



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:20 PM UTC

PDB ID : 8TBT / pdb_00008tbt
Title : Structure of human erythrocyte pyruvate kinase in complex with an allosteric activator Compound 2
Authors : Jin, L.; Padyana, A.
Deposited on : 2023-06-29
Resolution : 2.34 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

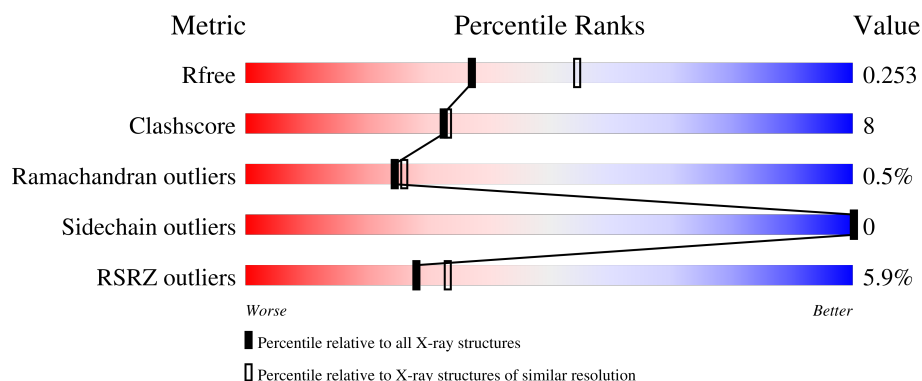
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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	544	5% 82% 13% 6%
1	B	544	3% 81% 14% 5%
1	C	544	8% 73% 19% 8%
1	D	544	6% 81% 13% 6%

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
3	PYR	A	602	-	X	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 16072 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 Pyruvate kinase PKLR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	514	Total	C	N	O	S	0	0	0
			3894	2446	705	725	18			
1	B	518	Total	C	N	O	S	0	0	0
			3918	2460	710	730	18			
1	C	503	Total	C	N	O	S	0	0	0
			3806	2391	688	709	18			
1	D	512	Total	C	N	O	S	0	0	0
			3877	2435	702	722	18			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	31	MET	-	initiating methionine	UNP P30613
A	32	GLY	-	expression tag	UNP P30613
A	33	SER	-	expression tag	UNP P30613
A	34	SER	-	expression tag	UNP P30613
A	35	HIS	-	expression tag	UNP P30613
A	36	HIS	-	expression tag	UNP P30613
A	37	HIS	-	expression tag	UNP P30613
A	38	HIS	-	expression tag	UNP P30613
A	39	HIS	-	expression tag	UNP P30613
A	40	HIS	-	expression tag	UNP P30613
A	41	SER	-	expression tag	UNP P30613
A	42	SER	-	expression tag	UNP P30613
A	43	GLY	-	expression tag	UNP P30613
A	44	LEU	-	expression tag	UNP P30613
A	45	VAL	-	expression tag	UNP P30613
A	46	PRO	-	expression tag	UNP P30613
A	47	ARG	-	expression tag	UNP P30613
A	48	GLY	-	expression tag	UNP P30613
A	49	SER	-	expression tag	UNP P30613
B	31	MET	-	initiating methionine	UNP P30613
B	32	GLY	-	expression tag	UNP P30613

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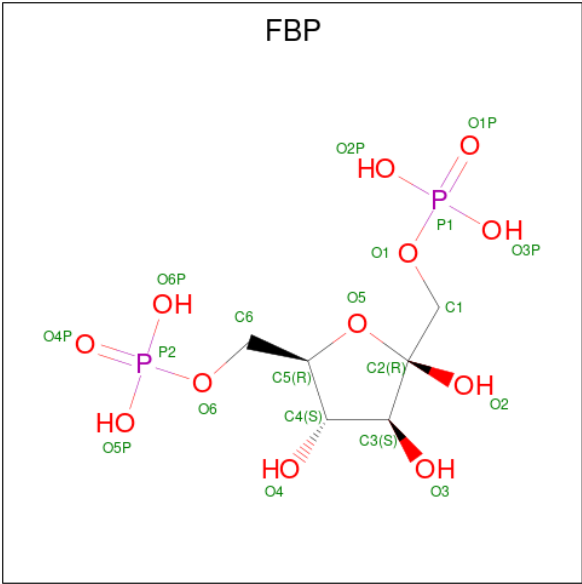
Chain	Residue	Modelled	Actual	Comment	Reference
B	33	SER	-	expression tag	UNP P30613
B	34	SER	-	expression tag	UNP P30613
B	35	HIS	-	expression tag	UNP P30613
B	36	HIS	-	expression tag	UNP P30613
B	37	HIS	-	expression tag	UNP P30613
B	38	HIS	-	expression tag	UNP P30613
B	39	HIS	-	expression tag	UNP P30613
B	40	HIS	-	expression tag	UNP P30613
B	41	SER	-	expression tag	UNP P30613
B	42	SER	-	expression tag	UNP P30613
B	43	GLY	-	expression tag	UNP P30613
B	44	LEU	-	expression tag	UNP P30613
B	45	VAL	-	expression tag	UNP P30613
B	46	PRO	-	expression tag	UNP P30613
B	47	ARG	-	expression tag	UNP P30613
B	48	GLY	-	expression tag	UNP P30613
B	49	SER	-	expression tag	UNP P30613
C	31	MET	-	initiating methionine	UNP P30613
C	32	GLY	-	expression tag	UNP P30613
C	33	SER	-	expression tag	UNP P30613
C	34	SER	-	expression tag	UNP P30613
C	35	HIS	-	expression tag	UNP P30613
C	36	HIS	-	expression tag	UNP P30613
C	37	HIS	-	expression tag	UNP P30613
C	38	HIS	-	expression tag	UNP P30613
C	39	HIS	-	expression tag	UNP P30613
C	40	HIS	-	expression tag	UNP P30613
C	41	SER	-	expression tag	UNP P30613
C	42	SER	-	expression tag	UNP P30613
C	43	GLY	-	expression tag	UNP P30613
C	44	LEU	-	expression tag	UNP P30613
C	45	VAL	-	expression tag	UNP P30613
C	46	PRO	-	expression tag	UNP P30613
C	47	ARG	-	expression tag	UNP P30613
C	48	GLY	-	expression tag	UNP P30613
C	49	SER	-	expression tag	UNP P30613
D	31	MET	-	initiating methionine	UNP P30613
D	32	GLY	-	expression tag	UNP P30613
D	33	SER	-	expression tag	UNP P30613
D	34	SER	-	expression tag	UNP P30613
D	35	HIS	-	expression tag	UNP P30613
D	36	HIS	-	expression tag	UNP P30613

