



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:17 AM UTC

PDB ID : 1T9D / pdb_00001t9d
Title : Crystal Structure Of Yeast Acetohydroxyacid Synthase In Complex With A Sulfonylurea Herbicide, Metsulfuron methyl
Authors : McCourt, J.A.; Pang, S.S.; Guddat, L.W.; Duggleby, R.G.
Deposited on : 2004-05-16
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

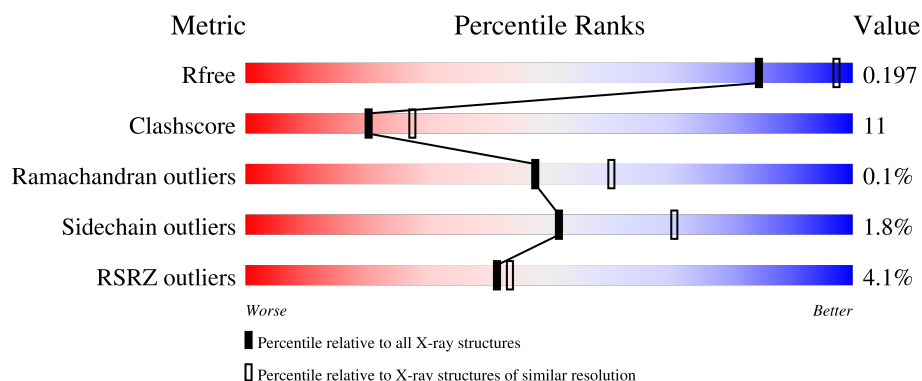
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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

Mol	Chain	Length	Quality of chain
1	A	677	0.1% 73% 14% • 12%
1	B	677	2% 66% 19% • 14%
1	C	677	11% 62% 22% • 15%
1	D	677	0.1% 73% 14% • 12%

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
4	1MM	A	695	-	X	-	-
4	1MM	B	1695	-	X	-	-
4	1MM	C	2695	-	X	-	-
4	1MM	C	3695	-	X	-	-
6	PYD	A	703	-	X	-	-
6	PYD	B	1703	-	X	-	-
6	PYD	C	2703	-	X	-	-
6	PYD	D	3703	-	X	-	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 20377 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 Acetolactate synthase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	596	Total	C	N	O	S	0	6	0
			4548	2879	783	865	21			
1	B	582	Total	C	N	O	S	0	3	0
			4392	2788	758	827	19			
1	C	575	Total	C	N	O	S	0	2	0
			4323	2742	746	816	19			
1	D	596	Total	C	N	O	S	0	5	0
			4528	2867	781	860	20			

There are 188 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	MET	-	cloning artifact	UNP P07342
A	12	HIS	-	cloning artifact	UNP P07342
A	13	HIS	-	cloning artifact	UNP P07342
A	14	HIS	-	cloning artifact	UNP P07342
A	15	HIS	-	cloning artifact	UNP P07342
A	16	HIS	-	cloning artifact	UNP P07342
A	17	HIS	-	cloning artifact	UNP P07342
A	18	SER	-	cloning artifact	UNP P07342
A	19	SER	-	cloning artifact	UNP P07342
A	20	GLY	-	cloning artifact	UNP P07342
A	21	LEU	-	cloning artifact	UNP P07342
A	22	VAL	-	cloning artifact	UNP P07342
A	23	PRO	-	cloning artifact	UNP P07342
A	24	ARG	-	cloning artifact	UNP P07342
A	25	GLY	-	cloning artifact	UNP P07342
A	26	SER	-	cloning artifact	UNP P07342
A	27	GLY	-	cloning artifact	UNP P07342
A	28	MET	-	cloning artifact	UNP P07342
A	29	LYS	-	cloning artifact	UNP P07342
A	30	GLU	-	cloning artifact	UNP P07342
A	31	THR	-	cloning artifact	UNP P07342

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Chain	Residue	Modelled	Actual	Comment	Reference
A	32	ALA	-	cloning artifact	UNP P07342
A	33	ALA	-	cloning artifact	UNP P07342
A	34	ALA	-	cloning artifact	UNP P07342
A	35	LYS	-	cloning artifact	UNP P07342
A	36	PHE	-	cloning artifact	UNP P07342
A	37	GLU	-	cloning artifact	UNP P07342
A	38	ARG	-	cloning artifact	UNP P07342
A	39	GLN	-	cloning artifact	UNP P07342
A	40	HIS	-	cloning artifact	UNP P07342
A	41	MET	-	cloning artifact	UNP P07342
A	42	ASP	-	cloning artifact	UNP P07342
A	43	SER	-	cloning artifact	UNP P07342
A	44	PRO	-	cloning artifact	UNP P07342
A	45	ASP	-	cloning artifact	UNP P07342
A	46	LEU	-	cloning artifact	UNP P07342
A	47	GLY	-	cloning artifact	UNP P07342
A	48	THR	-	cloning artifact	UNP P07342
A	49	ASP	-	cloning artifact	UNP P07342
A	50	ASP	-	cloning artifact	UNP P07342
A	51	ASP	-	cloning artifact	UNP P07342
A	52	ASP	-	cloning artifact	UNP P07342
A	53	LYS	-	cloning artifact	UNP P07342
A	54	ALA	-	cloning artifact	UNP P07342
A	55	MET	-	cloning artifact	UNP P07342
A	56	GLY	-	cloning artifact	UNP P07342
A	57	SER	-	cloning artifact	UNP P07342
B	11	MET	-	cloning artifact	UNP P07342
B	12	HIS	-	cloning artifact	UNP P07342
B	13	HIS	-	cloning artifact	UNP P07342
B	14	HIS	-	cloning artifact	UNP P07342
B	15	HIS	-	cloning artifact	UNP P07342
B	16	HIS	-	cloning artifact	UNP P07342
B	17	HIS	-	cloning artifact	UNP P07342
B	18	SER	-	cloning artifact	UNP P07342
B	19	SER	-	cloning artifact	UNP P07342
B	20	GLY	-	cloning artifact	UNP P07342
B	21	LEU	-	cloning artifact	UNP P07342
B	22	VAL	-	cloning artifact	UNP P07342
B	23	PRO	-	cloning artifact	UNP P07342
B	24	ARG	-	cloning artifact	UNP P07342
B	25	GLY	-	cloning artifact	UNP P07342
B	26	SER	-	cloning artifact	UNP P07342

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Chain	Residue	Modelled	Actual	Comment	Reference
B	27	GLY	-	cloning artifact	UNP P07342
B	28	MET	-	cloning artifact	UNP P07342
B	29	LYS	-	cloning artifact	UNP P07342
B	30	GLU	-	cloning artifact	UNP P07342
B	31	THR	-	cloning artifact	UNP P07342
B	32	ALA	-	cloning artifact	UNP P07342
B	33	ALA	-	cloning artifact	UNP P07342
B	34	ALA	-	cloning artifact	UNP P07342
B	35	LYS	-	cloning artifact	UNP P07342
B	36	PHE	-	cloning artifact	UNP P07342
B	37	GLU	-	cloning artifact	UNP P07342
B	38	ARG	-	cloning artifact	UNP P07342
B	39	GLN	-	cloning artifact	UNP P07342
B	40	HIS	-	cloning artifact	UNP P07342
B	41	MET	-	cloning artifact	UNP P07342
B	42	ASP	-	cloning artifact	UNP P07342
B	43	SER	-	cloning artifact	UNP P07342
B	44	PRO	-	cloning artifact	UNP P07342
B	45	ASP	-	cloning artifact	UNP P07342
B	46	LEU	-	cloning artifact	UNP P07342
B	47	GLY	-	cloning artifact	UNP P07342
B	48	THR	-	cloning artifact	UNP P07342
B	49	ASP	-	cloning artifact	UNP P07342
B	50	ASP	-	cloning artifact	UNP P07342
B	51	ASP	-	cloning artifact	UNP P07342
B	52	ASP	-	cloning artifact	UNP P07342
B	53	LYS	-	cloning artifact	UNP P07342
B	54	ALA	-	cloning artifact	UNP P07342
B	55	MET	-	cloning artifact	UNP P07342
B	56	GLY	-	cloning artifact	UNP P07342
B	57	SER	-	cloning artifact	UNP P07342
C	11	MET	-	cloning artifact	UNP P07342
C	12	HIS	-	cloning artifact	UNP P07342
C	13	HIS	-	cloning artifact	UNP P07342
C	14	HIS	-	cloning artifact	UNP P07342
C	15	HIS	-	cloning artifact	UNP P07342
C	16	HIS	-	cloning artifact	UNP P07342
C	17	HIS	-	cloning artifact	UNP P07342
C	18	SER	-	cloning artifact	UNP P07342
C	19	SER	-	cloning artifact	UNP P07342
C	20	GLY	-	cloning artifact	UNP P07342
C	21	LEU	-	cloning artifact	UNP P07342

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Chain	Residue	Modelled	Actual	Comment	Reference
C	22	VAL	-	cloning artifact	UNP P07342
C	23	PRO	-	cloning artifact	UNP P07342
C	24	ARG	-	cloning artifact	UNP P07342
C	25	GLY	-	cloning artifact	UNP P07342
C	26	SER	-	cloning artifact	UNP P07342
C	27	GLY	-	cloning artifact	UNP P07342
C	28	MET	-	cloning artifact	UNP P07342
C	29	LYS	-	cloning artifact	UNP P07342
C	30	GLU	-	cloning artifact	UNP P07342
C	31	THR	-	cloning artifact	UNP P07342
C	32	ALA	-	cloning artifact	UNP P07342
C	33	ALA	-	cloning artifact	UNP P07342
C	34	ALA	-	cloning artifact	UNP P07342
C	35	LYS	-	cloning artifact	UNP P07342
C	36	PHE	-	cloning artifact	UNP P07342
C	37	GLU	-	cloning artifact	UNP P07342
C	38	ARG	-	cloning artifact	UNP P07342
C	39	GLN	-	cloning artifact	UNP P07342
C	40	HIS	-	cloning artifact	UNP P07342
C	41	MET	-	cloning artifact	UNP P07342
C	42	ASP	-	cloning artifact	UNP P07342
C	43	SER	-	cloning artifact	UNP P07342
C	44	PRO	-	cloning artifact	UNP P07342
C	45	ASP	-	cloning artifact	UNP P07342
C	46	LEU	-	cloning artifact	UNP P07342
C	47	GLY	-	cloning artifact	UNP P07342
C	48	THR	-	cloning artifact	UNP P07342
C	49	ASP	-	cloning artifact	UNP P07342
C	50	ASP	-	cloning artifact	UNP P07342
C	51	ASP	-	cloning artifact	UNP P07342
C	52	ASP	-	cloning artifact	UNP P07342
C	53	LYS	-	cloning artifact	UNP P07342
C	54	ALA	-	cloning artifact	UNP P07342
C	55	MET	-	cloning artifact	UNP P07342
C	56	GLY	-	cloning artifact	UNP P07342
C	57	SER	-	cloning artifact	UNP P07342
D	11	MET	-	cloning artifact	UNP P07342
D	12	HIS	-	cloning artifact	UNP P07342
D	13	HIS	-	cloning artifact	UNP P07342
D	14	HIS	-	cloning artifact	UNP P07342
D	15	HIS	-	cloning artifact	UNP P07342
D	16	HIS	-	cloning artifact	UNP P07342

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Chain	Residue	Modelled	Actual	Comment	Reference
D	17	HIS	-	cloning artifact	UNP P07342
D	18	SER	-	cloning artifact	UNP P07342
D	19	SER	-	cloning artifact	UNP P07342
D	20	GLY	-	cloning artifact	UNP P07342
D	21	LEU	-	cloning artifact	UNP P07342
D	22	VAL	-	cloning artifact	UNP P07342
D	23	PRO	-	cloning artifact	UNP P07342
D	24	ARG	-	cloning artifact	UNP P07342
D	25	GLY	-	cloning artifact	UNP P07342
D	26	SER	-	cloning artifact	UNP P07342
D	27	GLY	-	cloning artifact	UNP P07342
D	28	MET	-	cloning artifact	UNP P07342
D	29	LYS	-	cloning artifact	UNP P07342
D	30	GLU	-	cloning artifact	UNP P07342
D	31	THR	-	cloning artifact	UNP P07342
D	32	ALA	-	cloning artifact	UNP P07342
D	33	ALA	-	cloning artifact	UNP P07342
D	34	ALA	-	cloning artifact	UNP P07342
D	35	LYS	-	cloning artifact	UNP P07342
D	36	PHE	-	cloning artifact	UNP P07342
D	37	GLU	-	cloning artifact	UNP P07342
D	38	ARG	-	cloning artifact	UNP P07342
D	39	GLN	-	cloning artifact	UNP P07342
D	40	HIS	-	cloning artifact	UNP P07342
D	41	MET	-	cloning artifact	UNP P07342
D	42	ASP	-	cloning artifact	UNP P07342
D	43	SER	-	cloning artifact	UNP P07342
D	44	PRO	-	cloning artifact	UNP P07342
D	45	ASP	-	cloning artifact	UNP P07342
D	46	LEU	-	cloning artifact	UNP P07342
D	47	GLY	-	cloning artifact	UNP P07342
D	48	THR	-	cloning artifact	UNP P07342
D	49	ASP	-	cloning artifact	UNP P07342
D	50	ASP	-	cloning artifact	UNP P07342
D	51	ASP	-	cloning artifact	UNP P07342
D	52	ASP	-	cloning artifact	UNP P07342
D	53	LYS	-	cloning artifact	UNP P07342
D	54	ALA	-	cloning artifact	UNP P07342
D	55	MET	-	cloning artifact	UNP P07342
D	56	GLY	-	cloning artifact	UNP P07342
D	57	SER	-	cloning artifact	UNP P07342

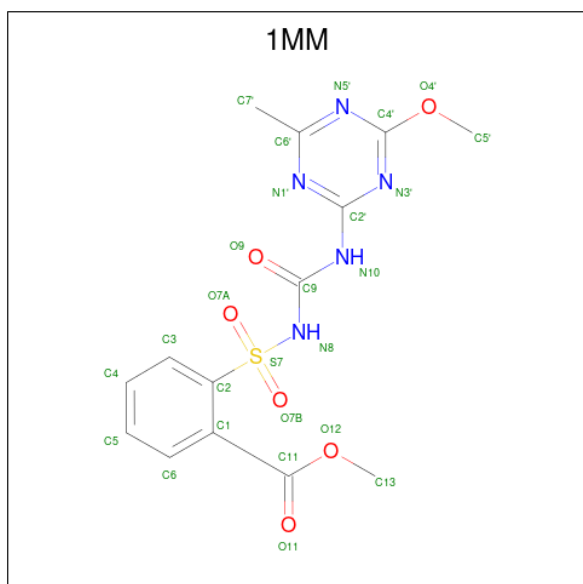
- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

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

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is METHYL 2-[(4-METHOXY-6-METHYL-1,3,5-TRIAZIN-2-YL)AMINO]CARBONYL]AMINO)SULFONYLBENZOATE (CCD ID: 1MM) (formula: C₁₄H₁₅N₅O₆S).



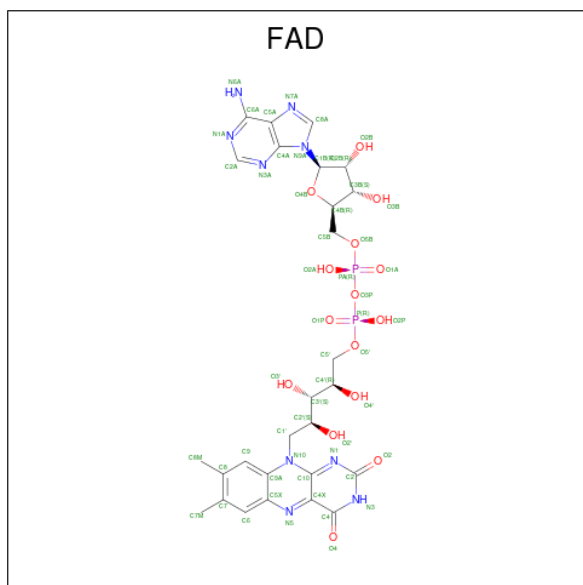
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O S 26 14 5 6 1	0	0

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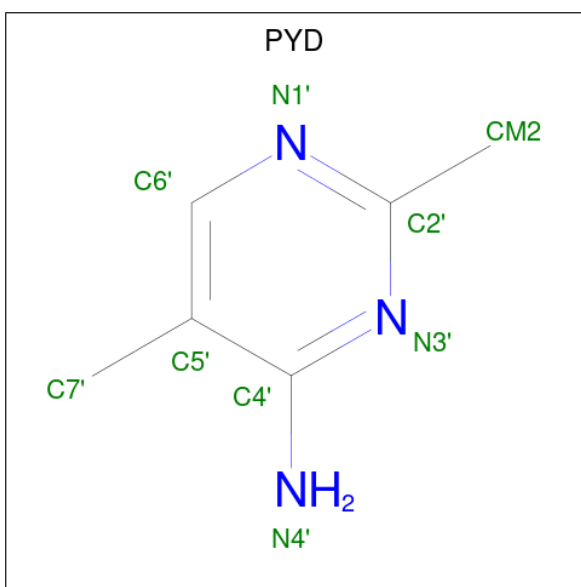
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	S	0	0
			26	14	5	6	1		
4	C	1	Total	C	N	O	S	0	0
			26	14	5	6	1		
4	C	1	Total	C	N	O	S	0	0
			26	14	5	6	1		

- Molecule 5 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



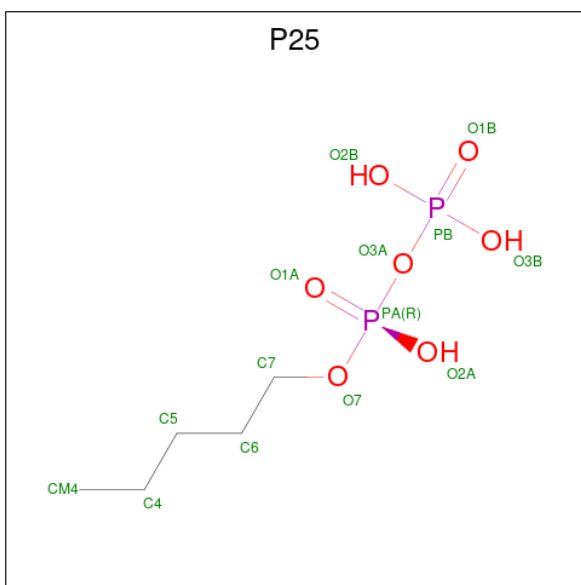
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
5	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
5	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
5	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 6 is 2,5-DIMETHYL-PYRIMIDIN-4-YLAMINE (CCD ID: PYD) (formula: $C_6H_9N_3$).



