0.25	5 (0%) 84 83	17, 29, 46, 88	0
1	B	549/580 (94%)	-0.22	5 (0%) 84 83	17, 30, 48, 77	0
1	C	549/580 (94%)	0.13	26 (4%) 31 30	25, 39, 59, 102	0
1	D	549/580 (94%)	0.22	34 (6%) 20 19	25, 40, 60, 96	0
All	All	2196/2320 (94%)	-0.03	70 (3%) 47 46	17, 35, 56, 102	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	278	GLU	5.0
1	D	279	ARG	4.5
1	C	279	ARG	4.3
1	D	280	CYS	4.2
1	C	278	GLU	4.0
1	A	12	LYS	4.0
1	D	39	ILE	3.9
1	D	12	LYS	3.8
1	C	12	LYS	3.6
1	A	278	GLU	3.5
1	D	382	THR	3.5
1	D	521	VAL	3.4
1	C	426	LEU	3.3
1	D	41	ALA	3.3
1	D	14	PRO	3.2
1	C	406	ASP	3.1
1	D	381	GLU	3.0
1	B	380	PRO	2.9
1	C	381	GLU	2.9
1	C	521	VAL	2.7
1	C	14	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	287	THR	2.7
1	C	427	PRO	2.7
1	D	307	ALA	2.7
1	D	42	ALA	2.6
1	D	498	VAL	2.6
1	C	405	GLY	2.6
1	D	406	ASP	2.6
1	C	520	ALA	2.6
1	B	407	ASP	2.6
1	C	380	PRO	2.5
1	D	229	GLY	2.5
1	D	426	LEU	2.4
1	D	13	LYS	2.4
1	C	517	VAL	2.4
1	D	281	ALA	2.4
1	C	382	THR	2.3
1	D	520	ALA	2.3
1	D	286	GLY	2.3
1	C	42	ALA	2.3
1	C	511	SER	2.3
1	A	279	ARG	2.3
1	B	12	LYS	2.3
1	C	55	ILE	2.3
1	D	160	THR	2.3
1	A	406	ASP	2.2
1	D	513	THR	2.2
1	C	13	LYS	2.2
1	D	541	ALA	2.2
1	A	427	PRO	2.2
1	C	280	CYS	2.2
1	D	285	GLU	2.2
1	D	540	HIS	2.1
1	C	506	ALA	2.1
1	D	516	TRP	2.1
1	B	278	GLU	2.1
1	C	513	THR	2.1
1	D	511	SER	2.1
1	C	525	LEU	2.1
1	D	552	GLU	2.1
1	D	358	THR	2.1
1	C	285	GLU	2.0
1	C	507	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	542	ASP	2.0
1	B	427	PRO	2.0
1	D	44	HIS	2.0
1	D	50	ILE	2.0
1	D	55	ILE	2.0
1	C	408	LYS	2.0
1	D	435	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](#)

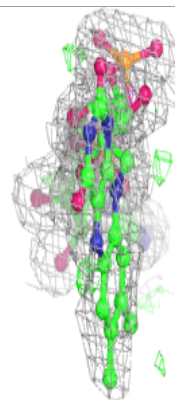
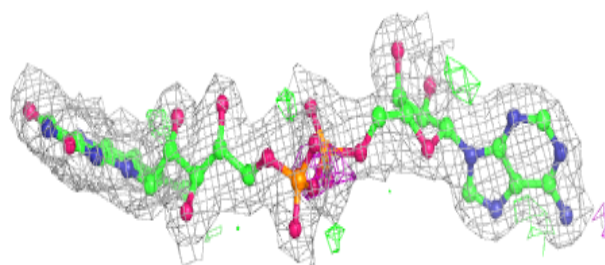
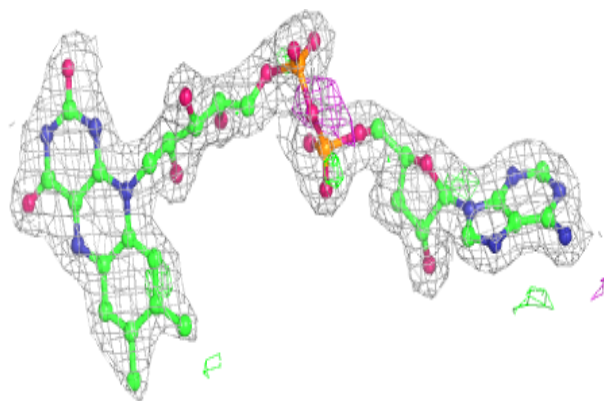
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	ORN	D	602	9/9	0.87	0.31	45,51,57,58	0
3	ORN	B	602	9/9	0.90	0.27	40,48,51,55	0
3	ORN	C	602	9/9	0.91	0.28	42,49,54,55	0
3	ORN	C	603	9/9	0.91	0.14	27,30,33,35	0
3	ORN	A	602	9/9	0.92	0.29	39,45,58,68	0
3	ORN	B	603	9/9	0.93	0.10	32,36,38,39	0
3	ORN	D	603	9/9	0.94	0.10	28,29,32,33	0
2	FAD	D	601	53/53	0.94	0.13	27,33,45,51	0
3	ORN	A	603	9/9	0.94	0.10	31,34,39,39	0
2	FAD	C	601	53/53	0.95	0.14	27,33,45,54	0
2	FAD	A	601	53/53	0.96	0.14	18,23,27,31	0
2	FAD	B	601	53/53	0.97	0.14	16,21,27,33	0