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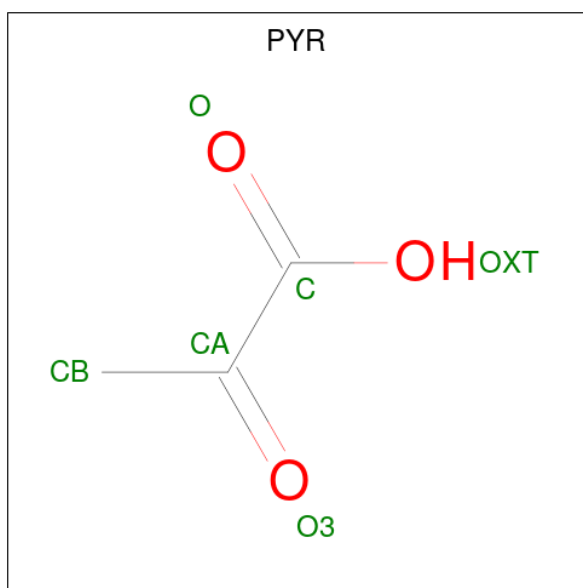
Chain	Residue	Modelled	Actual	Comment	Reference
D	37	HIS	-	expression tag	UNP P30613
D	38	HIS	-	expression tag	UNP P30613
D	39	HIS	-	expression tag	UNP P30613
D	40	HIS	-	expression tag	UNP P30613
D	41	SER	-	expression tag	UNP P30613
D	42	SER	-	expression tag	UNP P30613
D	43	GLY	-	expression tag	UNP P30613
D	44	LEU	-	expression tag	UNP P30613
D	45	VAL	-	expression tag	UNP P30613
D	46	PRO	-	expression tag	UNP P30613
D	47	ARG	-	expression tag	UNP P30613
D	48	GLY	-	expression tag	UNP P30613
D	49	SER	-	expression tag	UNP P30613

- Molecule 2 is 1,6-di-O-phosphono-beta-D-fructofuranose (CCD ID: FBP) (formula: $C_6H_{14}O_{12}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			20	6	12	2		
2	B	1	Total	C	O	P	0	0
			20	6	12	2		
2	C	1	Total	C	O	P	0	0
			20	6	12	2		
2	D	1	Total	C	O	P	0	0
			20	6	12	2		

- Molecule 3 is PYRUVIC ACID (CCD ID: PYR) (formula: $C_3H_4O_3$).



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

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

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

- Molecule 5 is POTASSIUM ION (CCD ID: K) (formula: K).

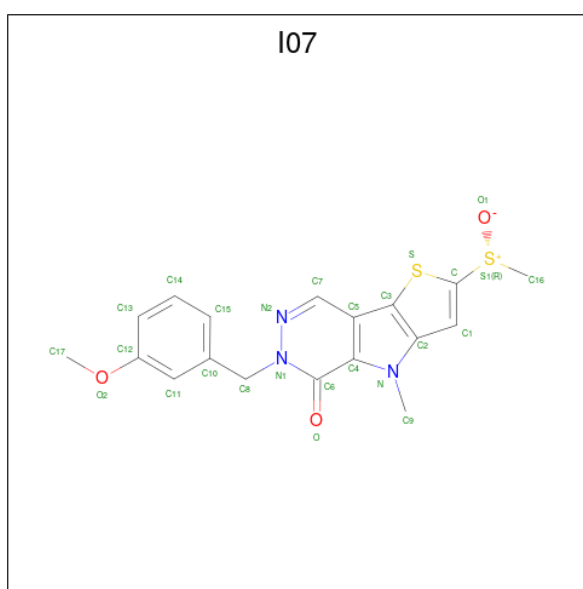
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	K	0	0
			1	1		