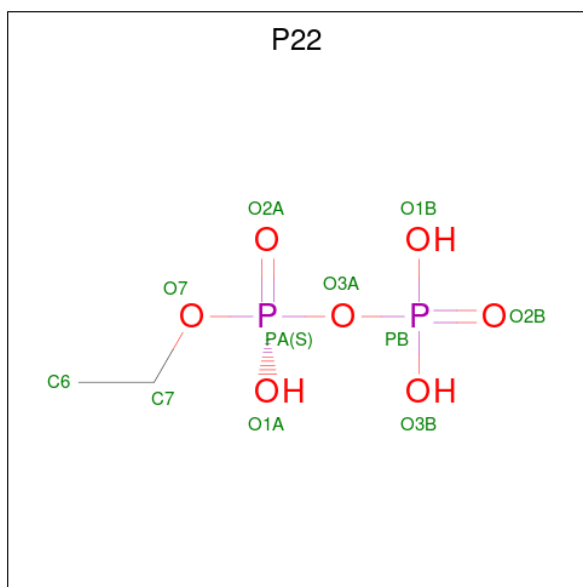
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	N	0	0
			9	6	3		
6	B	1	Total	C	N	0	0
			9	6	3		
6	C	1	Total	C	N	0	0
			9	6	3		
6	D	1	Total	C	N	0	0
			9	6	3		

- Molecule 7 is PENTYL TRIHYDROGEN DIPHOSPHATE (CCD ID: P25) (formula: C₅H₁₄O₇P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	O	P	0	0
			14	5	7	2		
7	B	1	Total	C	O	P	0	0
			14	5	7	2		

- Molecule 8 is ETHYL DIHYDROGEN DIPHOSPHATE (CCD ID: P22) (formula: $C_2H_8O_7P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	C	1	Total	C	O	P	0	0
			11	2	7	2		
8	D	1	Total	C	O	P	0	0
			11	2	7	2		

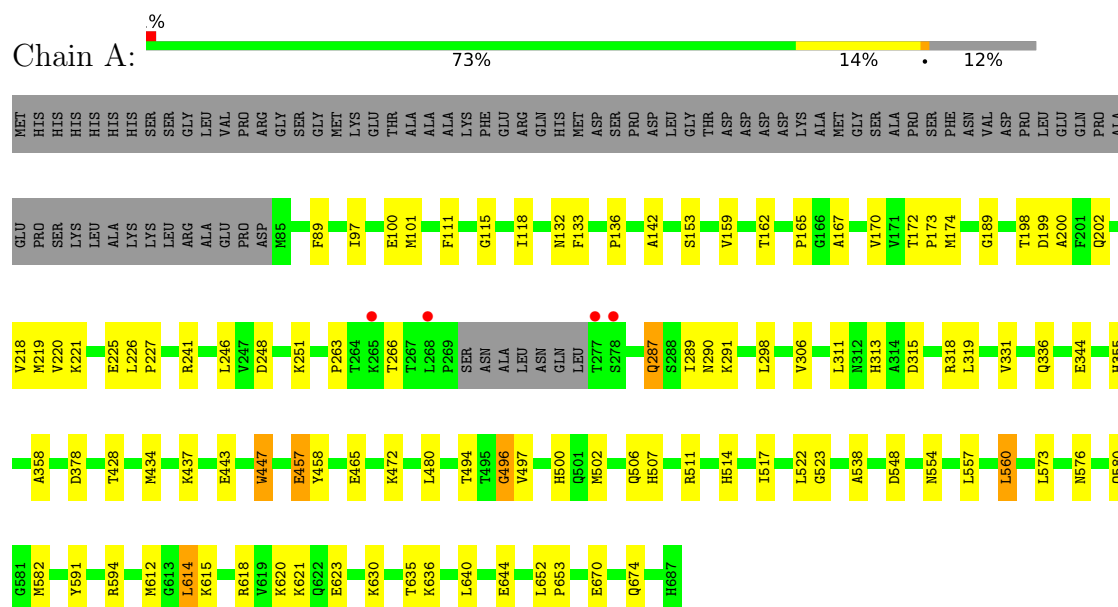
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	671	Total	O	1	0
			671	671		
9	B	468	Total	O	0	0
			468	468		
9	C	400	Total	O	0	0
			400	400		
9	D	637	Total	O	0	0
			637	637		

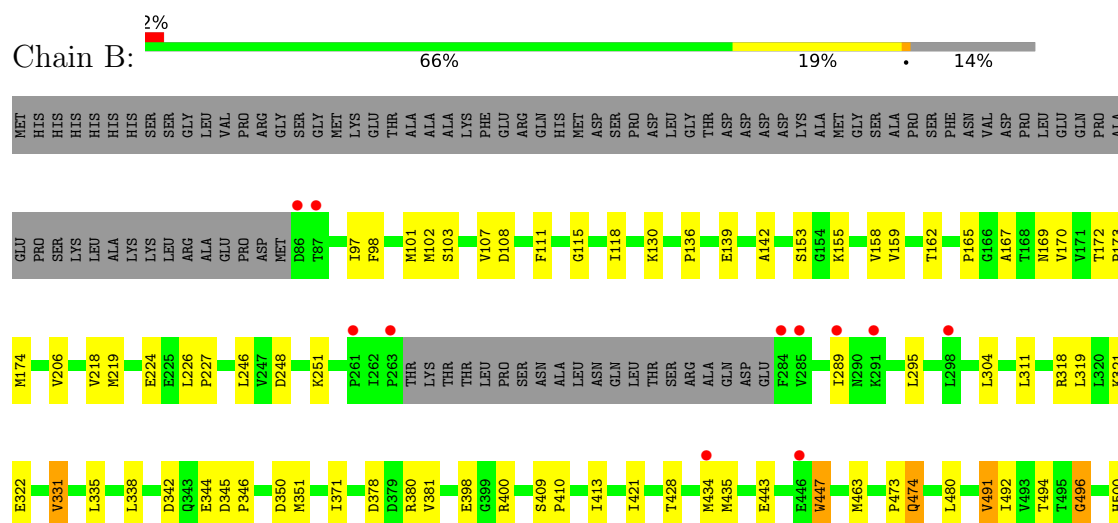
3 Residue-property plots

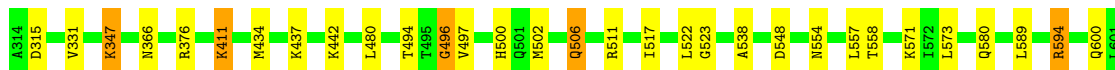
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Acetolactate synthase, mitochondrial



• Molecule 1: Acetolactate synthase, mitochondrial







4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	218.35Å 218.35Å 361.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 50.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	80.1 (50.00-2.30) 80.0 (50.00-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.73 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.164 , 0.195 0.167 , 0.197	Depositor DCC
R_{free} test set	15994 reflections (9.96%)	wwPDB-VP
Wilson B-factor (Å ²)	28.5	Xtriage
Anisotropy	0.342	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 64.0	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	0.96	EDS
Total number of atoms	20377	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, P25, 1MM, P22, PYD, FAD, K

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	0.38	0/4666	0.89	9/6332 (0.1%)
1	B	0.35	0/4497	0.85	4/6111 (0.1%)
1	C	0.32	0/4423	0.84	3/6014 (0.0%)
1	D	0.37	0/4642	0.89	7/6301 (0.1%)
All	All	0.36	0/18228	0.87	23/24758 (0.1%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	496	GLY	N-CA-C	-7.20	102.30	112.60
1	A	331	VAL	N-CA-C	7.18	118.24	107.75
1	D	496	GLY	N-CA-C	-7.11	102.44	112.60
1	D	306	VAL	N-CA-C	6.71	117.96	108.36
1	D	331	VAL	N-CA-C	6.61	117.37	108.11
1	A	580	GLN	N-CA-C	-6.29	100.18	109.62
1	D	506	GLN	N-CA-C	5.92	117.54	111.14
1	A	507	HIS	N-CA-C	5.82	120.22	113.18
1	B	491	VAL	N-CA-C	5.72	116.11	108.11
1	A	636	LYS	N-CA-C	-5.64	100.49	109.96
1	B	589	LEU	N-CA-C	5.61	117.47	111.36
1	C	189	GLY	N-CA-C	-5.57	106.21	112.33
1	A	241	ARG	N-CA-C	-5.56	101.51	109.24
1	B	331	VAL	N-CA-C	5.51	115.83	108.11
1	B	496	GLY	N-CA-C	-5.37	103.43	112.77
1	C	496	GLY	N-CA-C	-5.29	103.56	112.77
1	C	154	GLY	N-CA-C	-5.18	108.48	115.21
1	D	376	ARG	N-CA-C	5.16	119.26	112.92
1	A	189	GLY	N-CA-C	-5.14	106.68	112.33
1	A	306	VAL	N-CA-C	5.13	116.35	108.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	164	GLY	CA-C-N	5.09	124.88	119.28
1	D	164	GLY	C-N-CA	5.09	124.88	119.28
1	A	576	ASN	N-CA-C	5.05	116.50	108.67