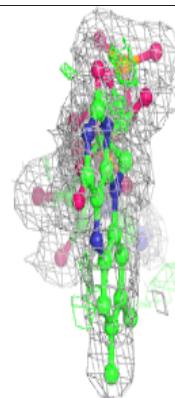
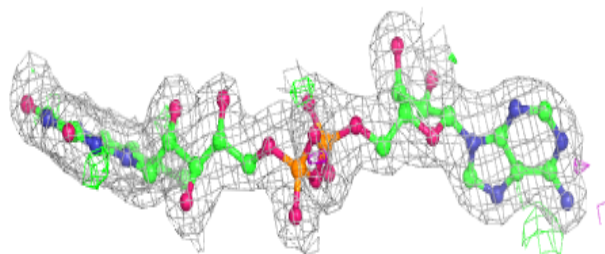
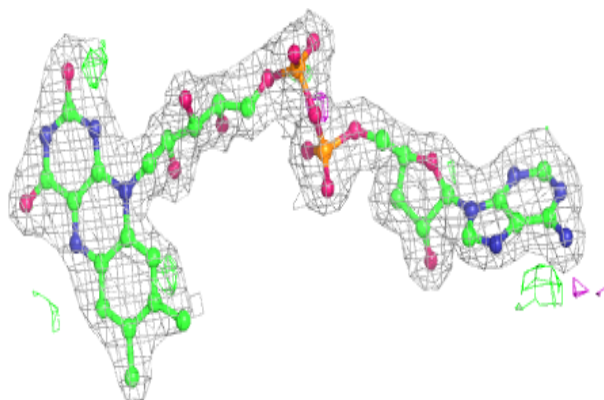
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD D 601:

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

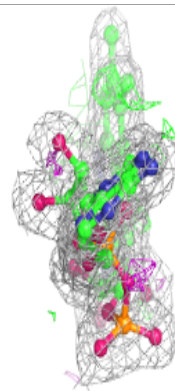
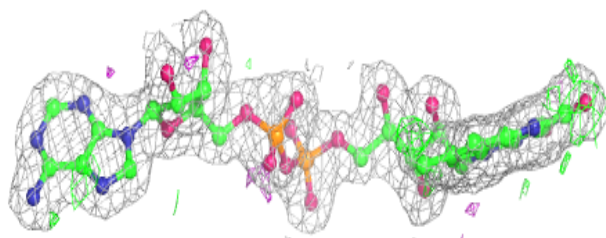
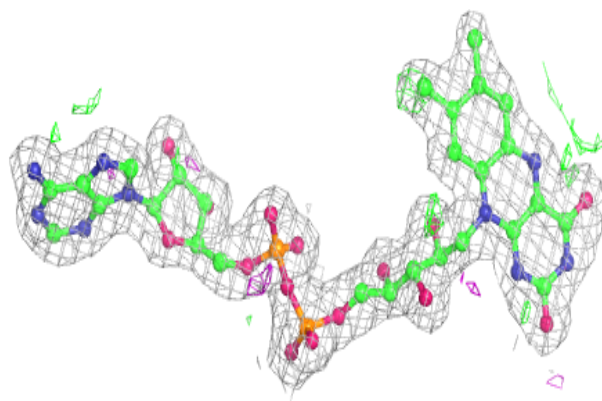
**Electron density around FAD C 601:**

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

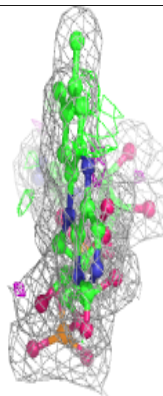
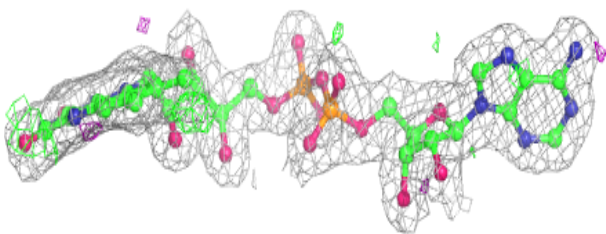
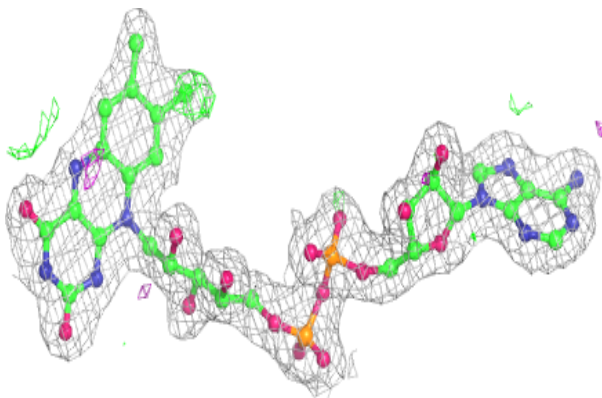


Electron density around FAD A 601:

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

**Electron density around FAD B 601:**

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



6.5 Other polymers ⓘ

There are no such residues in this entry.