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

- Molecule 6 is {6-[(3-methoxyphenyl)methyl]-4-methyl-5-oxo-5,6-dihydro-4H-thieno[2',3':4,5]pyrrolo[2,3-d]pyridazin-2-yl}(methyl)sulfaniumolate (CCD ID: I07) (formula: C₁₈H₁₇N₃O₃S₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total C N O S 52 36 6 6 4	0	1
6	D	1	Total C N O S 52 36 6 6 4	0	1

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	91	Total O 91 91	0	0
7	B	92	Total O 92 92	0	0
7	C	70	Total O 70 70	0	0

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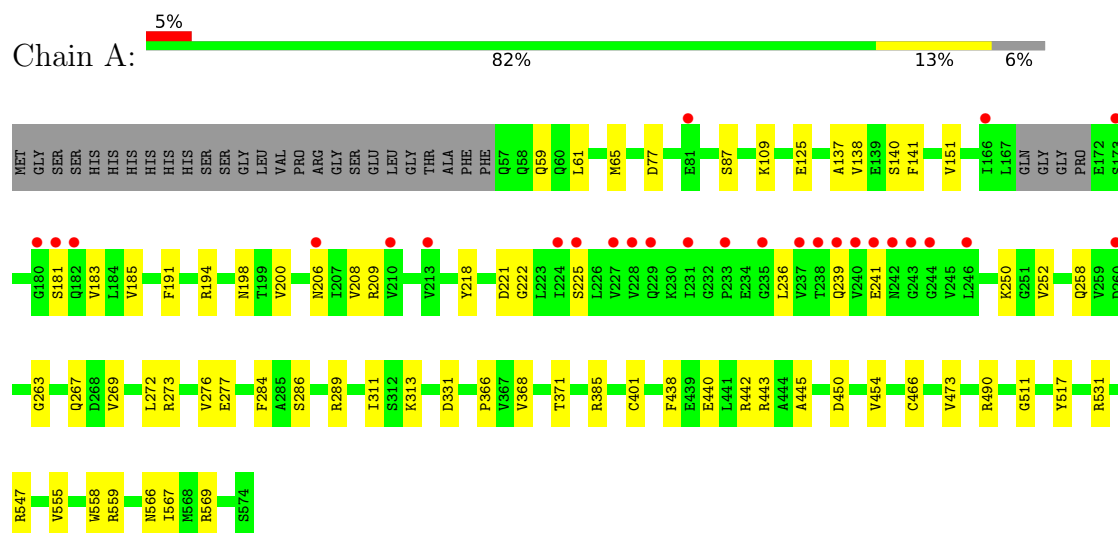
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	108	Total 108	O 108	0	0

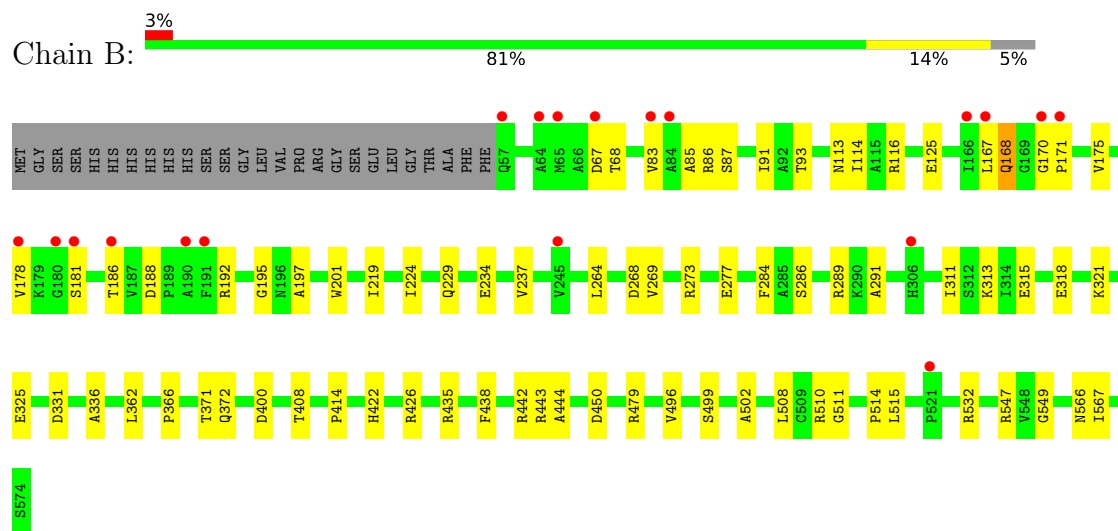
3 Residue-property plots [i](#)