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	4548	0	4521	83	0
1	B	4392	0	4352	114	0
1	C	4323	0	4263	143	0
1	D	4528	0	4494	85	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	26	0	15	1	0
4	B	26	0	15	4	0
4	C	52	0	30	2	0
5	A	53	0	31	0	0
5	B	53	0	31	1	0
5	C	53	0	31	0	0
5	D	53	0	31	0	0
6	A	9	0	9	1	0
6	B	9	0	9	0	0
6	C	9	0	9	0	0
6	D	9	0	9	0	0
7	A	14	0	11	2	0
7	B	14	0	11	1	0
8	C	11	0	5	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	D	11	0	5	1	0
9	A	671	0	0	9	0
9	B	468	0	0	13	0
9	C	400	0	0	19	0
9	D	637	0	0	12	0
All	All	20377	0	17882	409	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (409) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:594[A]:ARG:HG3	1:D:594[A]:ARG:HH11	1.16	1.04
1:C:114:PRO:HG2	1:D:580:GLN:HE22	1.25	1.00
1:B:335:LEU:HA	1:B:351:MET:HE2	1.49	0.94
1:C:102:MET:HE1	1:C:158:VAL:HG11	1.49	0.93
1:A:313:HIS:HD2	1:A:315:ASP:H	1.19	0.91
1:C:335:LEU:HA	1:C:351:MET:HE2	1.50	0.90
1:B:102:MET:HE1	1:B:158:VAL:HG11	1.55	0.89
1:C:619:VAL:HG22	1:C:628:LYS:HG3	1.54	0.89
1:B:580:GLN:HG3	7:B:698:P25:H72	1.54	0.87
1:D:313:HIS:HD2	1:D:315:ASP:H	1.21	0.86
1:B:619:VAL:HG22	1:B:628:LYS:HG3	1.60	0.82
1:B:103:SER:HB2	1:B:130:LYS:HD3	1.59	0.82
1:D:155:LYS:HE2	9:D:5257:HOH:O	1.78	0.82
1:C:114:PRO:HG2	1:D:580:GLN:NE2	1.95	0.81
1:D:101:MET:HE2	1:D:101:MET:HA	1.61	0.81
1:C:580:GLN:HE22	1:D:114:PRO:HG2	1.46	0.79
1:C:370:ILE:HB	1:C:403:ILE:CG2	2.13	0.79
1:B:579:GLU:OE1	1:B:584[A]:THR:HG21	1.83	0.79
1:A:355:HIS:HB2	1:A:502:MET:HE2	1.66	0.76
1:C:441:VAL:HG11	1:C:444:ARG:HD3	1.66	0.76
1:A:318:ARG:HD3	9:A:6062:HOH:O	1.86	0.76
1:C:398:GLU:HB2	1:C:400:ARG:HG2	1.68	0.75
1:D:594[A]:ARG:HG3	1:D:594[A]:ARG:NH1	1.93	0.75
1:C:385:ILE:HD13	1:C:416:VAL:HG12	1.68	0.75
1:D:411:LYS:NZ	1:D:411:LYS:HB3	2.02	0.75
1:B:289:ILE:HG23	1:B:434:MET:HB3	1.67	0.75
1:B:581:GLY:O	1:B:584[A]:THR:HG22	1.85	0.75
1:C:474:GLN:H	1:C:474:GLN:HE21	1.36	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:344:GLU:HG3	1:C:511:ARG:CZ	2.19	0.73
1:B:591:TYR:O	1:B:594:ARG:HG3	1.90	0.72
1:A:263:PRO:HB2	1:A:266:THR:HG23	1.72	0.71
1:B:600:GLN:CD	1:B:600:GLN:H	1.98	0.71
1:B:619:VAL:HG23	1:B:641:LEU:HD11	1.71	0.70
1:C:587:GLN:NE2	1:C:595:TYR:HA	2.06	0.70
1:C:394:ARG:HD2	1:C:400:ARG:HH12	1.57	0.69
1:C:366:ASN:HA	1:C:391:GLU:HG3	1.75	0.68
1:A:378:ASP:HB2	9:A:4126:HOH:O	1.93	0.68
1:A:221:LYS:HE2	9:A:5512:HOH:O	1.93	0.67
1:A:594[A]:ARG:HG3	1:A:594[A]:ARG:HH11	1.58	0.67
1:D:619:VAL:HG22	1:D:628:LYS:HG3	1.77	0.67
1:B:474:GLN:H	1:B:474:GLN:HE21	1.43	0.67
1:C:560:LEU:HD23	1:C:612:MET:HG3	1.77	0.66
1:C:352:LEU:HD23	1:C:352:LEU:H	1.60	0.66
1:C:370:ILE:HB	1:C:403:ILE:HG22	1.77	0.66
1:C:600:GLN:H	1:C:600:GLN:CD	2.04	0.66
1:A:621:LYS:HB3	1:A:623:GLU:OE2	1.95	0.66
1:C:102:MET:HE3	1:C:107:VAL:HG11	1.77	0.66
1:B:491:VAL:HA	9:B:5714:HOH:O	1.96	0.66
1:C:474:GLN:H	1:C:474:GLN:NE2	1.93	0.66
1:B:338:LEU:HD11	1:B:351:MET:HE3	1.77	0.65
1:B:335:LEU:HD12	1:B:351:MET:HE1	1.79	0.64
1:C:371:ILE:N	1:C:371:ILE:HD12	2.12	0.64
1:A:511:ARG:NH2	1:D:511:ARG:HH21	1.95	0.64
1:B:560:LEU:HD23	1:B:612:MET:HG3	1.80	0.64
1:A:251:LYS:HE2	9:B:5931:HOH:O	1.98	0.64
1:C:136:PRO:HG3	1:C:142:ALA:HB2	1.81	0.63
1:C:580:GLN:HE22	1:D:114:PRO:CG	2.12	0.63
1:B:338:LEU:HD11	1:B:351:MET:CE	2.28	0.63
1:C:451:ILE:HG22	1:C:455:LYS:HE2	1.80	0.63
1:C:554:ASN:HA	1:C:557[A]:LEU:HD12	1.79	0.63
1:D:411:LYS:HB3	1:D:411:LYS:HZ3	1.62	0.63
1:A:251:LYS:NZ	4:B:1695:1MM:HN8	1.97	0.63
1:B:496:GLY:H	1:B:500:HIS:HE1	1.48	0.62
1:C:102:MET:CE	1:C:158:VAL:HG11	2.28	0.62
1:C:292:ALA:HB2	1:C:421:ILE:HG21	1.80	0.62
1:A:496:GLY:H	1:A:500:HIS:HE1	1.48	0.62
1:A:511:ARG:NH2	1:D:511:ARG:NH2	2.47	0.62
1:B:304:LEU:HD23	1:B:371:ILE:HB	1.82	0.62
1:B:666:ASN:ND2	1:B:667:PHE:H	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:499:GLN:HE21	1:C:585:GLN:HE21	1.47	0.62
1:B:474:GLN:H	1:B:474:GLN:NE2	1.97	0.61
1:D:172:THR:HB	1:D:173:PRO:HD3	1.82	0.61
1:B:463:MET:HE3	1:B:463:MET:HA	1.82	0.61
1:C:396:ALA:HB2	1:C:402:GLY:HA2	1.82	0.61
1:D:594[A]:ARG:HD3	9:D:4998:HOH:O	1.99	0.61
1:A:313:HIS:CD2	1:A:315:ASP:H	2.11	0.61
1:C:522:LEU:HG	1:D:165:PRO:HD3	1.83	0.61
1:B:492:ILE:N	9:B:5714:HOH:O	2.29	0.61
1:B:623:GLU:CD	1:B:623:GLU:H	2.09	0.61
1:C:370:ILE:HB	1:C:403:ILE:HG23	1.82	0.61
1:C:499:GLN:HE21	1:C:585:GLN:NE2	1.99	0.61
1:A:594[A]:ARG:HG3	1:A:594[A]:ARG:NH1	2.16	0.61
1:B:102:MET:CE	1:B:158:VAL:HG11	2.29	0.61
1:B:136:PRO:HG3	1:B:142:ALA:HB2	1.82	0.61
1:A:172:THR:HB	1:A:173:PRO:HD3	1.83	0.60
1:B:600:GLN:H	1:B:600:GLN:NE2	1.99	0.60
1:C:347:LYS:HB2	1:C:347:LYS:NZ	2.15	0.60
1:C:293:ALA:HB2	1:C:434:MET:HG3	1.83	0.60
1:A:136:PRO:HG3	1:A:142:ALA:HB2	1.84	0.59
1:B:102:MET:HE3	1:B:107:VAL:HG11	1.83	0.59
1:D:170:VAL:HG23	1:D:174:MET:HE2	1.85	0.59
1:C:666:ASN:ND2	1:C:667:PHE:H	2.00	0.59
1:C:385:ILE:HB	9:C:5765:HOH:O	2.01	0.59
1:C:360:ALA:O	1:C:364:VAL:HG23	2.02	0.58
1:C:347:LYS:NZ	9:C:5476:HOH:O	2.33	0.58
1:D:621:LYS:HD2	9:D:5367:HOH:O	2.04	0.58
1:A:219:MET:HG3	1:A:248:ASP:HB3	1.84	0.58
1:A:226:LEU:HB3	1:A:227:PRO:HD3	1.85	0.58
1:C:619:VAL:HG23	1:C:641:LEU:HD11	1.85	0.58
1:C:512:ASN:HB2	9:C:5769:HOH:O	2.02	0.58
1:C:496:GLY:H	1:C:500:HIS:HE1	1.49	0.58
1:B:640:LEU:C	1:B:640:LEU:HD23	2.29	0.58
1:A:457:GLU:HG2	1:A:458:TYR:CE1	2.38	0.58
1:B:554:ASN:HA	1:B:557:LEU:HD13	1.86	0.57
1:A:251:LYS:HZ2	4:B:1695:1MM:HN8	1.51	0.57
1:C:435:MET:HG2	9:C:5209:HOH:O	2.04	0.57
1:D:289:ILE:HG23	1:D:434:MET:HB2	1.84	0.57
1:C:480:LEU:HD22	1:C:573:LEU:HD22	1.86	0.57
1:B:172:THR:HB	1:B:173:PRO:HD3	1.87	0.57
1:C:226:LEU:HB3	1:C:227:PRO:HD3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:502:MET:HE2	1:C:502:MET:HA	1.87	0.57
1:A:443:GLU:HG2	9:A:5554:HOH:O	2.05	0.56
1:C:580:GLN:NE2	9:C:4765:HOH:O	2.37	0.56
1:A:640:LEU:C	1:A:640:LEU:HD23	2.30	0.56
1:C:557[A]:LEU:HD11	1:D:558:THR:OG1	2.06	0.56
1:B:670:GLU:O	1:B:674:GLN:HG3	2.06	0.56
1:D:670:GLU:O	1:D:674:GLN:HG3	2.06	0.56
1:D:496:GLY:H	1:D:500:HIS:HE1	1.53	0.56
1:A:522:LEU:HG	1:B:165:PRO:HD3	1.87	0.56
1:A:132[B]:ASN:HD22	1:A:133:PHE:N	2.04	0.56
1:C:522:LEU:O	1:D:202:GLN:HG2	2.06	0.56
1:C:347:LYS:HE2	9:C:5476:HOH:O	2.06	0.55
4:C:3695:1MM:O7B	1:D:251:LYS:HD3	2.06	0.55
1:A:355:HIS:HB2	1:A:502:MET:CE	2.34	0.55
1:C:224:GLU:HG3	1:C:258:LEU:CD1	2.37	0.55
1:C:512:ASN:N	1:C:512:ASN:HD22	2.03	0.55
1:C:682:ARG:NH2	9:C:5797:HOH:O	2.38	0.55
1:D:594[B]:ARG:HD3	9:D:4998:HOH:O	2.06	0.55
1:C:429:THR:HG22	1:C:433:LYS:HE3	1.87	0.55
1:D:313:HIS:CD2	1:D:315:ASP:H	2.11	0.55
1:D:615:LYS:HE2	1:D:635:THR:HG21	1.88	0.55
1:C:557[B]:LEU:HD23	1:C:558:THR:N	2.22	0.55
9:A:4096:HOH:O	1:B:600:GLN:HB3	2.06	0.55
1:B:351:MET:HE1	1:B:505:ALA:HB1	1.88	0.55
1:B:410:PRO:HA	1:B:413:ILE:HD12	1.89	0.55
1:C:394:ARG:CD	1:C:400:ARG:HH12	2.20	0.55
1:D:89:PHE:CE1	1:D:100[A]:GLU:HG2	2.42	0.54
1:A:251:LYS:HD3	4:B:1695:1MM:O7B	2.07	0.54
1:C:565:GLN:HE22	1:D:600:GLN:HA	1.71	0.54
1:B:170:VAL:C	1:B:173:PRO:HD2	2.32	0.54
1:A:511:ARG:HH21	1:D:511:ARG:NH2	2.06	0.54
1:D:571:LYS:HB3	1:D:632:PHE:CZ	2.43	0.54
1:B:638:PRO:HD3	9:B:4328:HOH:O	2.07	0.54
1:C:554:ASN:HD21	1:C:602:ASN:ND2	2.05	0.53
1:C:494:THR:HA	1:C:517:ILE:O	2.08	0.53
1:A:89:PHE:CE2	1:A:100[A]:GLU:HG2	2.44	0.53
1:B:681:LYS:HB2	1:B:681:LYS:NZ	2.24	0.53
1:C:354:MET:HE1	1:C:380:ARG:NH1	2.23	0.53
1:D:226:LEU:HB3	1:D:227:PRO:HD3	1.89	0.53
1:C:624:GLU:HA	9:C:6033:HOH:O	2.07	0.53
1:D:170:VAL:C	1:D:173:PRO:HD2	2.34	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:ASP:HB3	9:C:5011:HOH:O	2.09	0.53
1:C:170:VAL:C	1:C:173:PRO:HD2	2.34	0.52
1:D:96:GLN:O	1:D:100[B]:GLU:HG2	2.09	0.52
1:A:170:VAL:C	1:A:173:PRO:HD2	2.34	0.52
1:B:554:ASN:HD21	1:B:602:ASN:ND2	2.07	0.52
1:C:103:SER:HB2	1:C:130:LYS:HD3	1.91	0.52
1:B:167:ALA:O	1:B:170:VAL:HG22	2.10	0.52
1:C:167:ALA:O	1:C:170:VAL:HG22	2.09	0.52
1:C:172:THR:HB	1:C:173:PRO:HD3	1.91	0.52
1:D:494:THR:HA	1:D:517:ILE:O	2.09	0.52
1:B:170:VAL:HG23	1:B:174:MET:HE2	1.91	0.52
1:A:344:GLU:HG3	9:A:6037:HOH:O	2.10	0.52
1:B:606:ILE:HD13	1:B:618:ARG:NH2	2.25	0.52
1:C:600:GLN:H	1:C:600:GLN:NE2	2.07	0.52
1:D:115:GLY:HA3	1:D:162:THR:HB	1.91	0.52
1:B:304:LEU:HB2	1:B:331:VAL:HG22	1.92	0.51
1:C:118:ILE:HG13	1:C:118:ILE:O	2.10	0.51
1:C:600:GLN:HB3	9:D:5772:HOH:O	2.10	0.51
1:C:335:LEU:HD12	1:C:351:MET:HE1	1.92	0.51
1:B:557:LEU:C	1:B:557:LEU:HD23	2.35	0.51
1:C:673:ARG:C	1:C:673:ARG:HD2	2.36	0.51
1:D:619:VAL:HG23	1:D:641:LEU:HD11	1.92	0.51
1:D:199:ASP:HA	9:D:6150:HOH:O	2.11	0.51
1:A:497:VAL:O	7:A:1698:P25:H62	2.10	0.51
1:B:295:LEU:HD12	1:B:421:ILE:HD12	1.93	0.51
1:C:306:VAL:HB	1:C:333:THR:HG22	1.92	0.51
1:C:347:LYS:CE	9:C:5476:HOH:O	2.59	0.50
1:A:591:TYR:O	1:A:594[A]:ARG:HG3	2.11	0.50
1:D:480:LEU:HD22	1:D:573:LEU:HD22	1.92	0.50
4:A:695:1MM:O7B	1:B:251:LYS:HD3	2.12	0.50
1:C:338:LEU:HD11	1:C:351:MET:CE	2.42	0.50
1:B:584[B]:THR:HG21	1:B:648:LYS:HD3	1.94	0.50
1:D:136:PRO:HG3	1:D:142:ALA:HB2	1.92	0.50
1:D:263:PRO:HB2	1:D:266:THR:HG23	1.94	0.50
1:A:582:MET:HG2	7:A:1698:P25:HM43	1.94	0.50
1:B:496:GLY:N	1:B:500:HIS:HE1	2.07	0.50
1:C:381:VAL:O	1:C:661:LEU:HD21	2.12	0.50
1:D:93:THR:H	1:D:96:GLN:HE21	1.59	0.50
1:B:606:ILE:O	1:B:610:GLU:HG3	2.12	0.49
1:A:480:LEU:HD22	1:A:573:LEU:HD22	1.94	0.49
1:D:141:GLY:O	1:D:145:MET:HG3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:584[A]:THR:HG23	9:B:4460:HOH:O	2.11	0.49
1:A:319:LEU:HD12	1:A:428:THR:HG23	1.95	0.49
1:B:322:GLU:OE2	1:B:435:MET:HG2	2.12	0.49
9:A:4937:HOH:O	1:D:347:LYS:HD3	2.12	0.49
1:C:111:PHE:O	1:C:159:VAL:HA	2.12	0.49
1:B:344:GLU:HG3	1:B:511:ARG:CZ	2.43	0.49
1:A:165:PRO:HB2	6:A:703:PYD:HM22	1.95	0.49
1:D:442:LYS:HE2	9:D:4193:HOH:O	2.13	0.49
1:D:640:LEU:HD23	1:D:640:LEU:C	2.38	0.49
1:B:342:ASP:OD2	1:B:344:GLU:HB2	2.13	0.49
1:B:512:ASN:HB2	1:B:515:THR:HG21	1.95	0.49
1:A:591:TYR:O	1:A:594[B]:ARG:HG3	2.13	0.49
1:B:118:ILE:HG13	1:B:118:ILE:O	2.13	0.49
1:B:622:GLN:HG3	9:B:5959:HOH:O	2.13	0.48
1:C:496:GLY:N	1:C:500:HIS:HE1	2.11	0.48
1:B:218:VAL:HG22	1:B:219:MET:N	2.29	0.48
1:B:494:THR:HA	1:B:517:ILE:O	2.12	0.48
1:C:122:TYR:HA	1:C:125:ILE:HG12	1.95	0.48
1:D:87:THR:HG22	1:D:261:PRO:HG3	1.96	0.48
1:A:298:LEU:C	1:A:298:LEU:HD12	2.39	0.48
1:B:111:PHE:O	1:B:159:VAL:HA	2.13	0.48
1:B:226:LEU:HB3	1:B:227:PRO:HD3	1.96	0.48
1:B:289:ILE:CG2	1:B:434:MET:HB3	2.39	0.48
1:C:351:MET:HE1	1:C:505:ALA:HB1	1.95	0.48
1:C:580:GLN:HE21	1:C:598:THR:HA	1.79	0.48
1:A:319:LEU:HD12	1:A:428:THR:CG2	2.43	0.48
1:D:502:MET:O	1:D:506:GLN:HG3	2.13	0.48
1:C:593:HIS:ND1	1:C:682:ARG:NH1	2.62	0.48
1:B:587:GLN:NE2	1:B:595:TYR:HA	2.29	0.47
1:C:465:GLU:HG3	1:C:472:LYS:HG2	1.96	0.47
1:D:554:ASN:HD21	1:D:602:ASN:ND2	2.11	0.47
1:D:580:GLN:NE2	9:D:4830:HOH:O	2.44	0.47
1:C:250:PRO:HB2	1:C:253:VAL:HG23	1.97	0.47
1:B:571:LYS:HB3	1:B:632:PHE:CZ	2.49	0.47
1:B:606:ILE:HD13	1:B:618:ARG:CZ	2.45	0.47
1:C:396:ALA:HB2	1:C:402:GLY:CA	2.44	0.47
1:B:98:PHE:O	1:B:102:MET:HG2	2.13	0.47
1:C:335:LEU:HA	1:C:351:MET:CE	2.35	0.47
1:A:494:THR:HA	1:A:517:ILE:O	2.14	0.47
1:B:581:GLY:HA2	1:B:584[A]:THR:HG22	1.96	0.47
1:D:167:ALA:O	1:D:170:VAL:HG22	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:594[A]:ARG:HD3	9:A:4406:HOH:O	2.14	0.47
1:C:378:ASP:OD2	1:C:380:ARG:HB2	2.15	0.47
1:C:623:GLU:CD	1:C:623:GLU:H	2.23	0.47
1:D:218:VAL:HG22	1:D:219:MET:N	2.30	0.46
1:C:557[B]:LEU:CD2	1:D:558:THR:HB	2.45	0.46
1:B:219:MET:HA	1:B:248:ASP:HB3	1.98	0.46
1:C:297:ASN:HA	9:C:5207:HOH:O	2.16	0.46
1:C:395:ALA:HB1	1:C:401:GLY:HA3	1.98	0.46
1:C:571:LYS:HB3	1:C:632:PHE:CZ	2.50	0.46
1:D:246:LEU:C	1:D:246:LEU:HD23	2.40	0.46
1:C:112:GLY:O	1:C:136:PRO:HD2	2.16	0.46
1:B:543:LEU:HG	1:B:543:LEU:O	2.16	0.46
1:A:200:ALA:HB1	4:B:1695:1MM:H5	1.97	0.46
1:C:103:SER:HA	9:C:5182:HOH:O	2.16	0.46
1:C:308:ALA:HB2	1:C:336:GLN:HB3	1.98	0.46
1:C:512:ASN:N	1:C:512:ASN:ND2	2.64	0.46
1:D:497:VAL:HG21	1:D:523:GLY:C	2.41	0.46
1:A:289:ILE:HG23	1:A:434:MET:HB2	1.98	0.46
1:B:344:GLU:HG3	1:B:511:ARG:NE	2.31	0.46
1:C:295:LEU:HD13	1:C:295:LEU:O	2.16	0.46
1:C:338:LEU:HD12	1:C:516:PHE:CE2	2.51	0.46
1:C:498:GLY:C	1:C:582:MET:HE2	2.41	0.46
1:C:640:LEU:C	1:C:640:LEU:HD23	2.41	0.46
1:B:447:TRP:HA	1:B:447:TRP:CE3	2.51	0.45
1:C:569:PRO:HD2	9:C:5452:HOH:O	2.16	0.45
1:A:170:VAL:HG23	1:A:174:MET:HE2	1.99	0.45
1:B:378:ASP:OD2	1:B:380:ARG:HB2	2.17	0.45
1:B:409:SER:HB3	9:B:4265:HOH:O	2.16	0.45
1:D:122:TYR:HA	1:D:125:ILE:HG12	1.96	0.45
1:A:618:ARG:HD3	1:A:620:LYS:HE2	1.99	0.45
1:C:395:ALA:HA	1:C:400:ARG:HG3	1.98	0.45
1:C:600:GLN:NE2	9:C:5050:HOH:O	2.49	0.45
1:D:129:ASP:HA	9:D:4852:HOH:O	2.16	0.45
1:D:571:LYS:HD2	1:D:639:VAL:CG1	2.46	0.45
1:D:606:ILE:HD11	1:D:617:LEU:O	2.16	0.45
1:A:447:TRP:HA	1:A:447:TRP:CE3	2.52	0.45
1:A:219:MET:HE2	1:A:221:LYS:HD3	1.99	0.45
1:A:290:ASN:OD1	1:A:437:LYS:HD3	2.17	0.45
1:D:219:MET:HG3	1:D:248:ASP:HB3	1.99	0.45
1:D:153:SER:HB3	1:D:538:ALA:HB1	1.99	0.45
1:A:115:GLY:HA3	1:A:162:THR:HB	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:ASP:OD2	1:C:155:LYS:HE2	2.17	0.45
1:C:369:LEU:HD11	1:C:404:ILE:HG13	1.98	0.45
1:C:447:TRP:CE3	1:C:447:TRP:HA	2.51	0.45
1:C:502:MET:O	1:C:506:GLN:HG3	2.17	0.45
1:D:228:LEU:HB2	1:D:266:THR:HB	1.99	0.45
1:A:132[B]:ASN:HD22	1:A:133:PHE:H	1.64	0.45
1:C:338:LEU:HD11	1:C:351:MET:HE3	1.99	0.45
1:D:118:ILE:O	1:D:118:ILE:HG13	2.17	0.45
1:B:543:LEU:HD21	1:B:633[A]:VAL:HG23	1.98	0.44
1:C:621:LYS:HB3	1:C:623:GLU:OE1	2.17	0.44
1:B:319:LEU:HD12	1:B:428:THR:HG23	1.99	0.44
1:C:347:LYS:HB2	1:C:347:LYS:HZ3	1.80	0.44
1:C:365:GLN:HG3	1:C:662:ASP:HB3	1.99	0.44
1:D:606:ILE:O	1:D:610[B]:GLU:HG3	2.17	0.44
1:A:670:GLU:O	1:A:674:GLN:HG3	2.18	0.44
1:D:89:PHE:HE1	1:D:100[A]:GLU:HG2	1.82	0.44
1:A:620:LYS:HE2	1:A:644:GLU:OE2	2.18	0.44
1:B:246:LEU:C	1:B:246:LEU:HD23	2.42	0.44
1:B:447:TRP:HA	1:B:447:TRP:HE3	1.82	0.44
1:D:496:GLY:N	1:D:500:HIS:HE1	2.14	0.44
1:A:167:ALA:O	1:A:170:VAL:HG22	2.18	0.44
1:A:246:LEU:C	1:A:246:LEU:HD23	2.43	0.44
1:B:139:GLU:HB2	1:B:169:ASN:HB3	1.99	0.44
1:B:473:PRO:HD3	1:B:645:VAL:HB	1.98	0.44
1:D:111:PHE:O	1:D:159:VAL:HA	2.18	0.44
1:B:206:VAL:HG21	9:B:4244:HOH:O	2.17	0.44
1:B:443:GLU:HA	9:B:4454:HOH:O	2.16	0.44
1:C:246:LEU:C	1:C:246:LEU:HD23	2.43	0.44
1:C:447:TRP:HA	1:C:447:TRP:HE3	1.83	0.44
1:C:478:LYS:HE3	1:C:507:HIS:ND1	2.32	0.44
1:C:653:PRO:HB2	9:C:5079:HOH:O	2.18	0.44
1:D:313:HIS:HD2	1:D:315:ASP:N	2.03	0.44
1:D:623:GLU:HG3	9:D:5898:HOH:O	2.17	0.44
1:A:165:PRO:HD3	1:B:522:LEU:HG	1.99	0.44
1:A:497:VAL:HG21	1:A:523:GLY:C	2.42	0.44
1:D:594[A]:ARG:NH1	1:D:594[A]:ARG:CG	2.70	0.44
1:B:480:LEU:HD22	1:B:573:LEU:HD22	2.00	0.44
1:B:600:GLN:CD	1:B:600:GLN:N	2.71	0.44
1:C:427:ALA:O	1:C:431:LEU:HB2	2.18	0.43
1:C:554:ASN:HA	1:C:557[B]:LEU:HD13	1.99	0.43
1:D:228:LEU:O	1:D:232:GLU:HG3	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:465:GLU:HG3	1:A:472:LYS:HG2	2.00	0.43
1:A:554:ASN:HA	1:A:557:LEU:HD23	2.00	0.43
1:C:342:ASP:OD1	1:C:344:GLU:HB2	2.18	0.43
1:A:97:ILE:O	1:A:101:MET:HG2	2.18	0.43
1:A:202:GLN:HG2	1:B:522:LEU:O	2.18	0.43
1:A:218:VAL:HG22	1:A:219:MET:N	2.34	0.43
1:B:318:ARG:HD3	9:B:4309:HOH:O	2.18	0.43
1:C:479:LYS:O	1:C:483:VAL:HG23	2.19	0.43
9:A:4194:HOH:O	1:D:442:LYS:HE3	2.17	0.43
1:C:494:THR:HG22	1:C:517:ILE:HB	2.00	0.43
1:D:621:LYS:HD3	9:D:6160:HOH:O	2.18	0.43
1:A:132[B]:ASN:ND2	1:A:133:PHE:N	2.67	0.43
1:B:115:GLY:HA3	1:B:162:THR:HB	2.01	0.43
1:B:621:LYS:HB3	1:B:623:GLU:OE1	2.18	0.43
1:C:200:ALA:HB1	4:C:2695:1MM:H5	2.00	0.43
1:D:93:THR:H	1:D:96:GLN:NE2	2.16	0.43
1:A:358:ALA:HB3	1:A:458:TYR:HB3	2.01	0.43
1:A:447:TRP:HA	1:A:447:TRP:HE3	1.83	0.43
1:B:311:LEU:O	1:B:514:HIS:HE1	2.01	0.43
1:B:581:GLY:C	1:B:584[A]:THR:HG22	2.44	0.43
1:C:165:PRO:HD3	1:D:522:LEU:HG	2.01	0.43
1:C:292:ALA:O	1:C:296:ILE:HG13	2.18	0.43
1:B:579:GLU:OE1	1:B:584[B]:THR:HG21	2.18	0.43
1:B:591:TYR:O	1:B:594:ARG:CG	2.65	0.43
1:C:474:GLN:HE21	1:C:474:GLN:N	2.11	0.43
1:D:139:GLU:HB2	1:D:169:ASN:HB3	2.00	0.43
1:D:497:VAL:O	8:D:2702:P22:H62	2.18	0.43
1:A:287[A]:GLN:NE2	1:A:291:LYS:NZ	2.66	0.43
1:D:554:ASN:HA	1:D:557:LEU:HD23	2.00	0.43
1:A:652:LEU:HA	1:A:653:PRO:C	2.43	0.43
1:C:497:VAL:HG21	1:C:523:GLY:C	2.44	0.43
1:B:681:LYS:HG2	9:B:4342:HOH:O	2.19	0.43
1:C:670:GLU:O	1:C:674:GLN:HG3	2.18	0.43
1:D:437:LYS:HG2	9:D:5562:HOH:O	2.19	0.43
1:C:652:LEU:HD13	1:C:667:PHE:HA	2.01	0.42
1:A:198:THR:O	1:A:199:ASP:HB2	2.20	0.42
1:B:295:LEU:HD12	1:B:421:ILE:CD1	2.49	0.42
1:B:512:ASN:N	1:B:512:ASN:ND2	2.65	0.42
1:C:407:GLU:O	1:C:424:GLU:HA	2.19	0.42
1:B:381:VAL:O	1:B:661:LEU:HD21	2.20	0.42
1:A:623:GLU:CD	1:A:623:GLU:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:LEU:O	1:A:514:HIS:HE1	2.02	0.42
1:A:336:GLN:HA	1:A:336:GLN:OE1	2.20	0.42
1:A:494:THR:HG22	1:A:517:ILE:HB	2.01	0.42
1:B:153:SER:HB3	1:B:538:ALA:HB1	2.02	0.42
1:D:301:LYS:HE2	1:D:366:ASN:O	2.19	0.42
1:B:398:GLU:HB2	1:B:400:ARG:HG2	2.01	0.42
1:C:398:GLU:HB2	1:C:400:ARG:CG	2.41	0.42
1:A:560:LEU:HG	1:A:614:LEU:HD21	2.02	0.42
1:B:679:ARG:HG2	1:B:687:HIS:OXT	2.20	0.42
1:C:352:LEU:HD13	1:C:364:VAL:HG21	2.02	0.42
1:D:115:GLY:HA3	1:D:162:THR:CB	2.50	0.42
1:C:129:ASP:HB3	9:C:5994:HOH:O	2.20	0.41
1:C:511:ARG:C	1:C:512:ASN:HD22	2.28	0.41
1:C:652:LEU:HA	1:C:653:PRO:C	2.45	0.41
1:A:496:GLY:N	1:A:500:HIS:HE1	2.13	0.41
1:B:582:MET:HE3	5:B:1701:FAD:HM71	2.02	0.41
1:C:318:ARG:HG2	1:C:322:GLU:OE2	2.20	0.41
1:B:502:MET:HA	1:B:502:MET:HE2	2.02	0.41
1:B:652:LEU:HA	1:B:653:PRO:C	2.44	0.41
1:C:498:GLY:O	1:C:501:GLN:HB3	2.21	0.41
1:A:615:LYS:HD3	1:A:635:THR:HG21	2.02	0.41
1:B:108:ASP:OD2	1:B:155:LYS:HE2	2.20	0.41
1:C:589:LEU:HG	9:C:6026:HOH:O	2.19	0.41
1:C:93:THR:OG1	1:C:96:GLN:HG3	2.20	0.41
1:C:619:VAL:CG2	1:C:641:LEU:HD11	2.48	0.41
1:A:594[A]:ARG:HH11	1:A:594[A]:ARG:CG	2.29	0.41
1:C:319:LEU:HD12	1:C:428:THR:CG2	2.51	0.41
1:A:118:ILE:HG13	1:A:118:ILE:O	2.21	0.41
1:A:153:SER:HB3	1:A:538:ALA:HB1	2.02	0.41
1:B:580:GLN:HG2	1:B:583:VAL:CG2	2.51	0.41
1:C:350:ASP:HB2	9:C:5038:HOH:O	2.19	0.41
1:D:101:MET:HA	1:D:101:MET:CE	2.42	0.41
1:A:111:PHE:O	1:A:159:VAL:HA	2.21	0.41
1:B:170:VAL:O	1:B:173:PRO:HD2	2.20	0.41
1:B:345:ASP:HA	1:B:346:PRO:HD3	1.92	0.41
1:C:681:LYS:HB2	1:C:681:LYS:NZ	2.36	0.41
1:A:115:GLY:HA3	1:A:162:THR:CB	2.51	0.41
1:B:543:LEU:C	9:B:5714:HOH:O	2.63	0.41
1:C:115:GLY:HA3	1:C:162:THR:HB	2.03	0.41
1:C:600:GLN:CD	9:C:5050:HOH:O	2.64	0.41
1:A:220:VAL:HG13	1:A:225:GLU:HG3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:606:ILE:HD11	1:B:618:ARG:HB2	2.03	0.41
1:A:560:LEU:HD23	1:A:612:MET:HG3	2.03	0.40
1:B:97:ILE:O	1:B:101:MET:HG2	2.21	0.40
1:B:494:THR:HG21	1:B:532:ALA:HA	2.02	0.40
1:D:281:GLN:O	1:D:285:VAL:HG23	2.21	0.40
1:A:502:MET:O	1:A:506:GLN:HG3	2.21	0.40
1:B:319:LEU:HD12	1:B:428:THR:CG2	2.51	0.40
1:D:571:LYS:HD2	1:D:639:VAL:HG12	2.02	0.40
1:B:321:LYS:NZ	9:B:4719:HOH:O	2.54	0.40
1:B:529:LEU:HB3	1:B:530:PRO:CD	2.52	0.40
1:B:606:ILE:HD12	1:B:606:ILE:N	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	598/677 (88%)	587 (98%)	11 (2%)	0	100	100
1	B	581/677 (86%)	568 (98%)	12 (2%)	1 (0%)	43	55
1	C	573/677 (85%)	551 (96%)	21 (4%)	1 (0%)	43	55
1	D	597/677 (88%)	586 (98%)	11 (2%)	0	100	100
All	All	2349/2708 (87%)	2292 (98%)	55 (2%)	2 (0%)	48	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	350	ASP
1	C	350	ASP

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	488/556 (88%)	480 (98%)	8 (2%)	55	73
1	B	464/556 (84%)	457 (98%)	7 (2%)	57	75
1	C	453/556 (82%)	441 (97%)	12 (3%)	40	59
1	D	483/556 (87%)	473 (98%)	10 (2%)	47	66
All	All	1888/2224 (85%)	1851 (98%)	37 (2%)	51	67

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	287[A]	GLN
1	A	287[B]	GLN
1	A	447	TRP
1	A	457	GLU
1	A	548	ASP
1	A	560	LEU
1	A	614	LEU
1	A	630	LYS
1	B	224	GLU
1	B	447	TRP
1	B	474	GLN
1	B	541	GLU
1	B	548	ASP
1	B	557	LEU
1	B	681	LYS
1	C	295	LEU
1	C	304	LEU
1	C	347	LYS
1	C	371	ILE
1	C	403	ILE
1	C	447	TRP
1	C	474	GLN
1	C	548	ASP
1	C	557[A]	LEU

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Mol	Chain	Res	Type
1	C	557[B]	LEU
1	C	560	LEU
1	C	681	LYS
1	D	136	PRO
1	D	298	LEU
1	D	347	LYS
1	D	411	LYS
1	D	548	ASP
1	D	589	LEU
1	D	594[A]	ARG
1	D	594[B]	ARG
1	D	614	LEU
1	D	681	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (59) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	99	ASN
1	A	202	GLN
1	A	312	ASN
1	A	313	HIS
1	A	500	HIS
1	A	512	ASN
1	A	554	ASN
1	A	666	ASN
1	B	106	ASN
1	B	126	HIS
1	B	132	ASN
1	B	169	ASN
1	B	312	ASN
1	B	328	GLN
1	B	452	ASN
1	B	474	GLN
1	B	499	GLN
1	B	500	HIS
1	B	512	ASN
1	B	514	HIS
1	B	554	ASN
1	B	587	GLN
1	B	600	GLN
1	B	602	ASN
1	B	666	ASN

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Mol	Chain	Res	Type
1	B	674	GLN
1	C	99	ASN
1	C	105	GLN
1	C	132	ASN
1	C	169	ASN
1	C	312	ASN
1	C	355	HIS
1	C	412	ASN
1	C	452	ASN
1	C	474	GLN
1	C	500	HIS
1	C	512	ASN
1	C	514	HIS
1	C	554	ASN
1	C	565	GLN
1	C	580	GLN
1	C	585	GLN
1	C	587	GLN
1	C	600	GLN
1	C	602	ASN
1	C	666	ASN
1	D	96	GLN
1	D	106	ASN
1	D	127	ASN
1	D	169	ASN
1	D	202	GLN
1	D	312	ASN
1	D	313	HIS
1	D	450	GLN
1	D	500	HIS
1	D	554	ASN
1	D	580	GLN
1	D	602	ASN
1	D	675	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 8 are monoatomic - leaving 16 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$
6	PYD	A	703	-	9,9,9	8.32	7 (77%)	11,12,12	3.95	7 (63%)
6	PYD	C	2703	-	9,9,9	8.54	6 (66%)	11,12,12	3.99	7 (63%)
5	FAD	A	701	-	58,58,58	2.66	24 (41%)	85,89,89	1.55	13 (15%)
7	P25	B	698	3	12,13,13	1.99	4 (33%)	16,18,18	3.17	5 (31%)
4	1MM	A	695	-	27,27,27	3.80	18 (66%)	37,38,38	4.08	17 (45%)
8	P22	C	3702	3	9,10,10	1.92	4 (44%)	13,15,15	1.98	2 (15%)
5	FAD	B	1701	-	58,58,58	2.81	25 (43%)	85,89,89	1.56	14 (16%)
4	1MM	B	1695	-	27,27,27	4.01	18 (66%)	37,38,38	4.13	17 (45%)
8	P22	D	2702	3	9,10,10	2.03	4 (44%)	13,15,15	1.89	1 (7%)
5	FAD	C	2701	-	58,58,58	2.88	27 (46%)	85,89,89	1.56	13 (15%)
4	1MM	C	3695	-	27,27,27	4.00	19 (70%)	37,38,38	4.17	16 (43%)
4	1MM	C	2695	-	27,27,27	3.86	18 (66%)	37,38,38	4.07	18 (48%)
5	FAD	D	3701	-	58,58,58	2.70	25 (43%)	85,89,89	1.54	14 (16%)
6	PYD	D	3703	-	9,9,9	8.42	7 (77%)	11,12,12	3.90	7 (63%)
7	P25	A	1698	3	12,13,13	1.97	4 (33%)	16,18,18	3.22	5 (31%)
6	PYD	B	1703	-	9,9,9	8.34	7 (77%)	11,12,12	4.00	7 (63%)

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
6	PYD	A	703	-	-	-	0/1/1/1
7	P25	B	698	3	-	0/13/13/13	-
5	FAD	A	701	-	-	2/34/50/50	0/6/6/6
6	PYD	B	1703	-	-	-	0/1/1/1
4	1MM	A	695	-	-	11/23/23/23	0/2/2/2
6	PYD	C	2703	-	-	-	0/1/1/1
5	FAD	B	1701	-	-	2/34/50/50	0/6/6/6
4	1MM	B	1695	-	-	12/23/23/23	0/2/2/2
8	P22	D	2702	3	-	2/10/10/10	-
5	FAD	C	2701	-	-	1/34/50/50	0/6/6/6
4	1MM	C	3695	-	-	10/23/23/23	0/2/2/2
4	1MM	C	2695	-	-	12/23/23/23	0/2/2/2
5	FAD	D	3701	-	-	1/34/50/50	0/6/6/6
7	P25	A	1698	3	-	2/13/13/13	-
6	PYD	D	3703	-	-	-	0/1/1/1
8	P22	C	3702	3	-	1/10/10/10	-