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: Pyruvate kinase PKLR

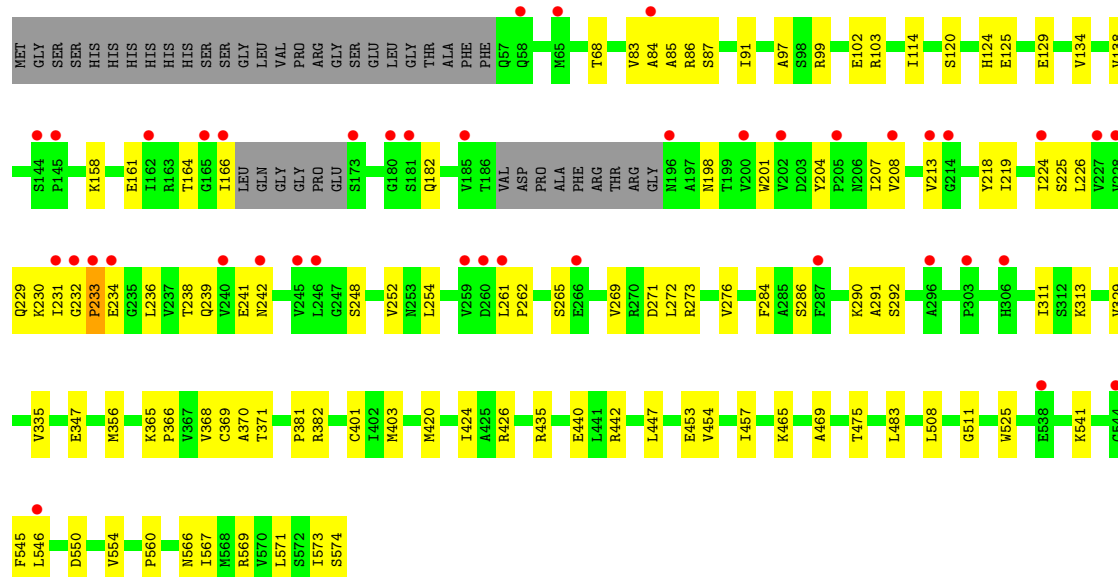


• Molecule 1: Pyruvate kinase PKLR

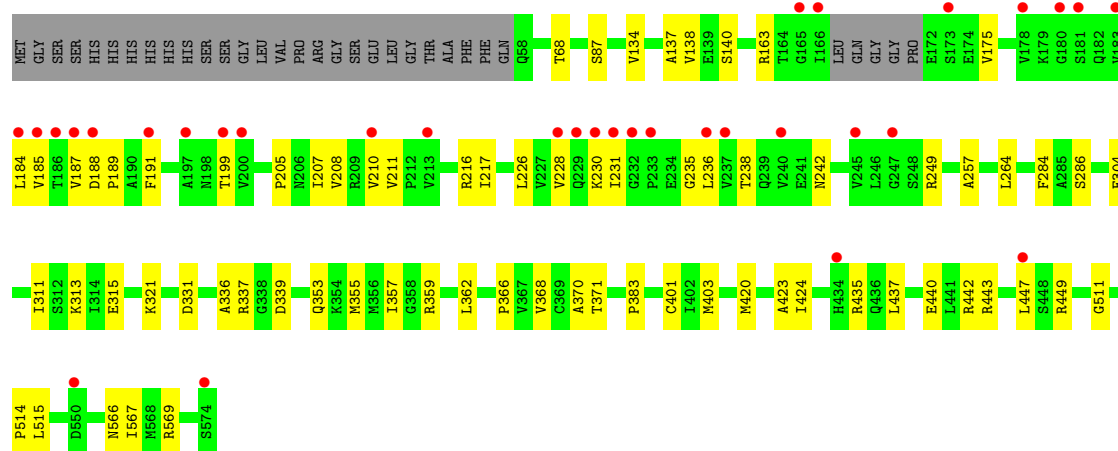
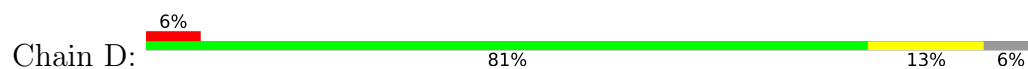


• Molecule 1: Pyruvate kinase PKLR





● Molecule 1: Pyruvate kinase PKLR



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.35Å 174.04Å 87.34Å 90.00° 94.29° 90.00°	Depositor
Resolution (Å)	33.77 – 2.34 33.77 – 2.34	Depositor EDS
% Data completeness (in resolution range)	97.2 (33.77-2.34) 97.2 (33.77-2.34)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.30 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.198 , 0.253 0.199 , 0.253	Depositor DCC
R_{free} test set	4551 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	45.3	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 30.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16072	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% 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: MN, PYR, FBP, K, I07

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	0.36	0/3956	0.56	0/5361
1	B	0.36	0/3982	0.56	0/5398
1	C	0.35	0/3865	0.55	0/5236
1	D	0.38	0/3939	0.57	0/5338
All	All	0.37	0/15742	0.56	0/21333

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3894	0	3974	48	0
1	B	3918	0	3996	57	0
1	C	3806	0	3886	109	0
1	D	3877	0	3955	63	0
2	A	20	0	10	0	0
2	B	20	0	10	0	0
2	C	20	0	10	2	0
2	D	20	0	10	0	0
3	A	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	6	0	0	0	0
3	C	6	0	0	0	0
3	D	6	0	0	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	1	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	B	52	0	0	0	0
6	D	52	0	0	1	0
7	A	91	0	0	1	0
7	B	92	0	0	1	0
7	C	70	0	0	0	0
7	D	108	0	0	2	0
All	All	16072	0	15851	250	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 8.

All (250) 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:C:290:LYS:CD	1:C:292:SER:H	1.31	1.40
1:C:290:LYS:HD3	1:C:292:SER:H	1.03	1.11
1:C:290:LYS:HD2	1:C:292:SER:HB3	1.30	1.07
1:C:290:LYS:CD	1:C:292:SER:N	2.17	1.06
1:C:229:GLN:OE1	1:C:238:THR:C	2.02	1.02
1:C:290:LYS:HD2	1:C:292:SER:H	1.21	0.99
1:C:290:LYS:HD3	1:C:292:SER:N	1.75	0.95
1:C:381:PRO:HB2	1:C:382:ARG:HH11	1.34	0.92
1:C:290:LYS:HD3	1:C:291:ALA:N	1.88	0.89
1:C:290:LYS:HD2	1:C:292:SER:CB	2.03	0.88
1:C:381:PRO:CB	1:C:382:ARG:HH11	1.86	0.87
1:C:381:PRO:CG	1:C:382:ARG:HH11	1.90	0.84
1:C:420:MET:HE2	1:C:424:ILE:HG13	1.60	0.84
1:C:381:PRO:CG	1:C:382:ARG:NH1	2.41	0.82
1:D:355:MET:SD	1:D:359:ARG:CZ	2.69	0.81
1:C:381:PRO:HB2	1:C:382:ARG:NH1	1.97	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:475:THR:HA	2:C:601:FBP:H61	1.64	0.79
1:C:261:LEU:HD12	1:C:262:PRO:HD2	1.64	0.79
1:B:567:ILE:HG12	1:D:569:ARG:HG2	1.64	0.78
1:A:263:GLY:HA2	1:A:289:ARG:HH22	1.47	0.78
1:C:290:LYS:HD2	1:C:292:SER:N	1.89	0.77
1:B:549:GLY:O	1:D:449:ARG:NH2	2.19	0.75
1:C:231:ILE:HG22	1:C:233:PRO:HD3	1.71	0.73
1:A:286:SER:HA	1:A:313:LYS:HE2	1.68	0.73
1:C:381:PRO:CB	1:C:382:ARG:NH1	2.52	0.73
1:C:442:ARG:HG2	1:C:457:ILE:HD11	1.73	0.69
1:C:213:VAL:HG11	1:C:230:LYS:HA	1.73	0.69
1:B:186:THR:HG22	1:B:188:ASP:H	1.55	0.68
1:D:230:LYS:HZ2	1:D:231:ILE:H	1.42	0.68
1:B:178:VAL:HG13	1:B:181:SER:HB2	1.75	0.67
1:C:541:LYS:NZ	1:C:574:SER:O	2.27	0.67
1:C:83:VAL:HG12	1:C:85:ALA:H	1.58	0.66
1:D:228:VAL:HA	1:D:238:THR:HG22	1.78	0.66
1:A:206:ASN:HD21	1:A:209:ARG:HB2	1.62	0.65
1:D:315:GLU:HG2	1:D:336:ALA:HB3	1.79	0.65
1:B:435:ARG:NH2	1:D:443:ARG:NH1	2.44	0.65
1:A:225:SER:HB2	1:A:241:GLU:HB3	1.79	0.65
1:C:368:VAL:HG22	1:C:401:CYS:HB2	1.78	0.64
1:C:225:SER:OG	1:C:241:GLU:OE1	2.16	0.64
1:C:229:GLN:OE1	1:C:238:THR:CA	2.46	0.64
1:A:191:PHE:HA	1:A:194:ARG:HG3	1.79	0.63
1:A:206:ASN:ND2	1:A:209:ARG:HB2	2.13	0.63
1:C:265:SER:O	1:C:269:VAL:HG23	1.98	0.63
1:B:168:GLN:HG3	1:B:195:GLY:O	1.97	0.62
1:B:435:ARG:NH2	1:D:443:ARG:HH11	1.98	0.62
1:C:381:PRO:HG2	1:C:382:ARG:NH1	2.13	0.62
1:C:381:PRO:HG2	1:C:382:ARG:HH11	1.61	0.62
1:C:233:PRO:HB2	1:C:234:GLU:OE1	2.00	0.62
1:B:167:LEU:HD23	1:B:195:GLY:HA3	1.81	0.62
1:A:569:ARG:HG2	1:C:567:ILE:HG12	1.81	0.62
1:D:447:LEU:HD22	1:D:447:LEU:H	1.66	0.60
1:C:420:MET:HE3	1:C:420:MET:HA	1.82	0.60
1:B:234:GLU:H	1:B:234:GLU:CD	2.10	0.60
1:A:284:PHE:HD1	1:A:311:ILE:HB	1.67	0.59
1:D:187:VAL:HG11	1:D:205:PRO:HA	1.84	0.59
1:B:83:VAL:HG12	1:B:85:ALA:H	1.68	0.59
1:C:290:LYS:CD	1:C:292:SER:HB3	2.20	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:603:PYR:O	4:D:604:MN:MN	1.60	0.58