All (217) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	3703	PYD	C4'-N3'	13.82	1.53	1.35
6	C	2703	PYD	C4'-N3'	13.17	1.52	1.35
6	A	703	PYD	C4'-N3'	13.14	1.52	1.35
6	B	1703	PYD	C4'-N3'	13.12	1.52	1.35
6	C	2703	PYD	C5'-C4'	12.34	1.60	1.41
6	D	3703	PYD	C5'-C4'	11.54	1.59	1.41
6	B	1703	PYD	C5'-C4'	11.54	1.59	1.41
6	A	703	PYD	C5'-C4'	11.45	1.59	1.41
6	A	703	PYD	C2'-N1'	10.67	1.50	1.34
6	B	1703	PYD	C2'-N1'	10.62	1.50	1.34
6	D	3703	PYD	C6'-C5'	10.47	1.55	1.39
6	C	2703	PYD	C2'-N1'	10.42	1.50	1.34
6	B	1703	PYD	C6'-N1'	10.30	1.55	1.34
6	C	2703	PYD	C6'-C5'	10.24	1.54	1.39
6	A	703	PYD	C6'-N1'	10.12	1.55	1.34
6	C	2703	PYD	C6'-N1'	10.10	1.55	1.34
6	D	3703	PYD	C6'-N1'	10.02	1.54	1.34
6	A	703	PYD	C6'-C5'	9.69	1.54	1.39
6	B	1703	PYD	C6'-C5'	9.63	1.53	1.39
6	D	3703	PYD	C2'-N1'	8.99	1.48	1.34
5	C	2701	FAD	C4A-N3A	7.96	1.49	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1701	FAD	C4A-N3A	7.78	1.49	1.34
4	B	1695	1MM	C3-C2	7.39	1.47	1.39
5	D	3701	FAD	C4A-N3A	7.34	1.48	1.34
4	C	3695	1MM	C3-C2	7.34	1.47	1.39
4	C	2695	1MM	C4'-N5'	7.21	1.46	1.33
4	A	695	1MM	C4'-N5'	7.20	1.46	1.33
4	B	1695	1MM	C4'-N5'	7.19	1.46	1.33
4	C	3695	1MM	C4'-N5'	7.17	1.46	1.33
4	A	695	1MM	C3-C2	7.16	1.47	1.39
5	A	701	FAD	C4A-N3A	6.99	1.47	1.34
5	C	2701	FAD	C2A-N1A	6.94	1.46	1.33
4	C	3695	1MM	C6'-N1'	6.93	1.45	1.34
4	C	2695	1MM	C3-C2	6.83	1.47	1.39
4	B	1695	1MM	C6'-N1'	6.57	1.45	1.34
5	B	1701	FAD	C2A-N1A	6.52	1.45	1.33
4	C	3695	1MM	C6'-N5'	6.38	1.45	1.34
5	D	3701	FAD	C2A-N1A	6.38	1.45	1.33
4	A	695	1MM	C6'-N1'	6.36	1.45	1.34
4	B	1695	1MM	C6'-N5'	6.24	1.44	1.34
5	C	2701	FAD	C5A-C4A	6.15	1.50	1.39
5	A	701	FAD	C2A-N1A	6.12	1.44	1.33
4	C	2695	1MM	C6'-N5'	6.05	1.44	1.34
4	A	695	1MM	C6'-N5'	5.98	1.44	1.34
4	C	2695	1MM	C6'-N1'	5.98	1.44	1.34
4	C	3695	1MM	C4'-N3'	5.97	1.44	1.33
5	B	1701	FAD	C5A-C4A	5.96	1.49	1.39
4	B	1695	1MM	C1-C2	5.94	1.46	1.40
5	A	701	FAD	C8A-N9A	5.76	1.47	1.37
5	C	2701	FAD	C5A-C6A	5.72	1.56	1.41
5	B	1701	FAD	C5A-C6A	5.70	1.56	1.41
4	C	3695	1MM	C1-C2	5.70	1.46	1.40
4	B	1695	1MM	C4'-N3'	5.70	1.43	1.33
5	B	1701	FAD	C8A-N9A	5.70	1.47	1.37
5	D	3701	FAD	C5A-C4A	5.65	1.49	1.39
5	A	701	FAD	C5A-C4A	5.65	1.49	1.39
4	C	2695	1MM	C1-C2	5.60	1.46	1.40
5	D	3701	FAD	C8A-N9A	5.60	1.47	1.37
5	C	2701	FAD	C8A-N9A	5.57	1.47	1.37
5	A	701	FAD	C5A-C6A	5.55	1.56	1.41
5	D	3701	FAD	C5A-C6A	5.37	1.55	1.41
4	C	2695	1MM	C4'-N3'	5.33	1.43	1.33
4	B	1695	1MM	C2'-N10	-5.16	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	2695	1MM	O12-C11	5.12	1.44	1.33
5	C	2701	FAD	C1'-C2'	-5.12	1.45	1.52
4	B	1695	1MM	C4-C3	5.11	1.47	1.38
4	A	695	1MM	C2'-N10	-5.03	1.32	1.38
4	C	3695	1MM	C4-C3	4.99	1.47	1.38
4	C	2695	1MM	C2'-N10	-4.94	1.32	1.38
4	C	3695	1MM	O12-C11	4.91	1.44	1.33
4	A	695	1MM	C4'-N3'	4.91	1.42	1.33
4	A	695	1MM	C4-C3	4.83	1.47	1.38
4	B	1695	1MM	O12-C11	4.83	1.44	1.33
4	C	2695	1MM	C4-C3	4.81	1.47	1.38
4	A	695	1MM	C1-C2	4.79	1.45	1.40
4	A	695	1MM	O12-C11	4.77	1.44	1.33
5	C	2701	FAD	O4'-C4'	4.68	1.53	1.43
5	B	1701	FAD	C1'-C2'	-4.63	1.46	1.52
5	A	701	FAD	O4'-C4'	4.55	1.52	1.43
4	C	3695	1MM	C2'-N10	-4.52	1.32	1.38
5	B	1701	FAD	O4'-C4'	4.51	1.52	1.43
5	D	3701	FAD	C1'-C2'	-4.47	1.46	1.52
5	C	2701	FAD	C9A-C5X	4.40	1.48	1.41
5	D	3701	FAD	O4'-C4'	4.38	1.52	1.43
5	B	1701	FAD	C9A-C5X	4.33	1.48	1.41
5	C	2701	FAD	C2A-N3A	4.31	1.41	1.33
5	A	701	FAD	C1'-C2'	-4.26	1.46	1.52
5	A	701	FAD	C9A-C5X	4.17	1.47	1.41
5	D	3701	FAD	C9A-C5X	4.15	1.47	1.41
5	C	2701	FAD	C6A-N1A	4.06	1.47	1.35
6	D	3703	PYD	C2'-N3'	4.06	1.41	1.34
5	D	3701	FAD	C6A-N1A	4.05	1.47	1.35
7	B	698	P25	PA-O7	-4.05	1.43	1.59
5	C	2701	FAD	C5'-C4'	-4.04	1.46	1.51
5	B	1701	FAD	O3'-C3'	4.02	1.52	1.43
5	B	1701	FAD	C6A-N1A	4.00	1.47	1.35
5	D	3701	FAD	O3'-C3'	4.00	1.52	1.43
5	C	2701	FAD	O3'-C3'	3.95	1.52	1.43
5	D	3701	FAD	C5'-C4'	-3.94	1.46	1.51
4	B	1695	1MM	C2-S7	3.92	1.83	1.77
5	A	701	FAD	O3'-C3'	3.91	1.52	1.43
7	A	1698	P25	PA-O7	-3.91	1.44	1.59
4	C	3695	1MM	C6-C1	3.86	1.45	1.39
4	C	2695	1MM	C6-C1	3.84	1.45	1.39
4	B	1695	1MM	C6-C1	3.79	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1701	FAD	C2A-N3A	3.77	1.40	1.33
5	A	701	FAD	C6A-N1A	3.76	1.46	1.35
5	D	3701	FAD	C2A-N3A	3.69	1.40	1.33
5	B	1701	FAD	C4X-N5	3.67	1.38	1.30
5	B	1701	FAD	C5'-C4'	-3.62	1.46	1.51
5	A	701	FAD	C5'-C4'	-3.61	1.46	1.51
5	C	2701	FAD	C4X-N5	3.55	1.38	1.30
4	C	3695	1MM	C2-S7	3.55	1.82	1.77
5	A	701	FAD	C2A-N3A	3.49	1.40	1.33
4	C	3695	1MM	C2'-N1'	3.47	1.44	1.34
4	A	695	1MM	C6-C1	3.44	1.45	1.39
4	C	2695	1MM	C2'-N1'	3.39	1.44	1.34
4	C	2695	1MM	C2-S7	3.39	1.82	1.77
4	B	1695	1MM	C2'-N1'	3.36	1.44	1.34
5	B	1701	FAD	C10-N1	3.32	1.39	1.33
5	A	701	FAD	C4X-N5	3.31	1.37	1.30
4	A	695	1MM	C2'-N1'	3.30	1.43	1.34
6	C	2703	PYD	C2'-N3'	3.29	1.39	1.34
4	A	695	1MM	C2-S7	3.27	1.82	1.77
5	C	2701	FAD	C10-N1	3.26	1.39	1.33
5	D	3701	FAD	C10-N1	3.24	1.39	1.33
8	D	2702	P22	PA-O7	-3.22	1.46	1.59
5	D	3701	FAD	C4X-N5	3.22	1.37	1.30
7	A	1698	P25	O7-C7	3.19	1.57	1.44
4	A	695	1MM	C9-N10	-3.09	1.30	1.37
4	C	2695	1MM	C9-N10	-3.08	1.30	1.37
5	C	2701	FAD	C1B-N9A	3.07	1.54	1.46
4	C	3695	1MM	C2'-N3'	2.99	1.43	1.34
5	A	701	FAD	C3B-C4B	2.98	1.60	1.53
5	C	2701	FAD	C8-C7	2.97	1.48	1.40
7	B	698	P25	O7-C7	2.94	1.56	1.44
5	B	1701	FAD	C1B-N9A	2.94	1.54	1.46
4	C	2695	1MM	C2'-N3'	2.94	1.42	1.34
4	B	1695	1MM	C9-N10	-2.93	1.31	1.37
5	D	3701	FAD	C8-C7	2.89	1.47	1.40
4	B	1695	1MM	C2'-N3'	2.89	1.42	1.34
5	B	1701	FAD	C3B-C4B	2.88	1.60	1.53
4	C	2695	1MM	C5-C6	2.87	1.43	1.38
5	A	701	FAD	C10-N1	2.84	1.39	1.33
5	D	3701	FAD	C3B-C4B	2.84	1.60	1.53
4	C	3695	1MM	C5-C6	2.83	1.43	1.38
5	C	2701	FAD	C3B-C4B	2.81	1.60	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	701	FAD	C8-C7	2.79	1.47	1.40
4	A	695	1MM	C2'-N3'	2.78	1.42	1.34
5	D	3701	FAD	C6-C5X	2.77	1.44	1.40
4	B	1695	1MM	C9-N8	-2.72	1.33	1.39
4	C	3695	1MM	O4'-C4'	2.71	1.39	1.33
4	C	3695	1MM	C9-N10	-2.71	1.31	1.37
4	B	1695	1MM	C5-C6	2.69	1.43	1.38
4	B	1695	1MM	O4'-C4'	2.68	1.39	1.33
5	B	1701	FAD	C8-C7	2.68	1.47	1.40
5	A	701	FAD	C1B-N9A	2.66	1.53	1.46
5	D	3701	FAD	P-O2P	-2.66	1.43	1.55
5	A	701	FAD	C9-C9A	2.64	1.43	1.39
5	C	2701	FAD	P-O2P	-2.63	1.43	1.55
5	A	701	FAD	P-O2P	-2.63	1.43	1.55
5	B	1701	FAD	P-O2P	-2.61	1.43	1.55
7	A	1698	P25	PB-O2B	-2.61	1.45	1.54
5	B	1701	FAD	C6-C5X	2.61	1.43	1.40
7	B	698	P25	PB-O2B	-2.59	1.45	1.54
4	A	695	1MM	C5-C6	2.58	1.43	1.38
5	B	1701	FAD	C9-C9A	2.58	1.43	1.39
4	C	2695	1MM	O4'-C4'	2.56	1.39	1.33
5	B	1701	FAD	C6-C7	2.56	1.43	1.39
4	A	695	1MM	C9-N8	-2.56	1.33	1.39
7	B	698	P25	PA-O2A	-2.55	1.43	1.55
4	C	2695	1MM	C1-C11	2.55	1.55	1.50
4	C	3695	1MM	S7-N8	2.54	1.70	1.64
5	C	2701	FAD	C9-C9A	2.53	1.43	1.39
8	C	3702	P22	PA-O2A	-2.53	1.42	1.50
8	D	2702	P22	PB-O1B	-2.50	1.45	1.54
7	A	1698	P25	PA-O2A	-2.49	1.43	1.55
4	A	695	1MM	O4'-C4'	2.47	1.39	1.33
5	C	2701	FAD	C6-C5X	2.46	1.43	1.40
5	C	2701	FAD	C6-C7	2.40	1.42	1.39
5	D	3701	FAD	C1B-N9A	2.39	1.53	1.46
4	C	2695	1MM	C9-N8	-2.39	1.34	1.39
5	D	3701	FAD	PA-O2A	-2.38	1.44	1.55
5	D	3701	FAD	C6-C7	2.36	1.42	1.39
5	C	2701	FAD	C10-N10	2.33	1.42	1.37
6	A	703	PYD	C2'-N3'	2.32	1.38	1.34
5	B	1701	FAD	PA-O5B	-2.30	1.50	1.59
8	C	3702	P22	PA-O7	-2.29	1.50	1.59
5	B	1701	FAD	C10-N10	2.28	1.42	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	D	2702	P22	PA-O1A	-2.24	1.45	1.55
4	C	3695	1MM	C1-C11	2.23	1.54	1.50
8	C	3702	P22	PB-O1B	-2.22	1.46	1.54
5	C	2701	FAD	PA-O2A	-2.19	1.45	1.55
5	D	3701	FAD	C10-N10	2.18	1.42	1.37
5	C	2701	FAD	C9-C8	2.17	1.42	1.39
5	D	3701	FAD	PA-O5B	-2.17	1.50	1.59
5	C	2701	FAD	PA-O5B	-2.17	1.50	1.59
5	C	2701	FAD	C5A-N7A	-2.16	1.35	1.39
4	A	695	1MM	C1-C11	2.16	1.54	1.50
5	A	701	FAD	PA-O2A	-2.16	1.45	1.55
5	A	701	FAD	PA-O5B	-2.15	1.51	1.59
6	B	1703	PYD	C2'-N3'	2.15	1.37	1.34
5	D	3701	FAD	C9-C9A	2.14	1.43	1.39
4	B	1695	1MM	C1-C11	2.14	1.54	1.50
5	A	701	FAD	C6-C5X	2.12	1.43	1.40
5	B	1701	FAD	C9-C8	2.12	1.42	1.39
5	B	1701	FAD	PA-O2A	-2.12	1.45	1.55
5	A	701	FAD	C6-C7	2.11	1.42	1.39
4	C	3695	1MM	C9-N8	-2.10	1.34	1.39
5	C	2701	FAD	C4'-C3'	2.07	1.57	1.53
8	C	3702	P22	O7-C7	2.07	1.55	1.44
5	A	701	FAD	C9-C8	2.05	1.42	1.39
8	D	2702	P22	PA-O2A	-2.04	1.43	1.50
6	B	1703	PYD	CM2-C2'	-2.02	1.44	1.49
6	A	703	PYD	CM2-C2'	-2.01	1.44	1.49
6	D	3703	PYD	C7'-C5'	2.01	1.54	1.51
5	D	3701	FAD	C5A-N7A	-2.01	1.35	1.39

All (163) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	3695	1MM	C5'-O4'-C4'	14.32	140.19	117.56
4	B	1695	1MM	C5'-O4'-C4'	14.21	140.01	117.56
4	C	2695	1MM	C5'-O4'-C4'	14.13	139.88	117.56
4	A	695	1MM	C5'-O4'-C4'	13.96	139.62	117.56
4	C	2695	1MM	C3-C2-S7	-9.98	103.17	117.51
4	B	1695	1MM	C3-C2-S7	-9.76	103.49	117.51
4	A	695	1MM	C3-C2-S7	-9.73	103.53	117.51
4	C	3695	1MM	C4'-N3'-C2'	9.68	120.76	112.86
4	C	3695	1MM	C3-C2-S7	-9.65	103.65	117.51
4	A	695	1MM	C4'-N3'-C2'	9.35	120.49	112.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1695	1MM	C4'-N3'-C2'	9.25	120.41	112.86
4	C	2695	1MM	C4'-N3'-C2'	8.96	120.18	112.86
7	B	698	P25	O7-C7-C6	8.43	139.20	109.28
6	B	1703	PYD	CM2-C2'-N1'	8.33	126.07	117.20
6	C	2703	PYD	CM2-C2'-N1'	8.31	126.04	117.20
7	A	1698	P25	O7-C7-C6	8.30	138.71	109.28
6	A	703	PYD	CM2-C2'-N1'	8.23	125.96	117.20
6	D	3703	PYD	CM2-C2'-N1'	8.23	125.96	117.20
4	C	3695	1MM	C2-C1-C11	-6.09	116.12	124.25
6	B	1703	PYD	C5'-C6'-N1'	-6.06	119.02	125.09
7	A	1698	P25	C5-C6-C7	5.99	139.48	113.47
6	A	703	PYD	C5'-C6'-N1'	-5.94	119.14	125.09
4	A	695	1MM	C2-C1-C11	-5.90	116.38	124.25
4	B	1695	1MM	C2-C1-C11	-5.76	116.56	124.25
7	B	698	P25	C5-C6-C7	5.71	138.27	113.47
8	C	3702	P22	PA-O7-C7	5.65	151.88	121.12
6	C	2703	PYD	C5'-C6'-N1'	-5.56	119.52	125.09
8	D	2702	P22	PA-O7-C7	5.54	151.31	121.12
4	C	2695	1MM	C2-C1-C11	-5.34	117.12	124.25
7	A	1698	P25	PA-O7-C7	5.33	146.65	121.26
4	C	2695	1MM	C2'-N1'-C6'	5.01	122.08	114.61
7	B	698	P25	PA-O7-C7	5.01	145.11	121.26
4	C	3695	1MM	C2'-N1'-C6'	5.01	122.08	114.61
4	B	1695	1MM	C2'-N1'-C6'	5.00	122.06	114.61
5	C	2701	FAD	C4A-N9A-C1B	4.99	138.31	126.63
6	C	2703	PYD	C6'-N1'-C2'	4.96	124.22	116.07
4	A	695	1MM	C2'-N1'-C6'	4.95	122.00	114.61
6	D	3703	PYD	C6'-N1'-C2'	4.87	124.06	116.07
6	B	1703	PYD	C6'-N1'-C2'	4.85	124.04	116.07
6	A	703	PYD	C6'-N1'-C2'	4.84	124.02	116.07
5	B	1701	FAD	C4A-N9A-C1B	4.81	137.88	126.63
4	B	1695	1MM	N8-C9-N10	-4.79	107.27	114.87
6	D	3703	PYD	C5'-C6'-N1'	-4.71	120.36	125.09
4	C	3695	1MM	N8-C9-N10	-4.67	107.46	114.87
6	D	3703	PYD	N1'-C2'-N3'	-4.65	117.80	125.53
4	A	695	1MM	N8-C9-N10	-4.64	107.50	114.87
5	D	3701	FAD	C4A-N9A-C1B	4.61	137.41	126.63
5	A	701	FAD	C4A-N9A-C1B	4.56	137.29	126.63
4	C	2695	1MM	N8-C9-N10	-4.56	107.63	114.87
6	C	2703	PYD	N1'-C2'-N3'	-4.55	117.96	125.53
4	C	3695	1MM	N3'-C2'-N1'	-4.51	118.87	126.26
6	A	703	PYD	N1'-C2'-N3'	-4.39	118.22	125.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	1703	PYD	N1'-C2'-N3'	-4.37	118.25	125.53
4	C	2695	1MM	N5'-C6'-N1'	-4.34	118.12	125.77
4	A	695	1MM	N5'-C6'-N1'	-4.32	118.15	125.77
4	A	695	1MM	N3'-C2'-N1'	-4.29	119.22	126.26
4	B	1695	1MM	N3'-C2'-N1'	-4.29	119.23	126.26
4	B	1695	1MM	N5'-C6'-N1'	-4.24	118.28	125.77
4	C	2695	1MM	N3'-C2'-N1'	-4.19	119.39	126.26
4	B	1695	1MM	O9-C9-N10	4.16	131.04	123.64
4	C	3695	1MM	N5'-C6'-N1'	-4.15	118.44	125.77
4	C	3695	1MM	O12-C11-C1	-4.15	105.42	112.31
4	A	695	1MM	O9-C9-N10	4.07	130.87	123.64
4	C	2695	1MM	N3'-C4'-N5'	-4.06	120.28	127.65
4	C	2695	1MM	O12-C11-C1	-4.05	105.59	112.31
4	C	3695	1MM	N3'-C4'-N5'	-4.00	120.38	127.65
4	C	3695	1MM	O12-C11-O11	3.99	131.22	123.46
4	C	3695	1MM	O9-C9-N10	3.98	130.71	123.64
4	B	1695	1MM	N3'-C4'-N5'	-3.97	120.43	127.65
4	B	1695	1MM	O12-C11-C1	-3.96	105.74	112.31
4	B	1695	1MM	O12-C11-O11	3.93	131.11	123.46
4	A	695	1MM	O12-C11-O11	3.91	131.06	123.46
4	A	695	1MM	O12-C11-C1	-3.89	105.85	112.31
4	A	695	1MM	N3'-C4'-N5'	-3.89	120.59	127.65
4	C	2695	1MM	O9-C9-N10	3.87	130.52	123.64
4	B	1695	1MM	O7B-S7-C2	3.81	113.97	107.68
4	C	3695	1MM	O7B-S7-C2	3.73	113.84	107.68
4	C	2695	1MM	O12-C11-O11	3.67	130.59	123.46
5	C	2701	FAD	C4A-N9A-C8A	-3.66	101.89	105.74
5	B	1701	FAD	C4A-N9A-C8A	-3.56	102.00	105.74
5	A	701	FAD	C6A-C5A-C4A	-3.55	112.33	117.18
5	D	3701	FAD	C6A-C5A-C4A	-3.52	112.37	117.18
6	D	3703	PYD	C7'-C5'-C4'	3.52	125.44	121.74
5	A	701	FAD	C4A-N9A-C8A	-3.47	102.09	105.74
4	C	3695	1MM	C6-C1-C11	3.46	125.65	118.66
4	C	2695	1MM	O7B-S7-C2	3.44	113.37	107.68
4	A	695	1MM	O7B-S7-C2	3.44	113.37	107.68
5	B	1701	FAD	C6A-C5A-C4A	-3.44	112.48	117.18
6	C	2703	PYD	C7'-C5'-C4'	3.34	125.26	121.74
4	A	695	1MM	C6-C1-C11	3.31	125.36	118.66
5	C	2701	FAD	C1B-N9A-C8A	-3.30	119.77	127.09
5	D	3701	FAD	N3A-C2A-N1A	-3.29	123.60	128.58
6	D	3703	PYD	C2'-N3'-C4'	3.26	123.12	118.10
4	B	1695	1MM	C6-C1-C11	3.25	125.23	118.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	2701	FAD	O2B-C2B-C1B	3.23	121.23	110.10
5	B	1701	FAD	O2B-C2B-C1B	3.21	121.17	110.10
5	A	701	FAD	N3A-C2A-N1A	-3.21	123.73	128.58
5	A	701	FAD	C6A-C5A-N7A	3.21	138.27	132.09
7	A	1698	P25	CM4-C4-C5	3.18	134.84	113.36
5	C	2701	FAD	C6A-C5A-C4A	-3.17	112.85	117.18
5	D	3701	FAD	C4A-N9A-C8A	-3.17	102.41	105.74
5	B	1701	FAD	C1B-N9A-C8A	-3.15	120.09	127.09
5	D	3701	FAD	O2B-C2B-C1B	3.13	120.90	110.10
6	C	2703	PYD	C2'-N3'-C4'	3.13	122.92	118.10
4	C	2695	1MM	C6-C1-C11	3.13	124.98	118.66
5	D	3701	FAD	C1B-N9A-C8A	-3.12	120.17	127.09
5	B	1701	FAD	N3A-C2A-N1A	-3.10	123.89	128.58
5	A	701	FAD	O2B-C2B-C1B	3.10	120.77	110.10
5	B	1701	FAD	C6A-C5A-N7A	3.08	138.03	132.09
5	D	3701	FAD	C6A-C5A-N7A	3.07	138.00	132.09
5	C	2701	FAD	N3A-C2A-N1A	-3.04	123.98	128.58
7	B	698	P25	CM4-C4-C5	3.04	133.88	113.36
4	C	3695	1MM	C7'-C6'-N1'	3.01	121.63	117.13
4	A	695	1MM	C7'-C6'-N1'	3.00	121.62	117.13
6	B	1703	PYD	C2'-N3'-C4'	2.96	122.66	118.10
6	B	1703	PYD	C7'-C5'-C4'	2.96	124.85	121.74
6	A	703	PYD	C2'-N3'-C4'	2.93	122.61	118.10
5	A	701	FAD	C1B-N9A-C8A	-2.93	120.59	127.09
4	B	1695	1MM	C7'-C6'-N1'	2.92	121.50	117.13
6	A	703	PYD	C7'-C5'-C4'	2.90	124.79	121.74
4	C	2695	1MM	C7'-C6'-N1'	2.90	121.46	117.13
5	C	2701	FAD	C6A-C5A-N7A	2.87	137.62	132.09
7	A	1698	P25	C4-C5-C6	2.87	140.41	115.25
5	A	701	FAD	O2B-C2B-C3B	2.83	120.90	111.82
7	B	698	P25	C4-C5-C6	2.82	140.01	115.25
5	D	3701	FAD	O2B-C2B-C3B	2.75	120.62	111.82
5	B	1701	FAD	O2B-C2B-C3B	2.66	120.33	111.82
5	C	2701	FAD	O2-C2-N1	-2.64	117.42	121.80
5	B	1701	FAD	O2-C2-N1	-2.62	117.44	121.80
5	C	2701	FAD	O2B-C2B-C3B	2.62	120.21	111.82
5	A	701	FAD	O2-C2-N1	-2.58	117.52	121.80
4	B	1695	1MM	O7B-S7-N8	-2.57	99.42	106.77
6	D	3703	PYD	C6'-C5'-C4'	-2.50	112.17	115.49
5	D	3701	FAD	O2-C2-N1	-2.50	117.64	121.80
5	D	3701	FAD	O3'-C3'-C4'	-2.46	103.35	108.93
5	B	1701	FAD	O3'-C3'-C4'	-2.40	103.48	108.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	2695	1MM	O7A-S7-C2	2.39	111.63	107.68
5	B	1701	FAD	O4B-C4B-C5B	-2.39	101.69	109.33
4	C	3695	1MM	O7B-S7-N8	-2.38	99.96	106.77
5	A	701	FAD	O3'-C3'-C4'	-2.35	103.59	108.93
5	C	2701	FAD	O3'-C3'-C4'	-2.32	103.66	108.93
5	A	701	FAD	O4B-C4B-C5B	-2.27	102.06	109.33
5	D	3701	FAD	O4B-C4B-C5B	-2.24	102.15	109.33
6	C	2703	PYD	C6'-C5'-C4'	-2.22	112.55	115.49
4	C	2695	1MM	O7B-S7-N8	-2.21	100.44	106.77
4	C	2695	1MM	C7'-C6'-N5'	2.20	120.42	117.13
5	A	701	FAD	C1'-C2'-C3'	2.20	115.61	109.66
5	C	2701	FAD	O4B-C4B-C5B	-2.14	102.47	109.33
5	C	2701	FAD	O4-C4-C4X	-2.12	120.94	126.53
4	A	695	1MM	O7B-S7-N8	-2.11	100.73	106.77
5	D	3701	FAD	O4-C4-C4X	-2.11	120.97	126.53
5	C	2701	FAD	C1'-C2'-C3'	2.10	115.36	109.66
4	A	695	1MM	C7'-C6'-N5'	2.07	120.23	117.13
4	B	1695	1MM	C7'-C6'-N5'	2.06	120.22	117.13
5	B	1701	FAD	C1'-C2'-C3'	2.06	115.25	109.66
5	B	1701	FAD	O4-C4-C4X	-2.06	121.10	126.53
5	A	701	FAD	O4-C4-C4X	-2.05	121.12	126.53
8	C	3702	P22	O3A-PB-O2B	-2.04	100.30	111.04
5	D	3701	FAD	O3'-C3'-C2'	-2.02	104.34	108.93
6	B	1703	PYD	C6'-C5'-C4'	-2.01	112.82	115.49
6	A	703	PYD	C6'-C5'-C4'	-2.01	112.83	115.49
5	D	3701	FAD	C1'-C2'-C3'	2.00	115.09	109.66
5	B	1701	FAD	O2'-C2'-C3'	-2.00	104.56	109.25