1:B:273:ARG:O	1:B:277:GLU:HG3	2.04	0.58
1:D:87:SER:HB2	1:D:511:GLY:HA2	1.85	0.57
1:A:567:ILE:HG12	1:C:569:ARG:HG2	1.85	0.57
1:C:335:VAL:HG12	1:C:356:MET:HE3	1.86	0.57
1:C:347:GLU:HG2	1:D:423:ALA:HB1	1.85	0.57
1:B:435:ARG:NH2	1:D:443:ARG:HD2	2.20	0.57
1:A:442:ARG:HH22	1:C:442:ARG:HD3	1.70	0.56
1:A:263:GLY:HA2	1:A:289:ARG:NH2	2.17	0.56
1:B:87:SER:HB2	1:B:511:GLY:HA2	1.87	0.56
1:C:290:LYS:HD3	1:C:291:ALA:CA	2.36	0.56
1:B:435:ARG:HH22	1:D:443:ARG:NH1	2.02	0.56
1:C:366:PRO:HB3	1:C:508:LEU:O	2.06	0.56
1:D:566:ASN:OD1	1:D:567:ILE:HG13	2.06	0.56
1:A:517:TYR:OH	1:A:531:ARG:HD2	2.07	0.55
1:D:304:GLU:CD	1:D:304:GLU:H	2.14	0.55
1:B:86:ARG:HD3	1:B:422:HIS:ND1	2.22	0.54
1:D:184:LEU:HD13	1:D:184:LEU:C	2.32	0.54
1:B:566:ASN:OD1	1:B:567:ILE:HG13	2.08	0.54
1:A:440:GLU:CD	1:B:68:THR:HB	2.32	0.54
1:C:225:SER:HB3	1:C:242:ASN:N	2.22	0.54
1:C:290:LYS:HD2	1:C:292:SER:CA	2.37	0.54
1:D:210:VAL:HG22	1:D:257:ALA:HB1	1.89	0.54
1:B:219:ILE:HB	1:B:224:ILE:HB	1.90	0.53
1:B:67:ASP:OD1	7:B:701:HOH:O	2.18	0.53
1:A:61:LEU:O	1:A:65:MET:HG3	2.09	0.53
1:B:264:LEU:HD22	1:B:269:VAL:HG22	1.90	0.53
1:B:289:ARG:HB3	1:B:318:GLU:HG2	1.91	0.52
1:B:450:ASP:OD2	1:B:479:ARG:NH2	2.39	0.52
1:C:87:SER:HB2	1:C:511:GLY:HA2	1.92	0.52
1:D:207:ILE:HA	1:D:210:VAL:HG12	1.91	0.52
1:C:525:TRP:CE2	1:C:560:PRO:HG3	2.45	0.52
1:D:184:LEU:HD12	1:D:199:THR:HG23	1.92	0.51
1:A:125:GLU:CD	1:A:125:GLU:H	2.19	0.51
1:C:286:SER:HA	1:C:313:LYS:HE2	1.91	0.51
1:A:385:ARG:HG2	1:B:372:GLN:OE1	2.11	0.51
1:B:284:PHE:HD1	1:B:311:ILE:HB	1.75	0.51
1:C:83:VAL:O	1:C:426:ARG:HD2	2.11	0.51
1:C:208:VAL:HG12	1:C:236:LEU:HG	1.92	0.51
1:A:267:GLN:NE2	7:A:703:HOH:O	2.43	0.50
1:B:86:ARG:NH2	1:B:113:ASN:OD1	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:ASP:OD1	1:B:321:LYS:NZ	2.44	0.50
1:D:339:ASP:OD2	3:D:603:PYR:O	2.29	0.50
1:B:86:ARG:HB2	1:B:426:ARG:HG3	1.94	0.50
1:C:134:VAL:O	1:C:138:VAL:HG23	2.11	0.50
1:C:219:ILE:HG12	1:C:252:VAL:HG22	1.92	0.50
1:A:61:LEU:HD11	1:A:490:ARG:CZ	2.41	0.49
1:A:181:SER:O	1:A:239:GLN:HG3	2.12	0.49
1:C:370:ALA:HB1	1:C:403:MET:HE2	1.95	0.49
1:D:188:ASP:HB3	1:D:191:PHE:CD2	2.47	0.49
1:B:315:GLU:HG2	1:B:336:ALA:HB3	1.94	0.49
1:D:284:PHE:HD1	1:D:311:ILE:HB	1.77	0.49
1:B:435:ARG:HH21	1:D:443:ARG:HD2	1.77	0.49
1:C:120:SER:OG	1:C:161:GLU:OE2	2.30	0.49
1:C:290:LYS:HD3	1:C:291:ALA:C	2.37	0.49
1:D:163:ARG:HD3	1:D:249:ARG:NH1	2.26	0.49
1:A:218:TYR:HD2	1:A:222:GLY:HA2	1.78	0.49
1:C:420:MET:HE2	1:C:424:ILE:CG1	2.38	0.49
1:C:229:GLN:OE1	1:C:239:GLN:N	2.45	0.49
1:D:304:GLU:O	7:D:701:HOH:O	2.20	0.48