There are no chirality outliers.

All (56) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1698	P25	C6-C7-O7-PA
8	D	2702	P22	C6-C7-O7-PA
4	A	695	1MM	C1-C2-S7-O7A
4	B	1695	1MM	C1-C2-S7-O7A
4	C	2695	1MM	C1-C2-S7-O7A
4	C	3695	1MM	C1-C2-S7-O7A
7	A	1698	P25	C4-C5-C6-C7
4	B	1695	1MM	C3-C2-S7-N8
4	A	695	1MM	C1-C2-S7-O7B
4	B	1695	1MM	C1-C2-S7-O7B

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Mol	Chain	Res	Type	Atoms
4	C	2695	1MM	C1-C2-S7-O7B
4	C	3695	1MM	C1-C2-S7-O7B
4	A	695	1MM	C3-C2-S7-N8
4	A	695	1MM	C1-C2-S7-N8
4	B	1695	1MM	C1-C2-S7-N8
4	C	2695	1MM	C1-C2-S7-N8
4	C	3695	1MM	C1-C2-S7-N8
8	C	3702	P22	PA-O3A-PB-O2B
8	D	2702	P22	PA-O3A-PB-O2B
4	C	2695	1MM	C3-C2-S7-N8
4	C	3695	1MM	C3-C2-S7-N8
4	C	2695	1MM	C6-C1-C11-O12
4	A	695	1MM	C3-C2-S7-O7B
4	B	1695	1MM	C3-C2-S7-O7B
4	C	3695	1MM	C3-C2-S7-O7B
4	B	1695	1MM	C6-C1-C11-O12
4	A	695	1MM	C3-C2-S7-O7A
4	C	2695	1MM	C3-C2-S7-O7B
4	A	695	1MM	C6-C1-C11-O12
4	C	2695	1MM	C2-C1-C11-O12
4	C	3695	1MM	C6-C1-C11-O12
4	B	1695	1MM	C3-C2-S7-O7A
4	A	695	1MM	C2-C1-C11-O12
4	B	1695	1MM	C2-C1-C11-O12
4	C	2695	1MM	C3-C2-S7-O7A
4	C	3695	1MM	C3-C2-S7-O7A
4	A	695	1MM	C2-C1-C11-O11
4	C	2695	1MM	C2-C1-C11-O11
5	A	701	FAD	O4B-C4B-C5B-O5B
5	C	2701	FAD	O4B-C4B-C5B-O5B
5	D	3701	FAD	O4B-C4B-C5B-O5B
4	A	695	1MM	N8-C9-N10-C2'
4	B	1695	1MM	N8-C9-N10-C2'
4	C	2695	1MM	N8-C9-N10-C2'
4	C	3695	1MM	N8-C9-N10-C2'
4	C	2695	1MM	C6-C1-C11-O11
4	B	1695	1MM	C2-C1-C11-O11
5	A	701	FAD	C2B-C1B-N9A-C8A
5	B	1701	FAD	C2B-C1B-N9A-C8A
5	B	1701	FAD	O4B-C4B-C5B-O5B
4	B	1695	1MM	C6-C1-C11-O11
4	C	3695	1MM	C2-C1-C11-O12

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Mol	Chain	Res	Type	Atoms
4	A	695	1MM	O9-C9-N10-C2'
4	B	1695	1MM	O9-C9-N10-C2'
4	C	2695	1MM	O9-C9-N10-C2'
4	C	3695	1MM	O9-C9-N10-C2'

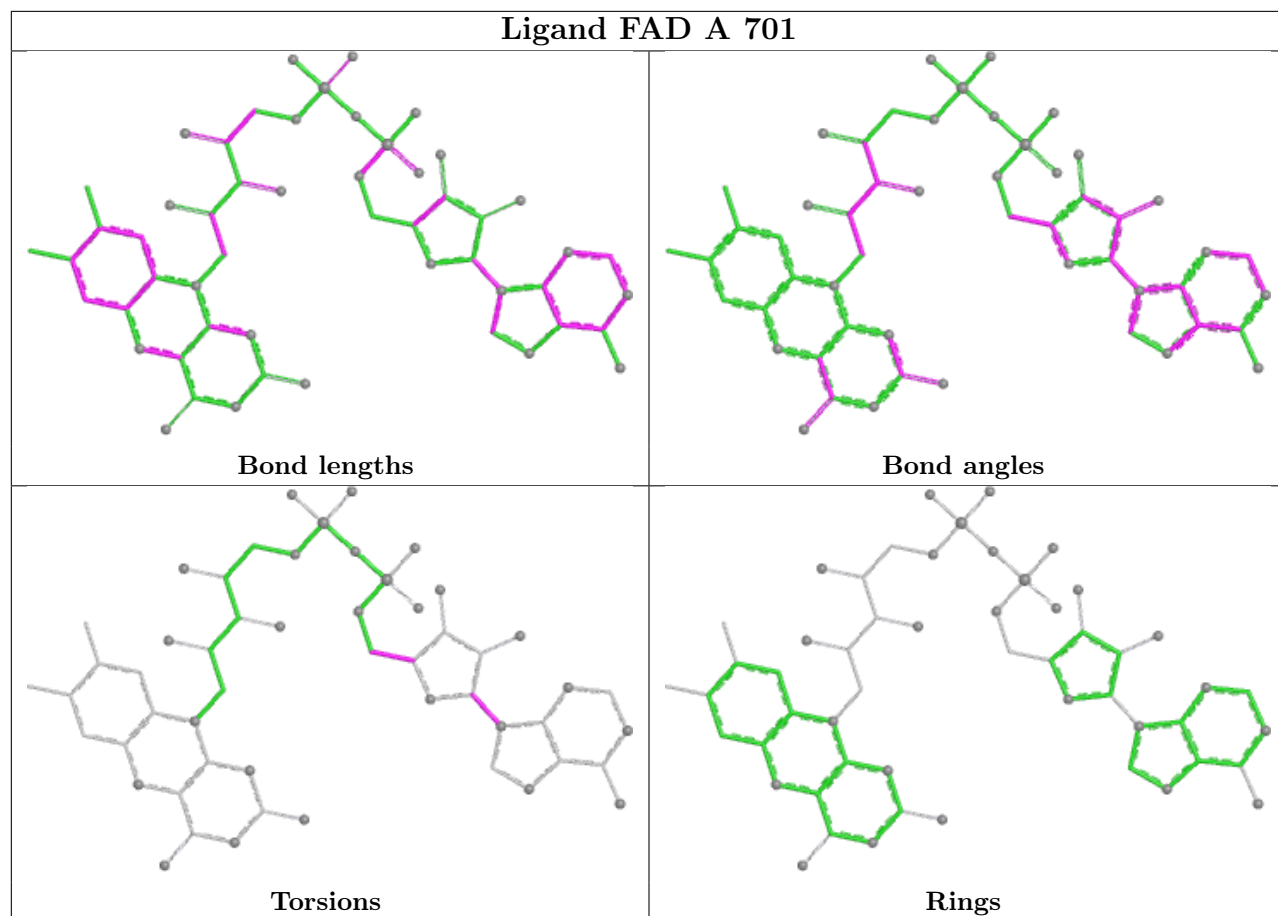
There are no ring outliers.

9 monomers are involved in 13 short contacts:

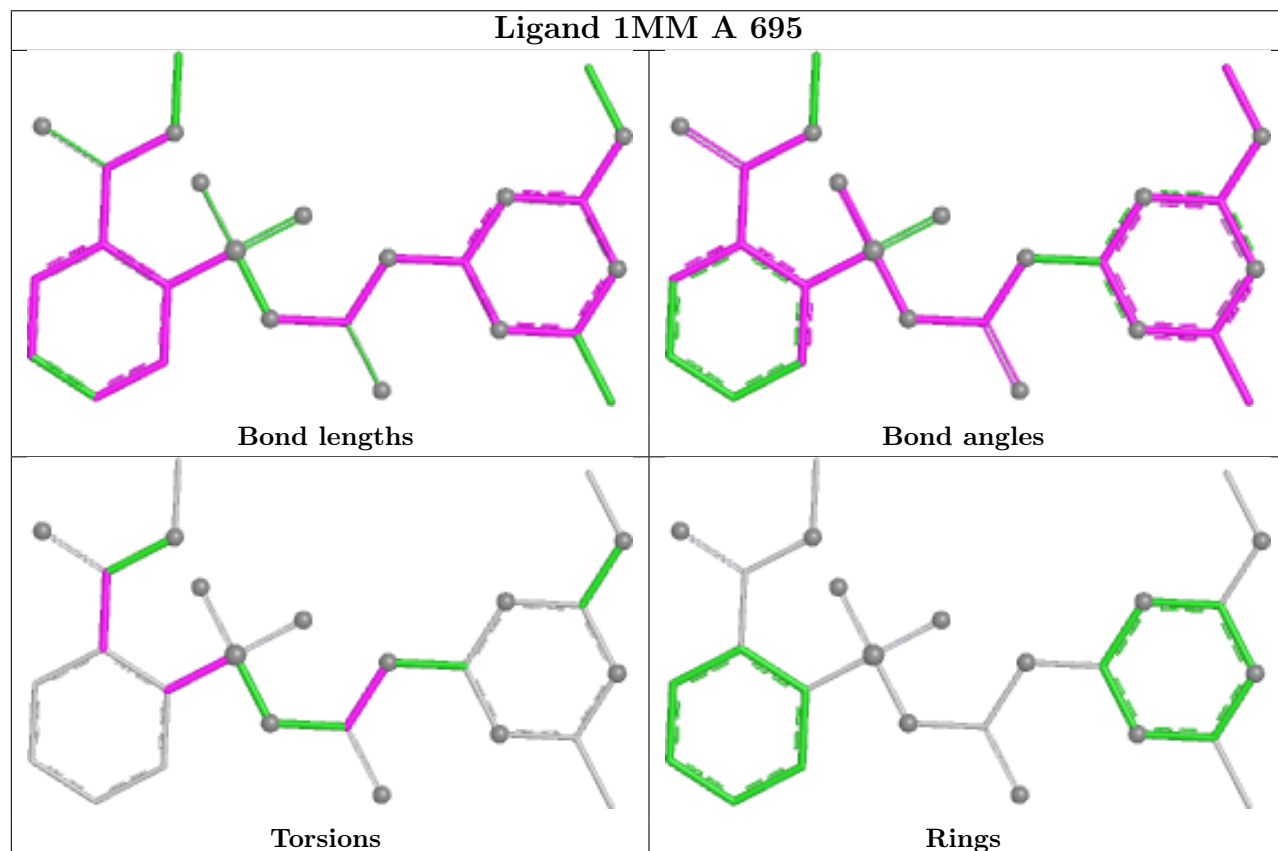
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	703	PYD	1	0
7	B	698	P25	1	0
4	A	695	1MM	1	0
5	B	1701	FAD	1	0
4	B	1695	1MM	4	0
8	D	2702	P22	1	0
4	C	3695	1MM	1	0
4	C	2695	1MM	1	0
7	A	1698	P25	2	0

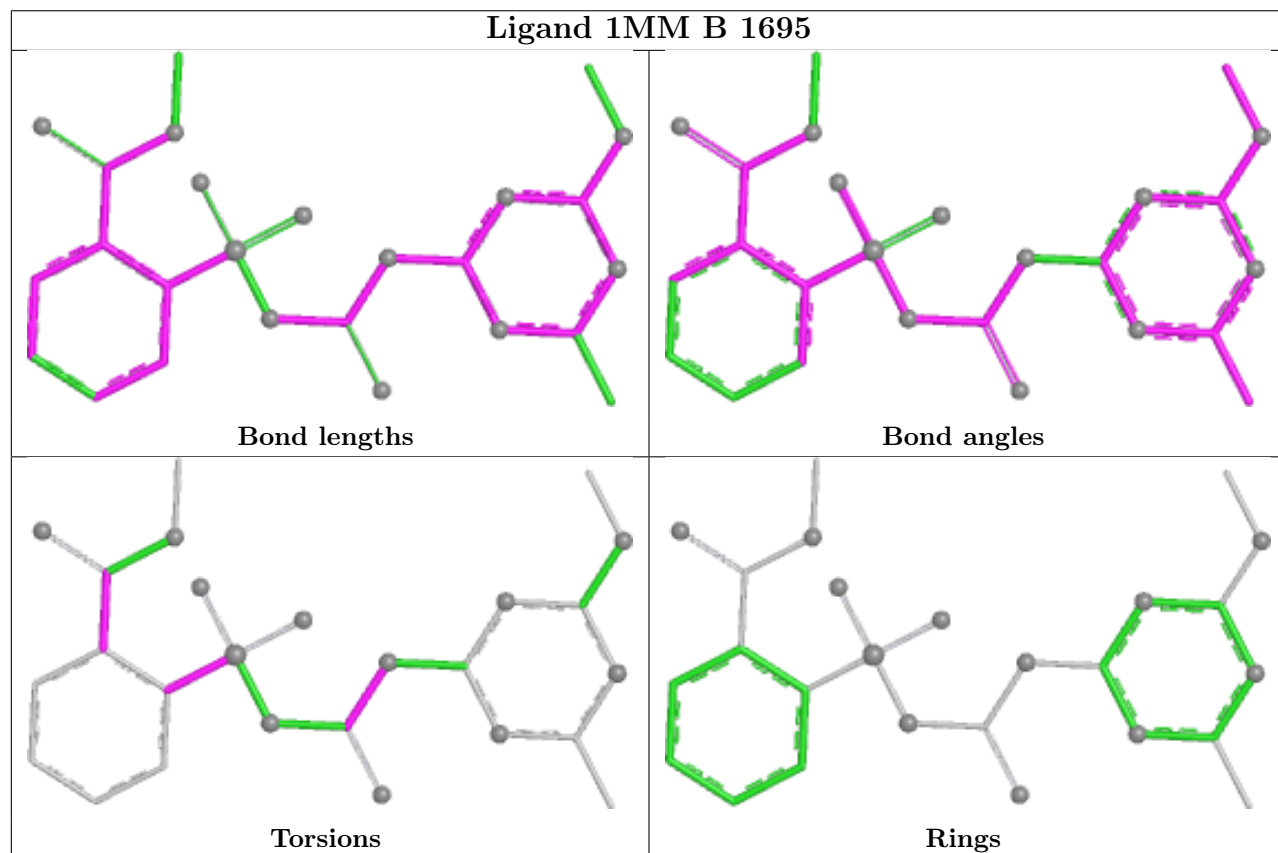
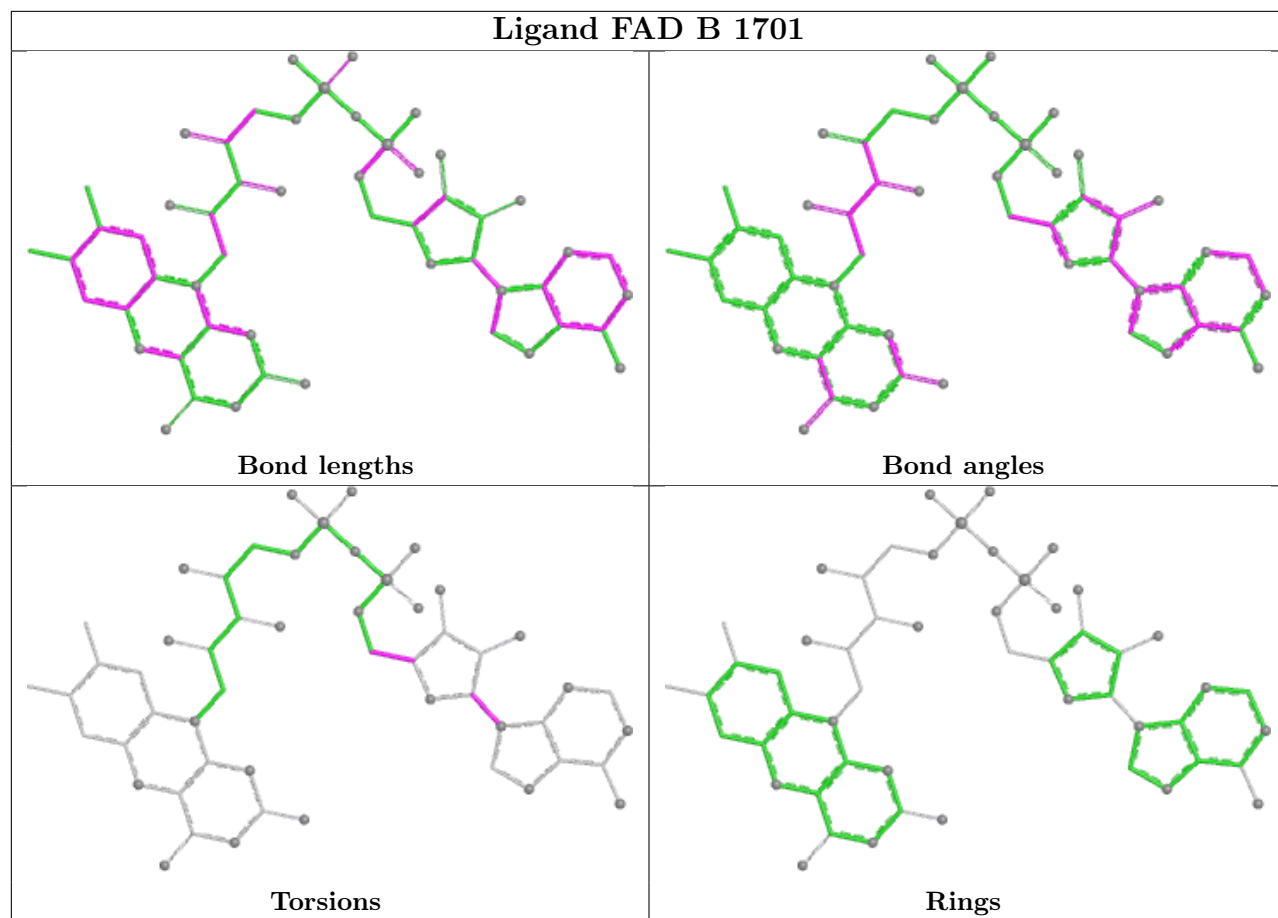
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

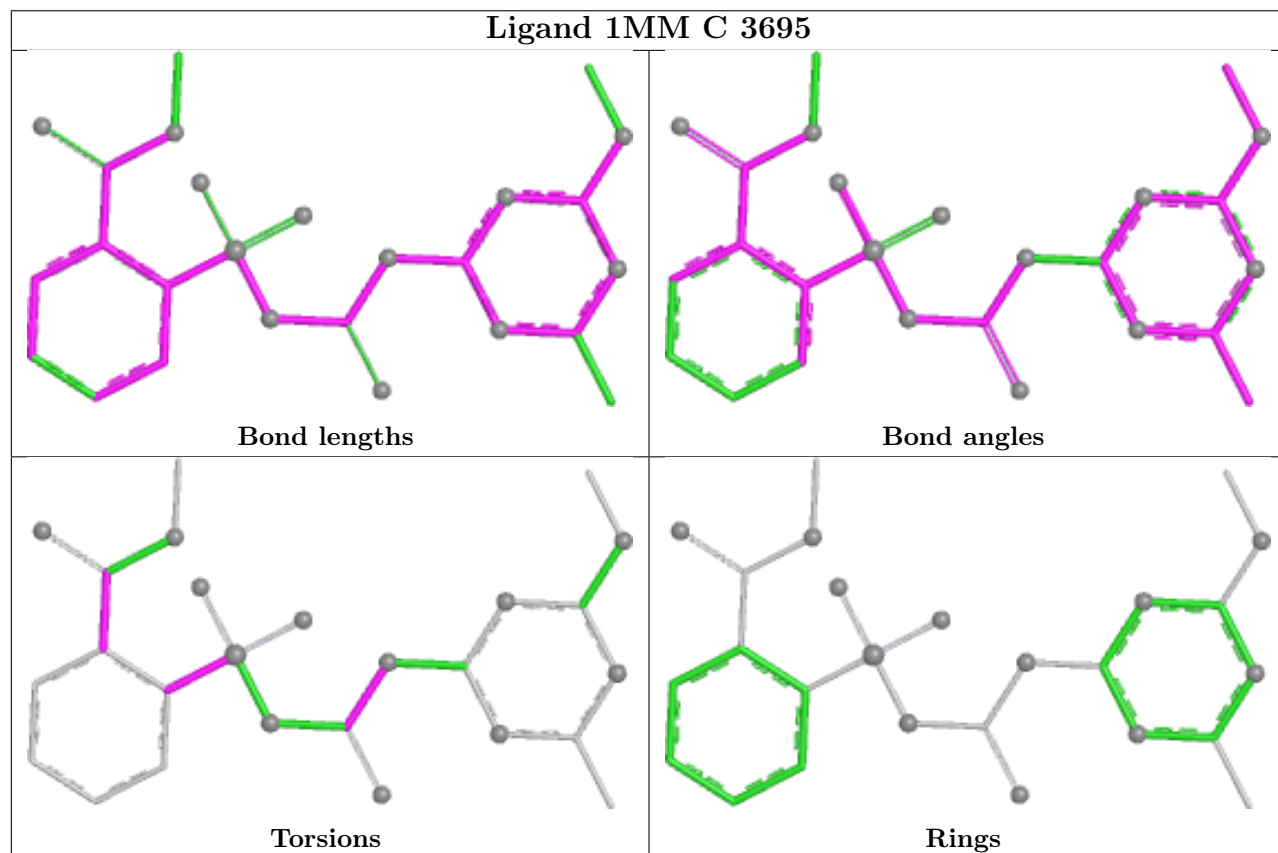
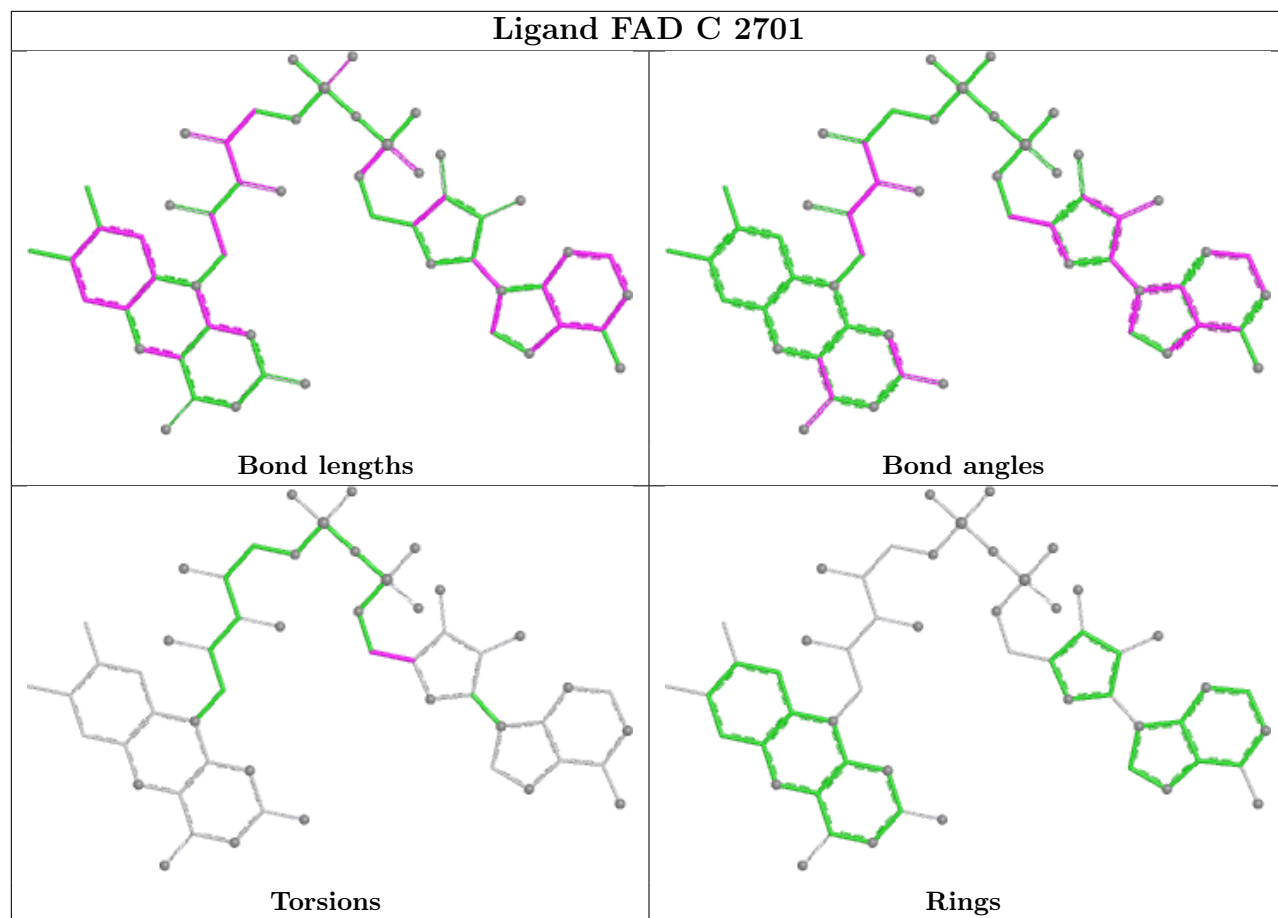
Ligand FAD A 701



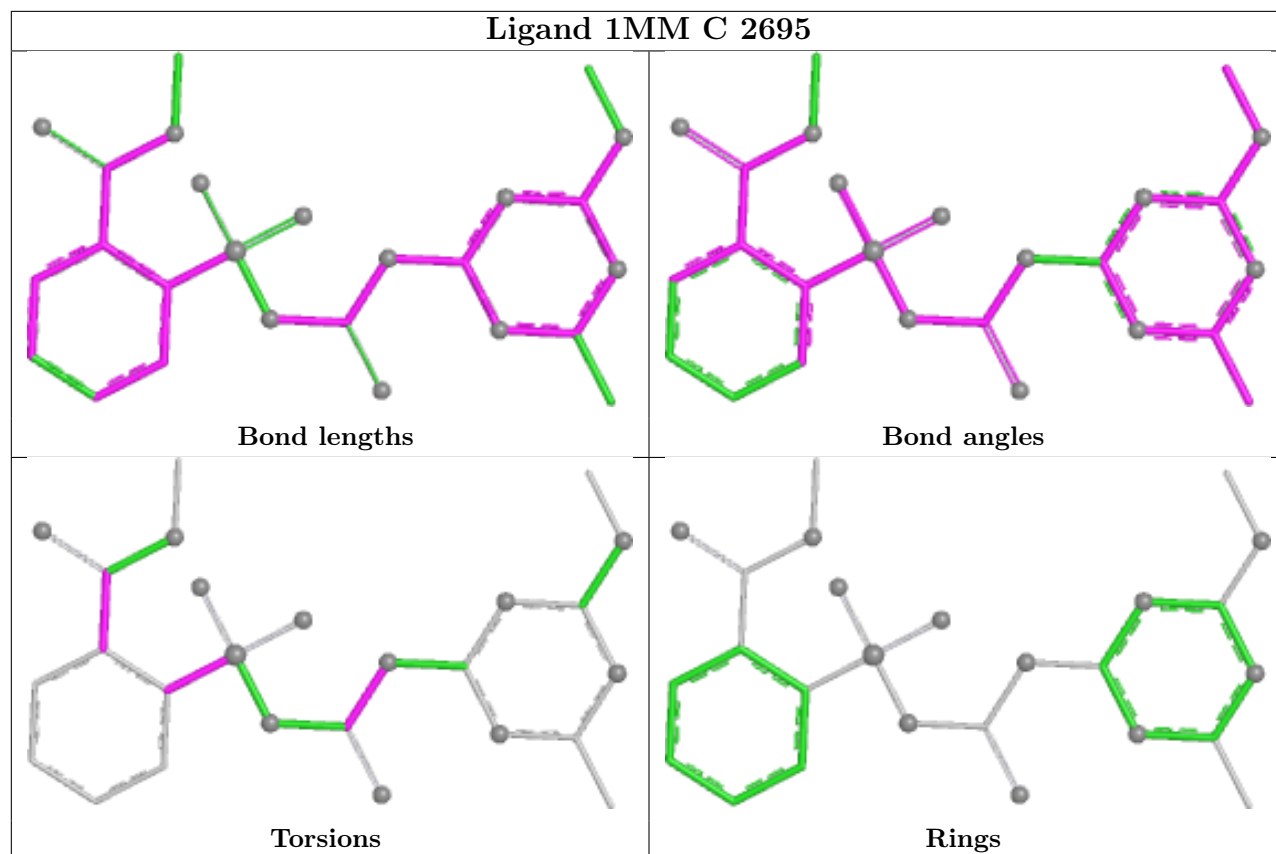
Ligand 1MM A 695



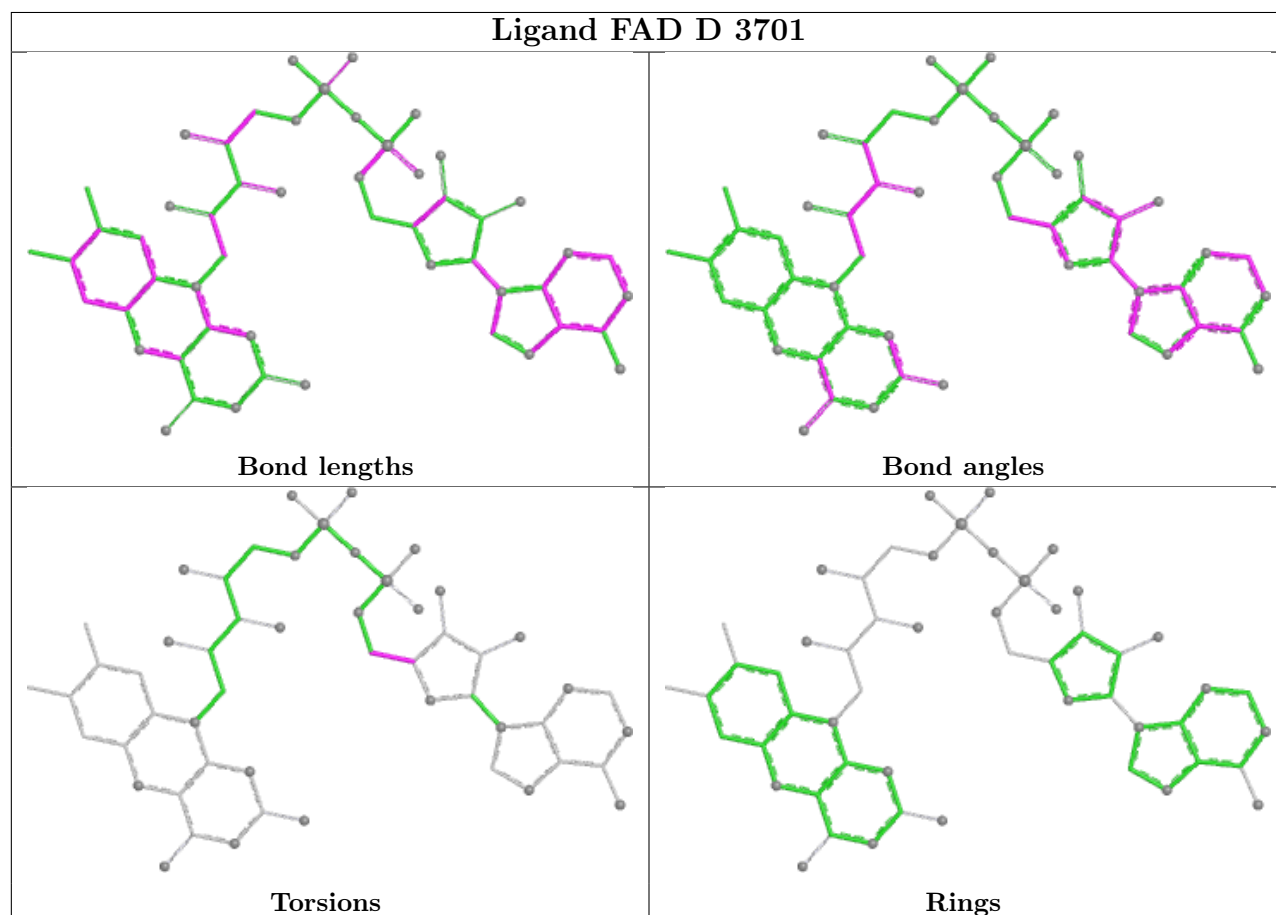




Ligand 1MM C 2695



Ligand FAD D 3701



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	596/677 (88%)	-0.54	4 (0%) 84 85	15, 25, 41, 76	6 (1%)
1	B	582/677 (85%)	0.02	12 (2%) 63 65	19, 38, 62, 84	3 (0%)
1	C	575/677 (84%)	0.68	75 (13%) 7 8	18, 49, 99, 114	2 (0%)
1	D	596/677 (88%)	-0.43	5 (0%) 82 83	16, 28, 43, 74	5 (0%)
All	All	2349/2708 (86%)	-0.07	96 (4%) 41 43	15, 32, 77, 114	16 (0%)

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	132[A]	ASN	7.6
1	A	277	THR	7.0
1	C	293	ALA	6.3
1	C	291	LYS	5.5
1	C	438	ILE	5.1
1	C	298	LEU	5.0
1	C	402	GLY	4.8
1	C	403	ILE	4.8
1	B	284	PHE	4.5
1	C	441	VAL	4.2
1	C	303	VAL	4.2
1	C	264	THR	4.2
1	C	446	GLU	4.0
1	C	440	PRO	4.0
1	C	437	LYS	4.0
1	C	295	LEU	4.0
1	C	290	ASN	3.9
1	C	447	TRP	3.8
1	C	439	PHE	3.7
1	C	454	TRP	3.6
1	C	306	VAL	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	445	SER	3.5
1	C	304	LEU	3.5
1	C	665	ILE	3.4
1	C	443	GLU	3.4
1	C	323	LEU	3.4
1	C	370	ILE	3.4
1	A	268	LEU	3.3
1	C	296	ILE	3.3
1	B	87	THR	3.3
1	C	262	ILE	3.3
1	C	265	LYS	3.3
1	B	263	PRO	3.3
1	D	278	SER	3.2
1	C	458	TYR	3.2
1	D	129	ASP	3.1
1	C	448	PHE	3.1
1	C	386	SER	3.0
1	C	383	GLY	3.0
1	C	444	ARG	2.9
1	C	442	LYS	2.9
1	C	434	MET	2.9
1	C	397	ALA	2.9
1	C	431	LEU	2.9
1	C	364	VAL	2.9
1	D	270	SER	2.8
1	C	422	ALA	2.8
1	C	371	ILE	2.8
1	C	292	ALA	2.8
1	C	666	ASN	2.7
1	C	314	ALA	2.7
1	C	406	PHE	2.7
1	B	434	MET	2.7
1	C	362	LEU	2.7
1	C	327	ALA	2.6
1	C	326	ARG	2.6
1	B	291	LYS	2.6
1	C	263	PRO	2.6
1	C	404	ILE	2.6
1	C	311	LEU	2.6
1	A	265	LYS	2.6
1	C	319	LEU	2.5
1	C	393	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	294	ASP	2.5
1	C	449	ALA	2.5
1	C	363	ALA	2.4
1	C	417	VAL	2.4
1	C	667	PHE	2.4
1	C	433	LYS	2.4
1	B	289	ILE	2.4
1	C	320	LEU	2.4
1	B	86	ASP	2.3
1	C	385	ILE	2.3
1	C	661	LEU	2.2
1	C	660	GLY	2.2
1	A	278	SER	2.2
1	C	423	VAL	2.2
1	C	350	ASP	2.2
1	C	430	ASN	2.2
1	C	432	GLY	2.2
1	C	372	ALA	2.2
1	B	298	LEU	2.1
1	C	329	ILE	2.1
1	C	351	MET	2.1
1	C	592	GLU	2.1
1	C	367	ALA	2.1
1	C	399	GLY	2.1
1	B	285	VAL	2.1
1	B	446	GLU	2.1
1	C	669	PRO	2.1
1	D	265	LYS	2.0
1	B	261	PRO	2.0
1	C	331	VAL	2.0
1	C	332	THR	2.0
1	B	580	GLN	2.0
1	C	328	GLN	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