1:B:192:ARG:HG2	1:B:192:ARG:HH11	1.78	0.48
1:A:208:VAL:HA	1:A:236:LEU:HD21	1.96	0.48
1:C:335:VAL:CG2	1:C:369:CYS:HA	2.44	0.48
1:D:514:PRO:O	1:D:515:LEU:HD23	2.13	0.48
1:B:438:PHE:CE2	1:B:442:ARG:HD2	2.49	0.48
1:D:368:VAL:HG22	1:D:401:CYS:HB2	1.95	0.48
1:A:138:VAL:HG21	1:A:151:VAL:HB	1.96	0.47
1:B:86:ARG:HD2	1:B:422:HIS:HA	1.97	0.47
1:B:125:GLU:CD	1:B:125:GLU:H	2.20	0.47
1:C:84:ALA:HA	1:C:426:ARG:HH11	1.79	0.47
1:C:99:ARG:NH2	1:C:129:GLU:HB2	2.29	0.47
1:C:225:SER:HB2	1:C:242:ASN:HB2	1.96	0.47
1:C:290:LYS:CE	1:C:292:SER:H	2.16	0.47
1:C:546:LEU:HD13	1:C:573:ILE:CD1	2.45	0.47
1:D:210:VAL:HG13	1:D:211:VAL:HG13	1.96	0.47
1:C:120:SER:HA	1:C:158:LYS:HG3	1.96	0.47
1:C:291:ALA:HB1	1:C:329:VAL:HG21	1.97	0.47
1:A:368:VAL:HG22	1:A:401:CYS:HB2	1.97	0.46
1:B:229:GLN:HB2	1:B:237:VAL:HG23	1.97	0.46
1:C:86:ARG:HB2	1:C:426:ARG:HG2	1.97	0.46
1:A:250:LYS:HA	1:A:250:LYS:HD3	1.61	0.46
1:A:438:PHE:CE2	1:A:442:ARG:HD3	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:TYR:CE1	1:C:261:LEU:HB2	2.50	0.46
1:C:382:ARG:NH2	1:D:242:ASN:CG	2.74	0.46
1:C:229:GLN:OE1	1:C:238:THR:HA	2.16	0.46
1:D:208:VAL:HA	1:D:236:LEU:HD11	1.96	0.46
1:A:109:LYS:HE3	1:A:141:PHE:HE1	1.79	0.46
1:B:186:THR:CG2	1:B:188:ASP:H	2.26	0.46
1:D:331:ASP:O	1:D:366:PRO:HD2	2.16	0.46
1:A:258:GLN:N	1:A:258:GLN:OE1	2.49	0.45
1:D:355:MET:SD	1:D:359:ARG:NH1	2.89	0.45
1:B:291:ALA:HB2	1:B:325:GLU:HG3	1.98	0.45
1:A:443:ARG:NE	1:C:435:ARG:HH11	2.14	0.45
1:C:218:TYR:O	1:C:252:VAL:HA	2.17	0.45
1:B:192:ARG:HA	1:B:201:TRP:CG	2.52	0.45
1:B:435:ARG:NH2	1:D:443:ARG:CD	2.80	0.45
1:B:499:SER:HB3	1:B:502:ALA:HB3	1.98	0.45
1:C:440:GLU:CD	1:D:68:THR:HB	2.41	0.45
1:B:93:THR:HA	1:B:116:ARG:HB3	1.97	0.45
1:C:91:ILE:HG12	1:C:114:ILE:HB	1.99	0.45
1:B:442:ARG:HH11	1:D:442:ARG:NH2	2.14	0.45
1:C:166:ILE:HA	1:C:248:SER:HB3	1.99	0.45
1:D:184:LEU:HD13	1:D:185:VAL:N	2.32	0.44
1:D:230:LYS:NZ	1:D:231:ILE:HG22	2.33	0.44
1:A:269:VAL:O	1:A:273:ARG:HG3	2.17	0.44
1:A:87:SER:HB2	1:A:511:GLY:HA2	2.00	0.44
1:B:83:VAL:CG1	1:B:85:ALA:H	2.30	0.44
1:C:102:GLU:CD	1:C:102:GLU:H	2.25	0.44
1:A:59:GLN:HB3	1:A:61:LEU:HD13	1.99	0.44
1:D:362:LEU:HD22	7:D:750:HOH:O	2.18	0.44
1:C:164:THR:HA	1:C:201:TRP:O	2.17	0.44
1:D:383:PRO:HG3	1:D:420:MET:HG2	1.99	0.44
1:A:185:VAL:HA	1:A:200:VAL:HG23	1.99	0.43
1:C:469:ALA:HB2	1:C:550:ASP:CG	2.43	0.43
1:D:353:GLN:O	1:D:357:ILE:HG13	2.18	0.43
1:A:273:ARG:O	1:A:277:GLU:HG3	2.17	0.43
1:A:331:ASP:O	1:A:366:PRO:HD2	2.18	0.43
1:C:382:ARG:HH22	1:D:242:ASN:CG	2.26	0.43
1:C:84:ALA:HA	1:C:426:ARG:NH1	2.34	0.43
1:C:232:GLY:N	1:C:233:PRO:HD3	2.33	0.43