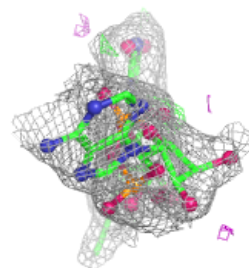
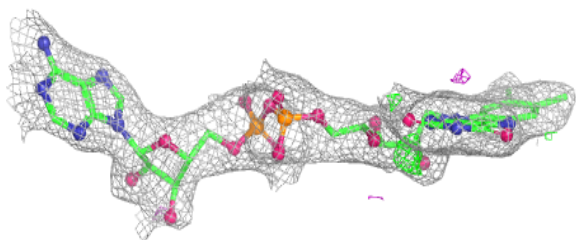
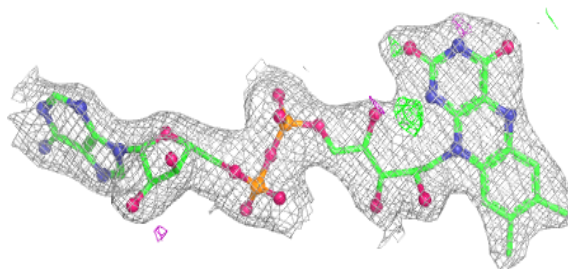
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	PYD	D	3703	9/9	0.83	0.20	43,46,49,51	0
6	PYD	B	1703	9/9	0.85	0.17	35,37,38,38	0
6	PYD	C	2703	9/9	0.86	0.16	39,40,42,44	0
6	PYD	A	703	9/9	0.89	0.16	33,37,39,39	0
5	FAD	C	2701	53/53	0.95	0.09	48,51,55,57	0
4	1MM	C	3695	26/26	0.96	0.10	48,51,53,56	0
7	P25	B	698	14/14	0.96	0.12	25,29,42,44	0
4	1MM	B	1695	26/26	0.97	0.07	30,35,38,39	0
4	1MM	C	2695	26/26	0.97	0.08	27,30,31,34	0
2	K	C	3696	1/1	0.97	0.06	61,61,61,61	0
5	FAD	B	1701	53/53	0.97	0.07	30,33,36,37	0
7	P25	A	1698	14/14	0.97	0.09	18,22,33,36	0
4	1MM	A	695	26/26	0.97	0.08	21,25,30,30	0
8	P22	C	3702	11/11	0.97	0.08	38,40,44,45	0
5	FAD	A	701	53/53	0.98	0.05	13,18,23,24	0
3	MG	C	3699	1/1	0.99	0.07	36,36,36,36	0
3	MG	D	2699	1/1	0.99	0.03	24,24,24,24	0
2	K	B	696	1/1	0.99	0.04	45,45,45,45	0
2	K	D	2696	1/1	0.99	0.03	26,26,26,26	0
5	FAD	D	3701	53/53	0.99	0.05	17,21,24,25	0
3	MG	B	699	1/1	0.99	0.03	25,25,25,25	0
8	P22	D	2702	11/11	0.99	0.06	21,24,27,30	0
3	MG	A	1699	1/1	1.00	0.02	19,19,19,19	0
2	K	A	1696	1/1	1.00	0.02	24,24,24,24	0

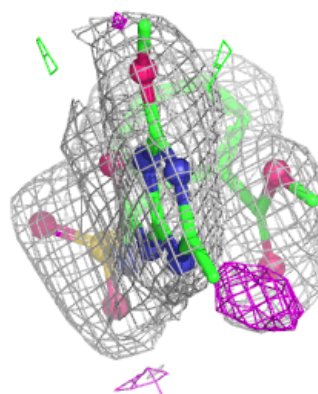
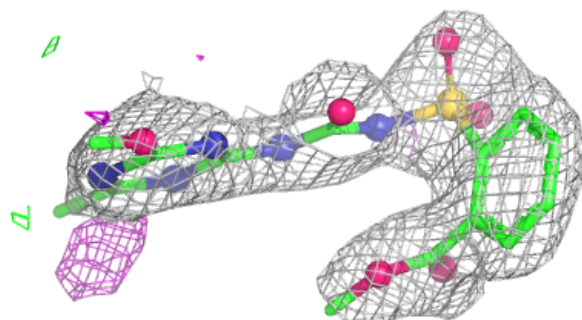
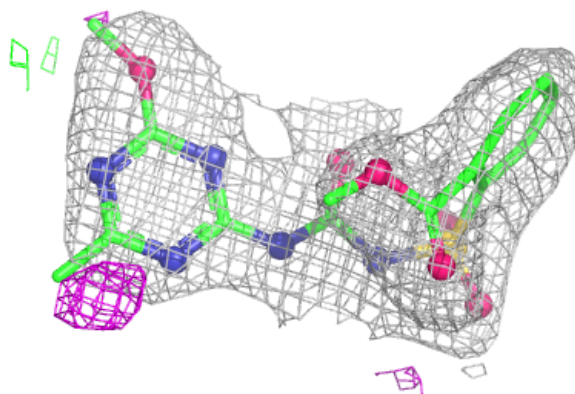
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD C 2701:

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

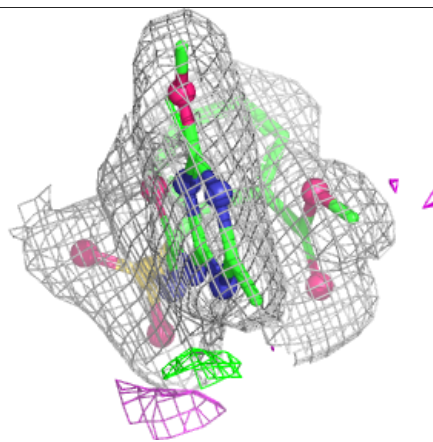
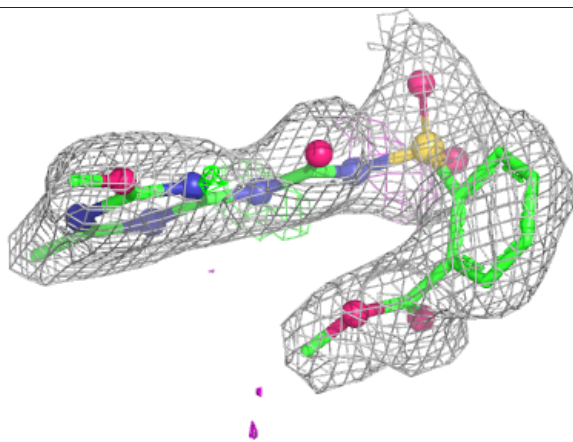
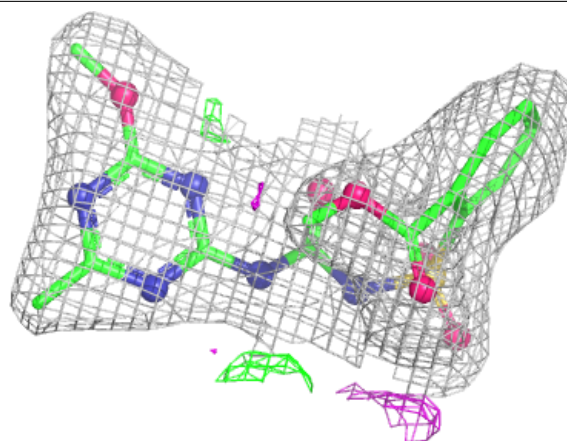
**Electron density around 1MM C 3695:**

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



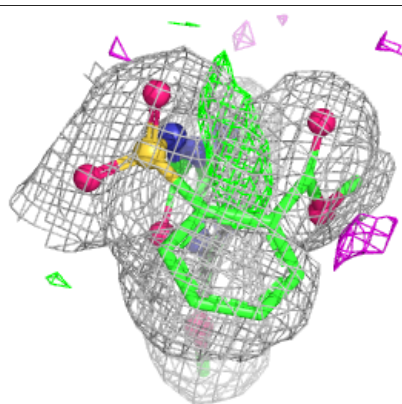
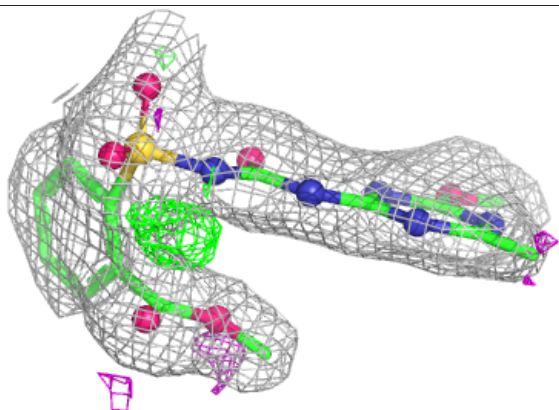
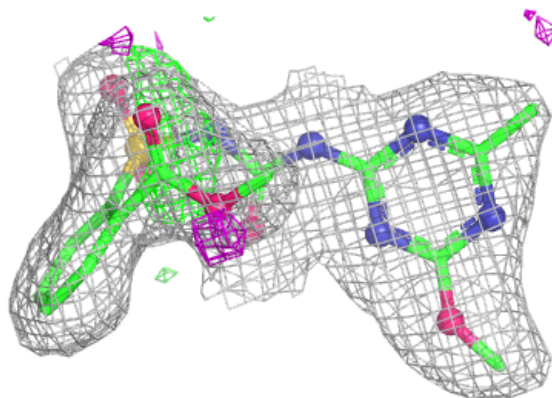
Electron density around 1MM B 1695:

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

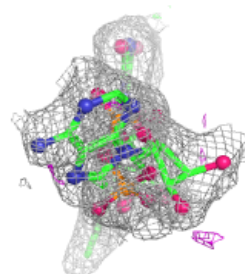
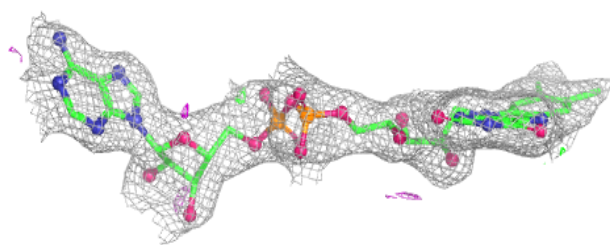
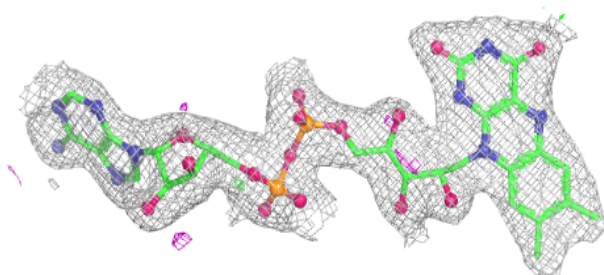


Electron density around 1MM C 2695:

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and green (positive)

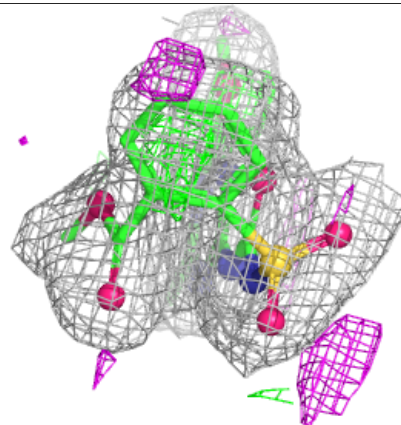
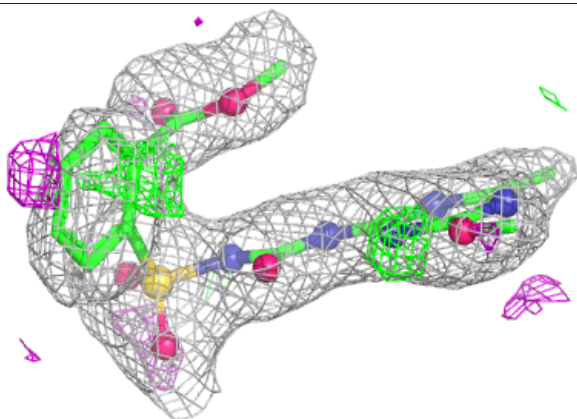
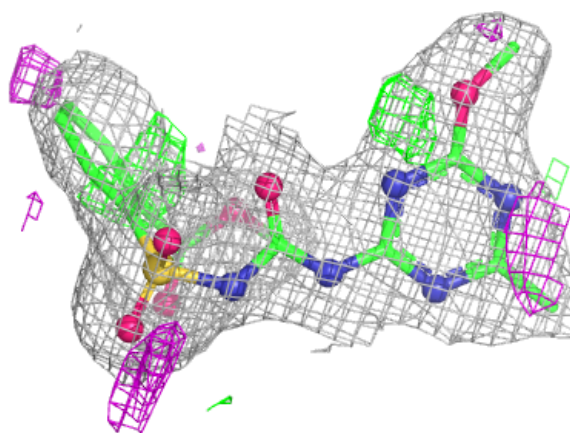
**Electron density around FAD B 1701:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

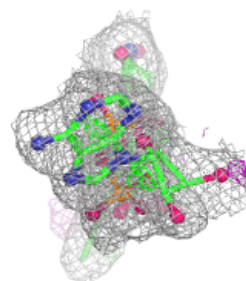
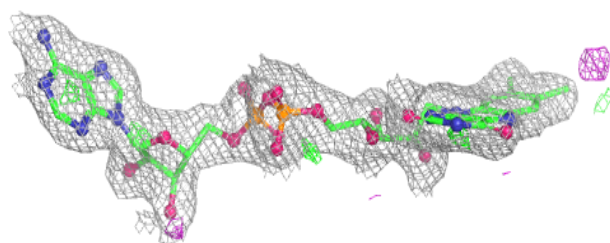
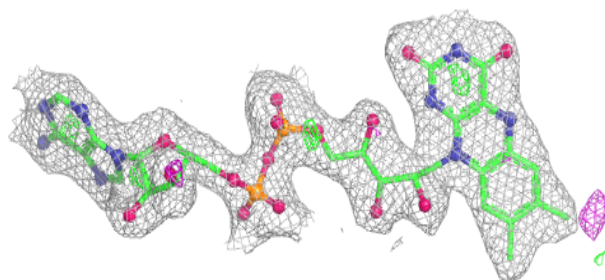


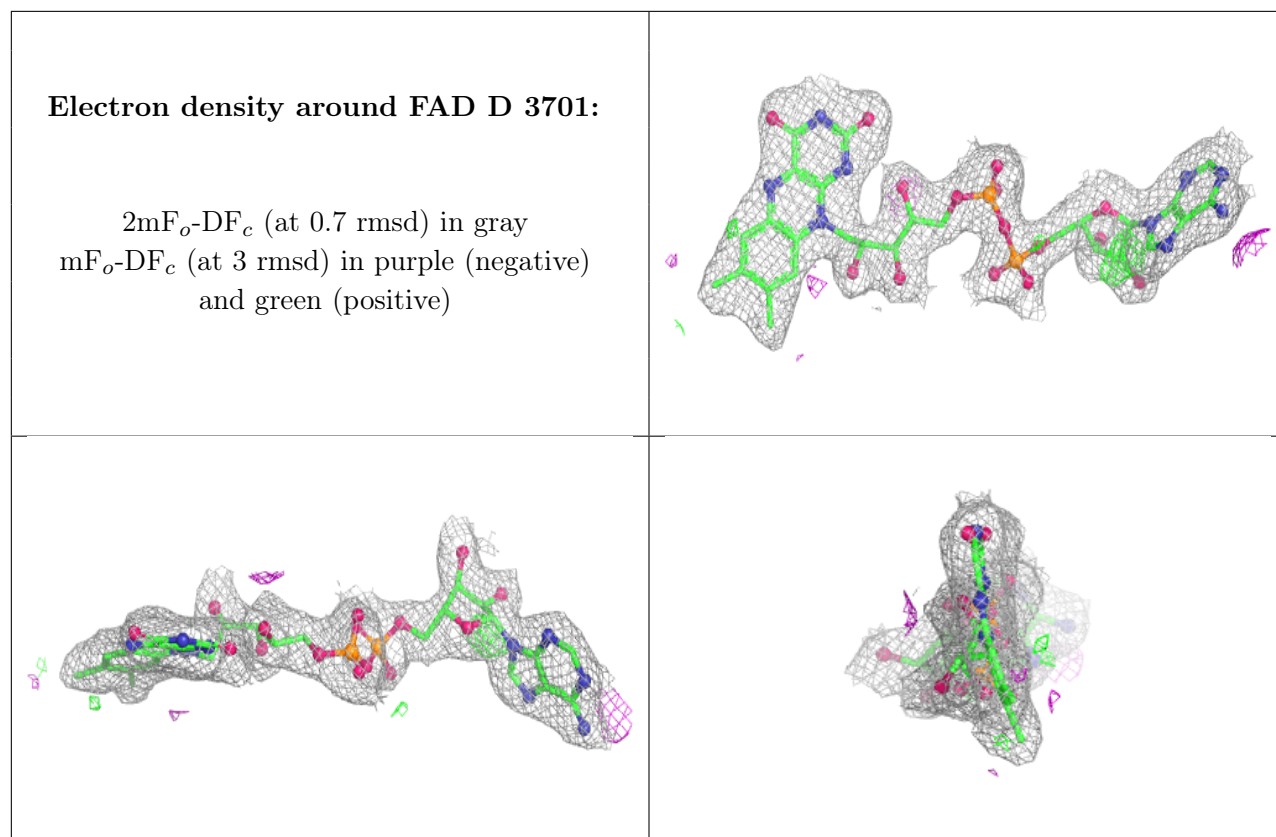
Electron density around 1MM A 695:

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

**Electron density around FAD A 701:**

$2mF_o-DF_c$ (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.