1:C:269:VAL:O	1:C:273:ARG:HG2	2.18	0.43
1:B:286:SER:HA	1:B:313:LYS:HD3	2.00	0.43
1:D:337:ARG:HD3	1:D:353:GLN:OE1	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:466:CYS:HB3	1:C:454:VAL:HG21	2.01	0.43
1:B:331:ASP:O	1:B:366:PRO:HD2	2.19	0.43
1:C:207:ILE:HG13	1:C:254:LEU:HD11	2.01	0.43
1:C:525:TRP:CD1	1:C:560:PRO:HG3	2.54	0.43
1:C:382:ARG:HH22	1:D:242:ASN:CB	2.32	0.43
1:A:185:VAL:HB	1:A:236:LEU:HB2	2.00	0.43
1:D:137:ALA:O	1:D:140:SER:HB3	2.19	0.43
1:A:442:ARG:HH22	1:C:442:ARG:CD	2.30	0.43
1:A:558:TRP:CE3	1:A:559:ARG:HB2	2.53	0.42
1:C:85:ALA:HB2	1:C:545:PHE:CD2	2.54	0.42
1:C:124:HIS:NE2	1:C:271:ASP:OD1	2.44	0.42
1:D:264:LEU:HD12	1:D:264:LEU:HA	1.85	0.42
1:B:83:VAL:HG12	1:B:85:ALA:N	2.32	0.42
1:D:188:ASP:CG	1:D:191:PHE:HD2	2.26	0.42
1:D:370:ALA:HB1	1:D:403:MET:HE2	2.01	0.42
1:B:443:ARG:HD3	1:D:435:ARG:HH11	1.83	0.42
1:D:420:MET:HE3	1:D:424:ILE:HD11	2.01	0.42
1:B:91:ILE:HG12	1:B:114:ILE:HB	2.02	0.42
1:B:496:VAL:HG11	1:B:532:ARG:HB3	2.01	0.42
1:B:408:THR:HA	1:B:414:PRO:HB3	2.01	0.42
1:A:183:VAL:HG22	1:A:198:ASN:HA	2.01	0.42
1:C:554:VAL:HG21	1:C:571:LEU:HD12	2.00	0.42
1:C:182:GLN:O	1:C:198:ASN:ND2	2.48	0.42
1:C:208:VAL:HA	1:C:236:LEU:HD21	2.01	0.42
1:B:234:GLU:OE2	1:B:234:GLU:N	2.45	0.42
1:D:134:VAL:O	1:D:138:VAL:HG23	2.19	0.42
1:A:221:ASP:N	1:A:221:ASP:OD1	2.49	0.42
1:A:272:LEU:O	1:A:276:VAL:HG23	2.20	0.42
1:C:83:VAL:CG1	1:C:85:ALA:H	2.29	0.42
1:C:225:SER:O	1:C:226:LEU:HD23	2.19	0.42
1:C:231:ILE:HG22	1:C:233:PRO:CD	2.46	0.42
1:A:218:TYR:O	1:A:252:VAL:HA	2.20	0.41
1:B:400:ASP:HA	1:B:510:ARG:HB2	2.01	0.41
1:C:125:GLU:H	1:C:125:GLU:CD	2.27	0.41
1:D:184:LEU:HD21	1:D:235:GLY:HA3	2.02	0.41
1:C:83:VAL:HG12	1:C:85:ALA:N	2.31	0.41
1:C:347:GLU:HG2	1:D:423:ALA:CB	2.50	0.41
1:B:175:VAL:HG11	1:B:197:ALA:N	2.35	0.41
1:C:447:LEU:HD22	1:C:447:LEU:H	1.85	0.41
1:D:87:SER:HB2	1:D:511:GLY:CA	2.49	0.41
1:D:286:SER:HA	1:D:313:LYS:HD3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:445:ALA:HB3	1:C:465:LYS:NZ	2.35	0.41
1:C:97:ALA:O	1:C:103:ARG:HD2	2.21	0.41
1:C:284:PHE:HD1	1:C:311:ILE:HB	1.85	0.41
1:C:365:LYS:HA	1:C:366:PRO:HD3	1.96	0.41
1:D:188:ASP:HA	1:D:189:PRO:HD3	1.91	0.41
1:C:382:ARG:HH22	1:D:242:ASN:HB3	1.86	0.41
1:D:211:VAL:HG21	1:D:217:ILE:HD11	2.02	0.41
1:B:264:LEU:HD23	1:B:268:ASP:HB2	2.03	0.41
1:C:442:ARG:CZ	1:C:442:ARG:HB3	2.51	0.41
1:A:137:ALA:O	1:A:140:SER:HB3	2.21	0.41
1:A:473:VAL:HG23	1:A:555:VAL:HB	2.03	0.41
1:A:547:ARG:HE	1:A:547:ARG:HB2	1.71	0.41
1:B:514:PRO:O	1:B:515:LEU:HD23	2.21	0.41
1:C:68:THR:HB	1:D:440:GLU:CD	2.45	0.41
1:C:224:ILE:HD13	1:C:224:ILE:HA	1.76	0.41
1:C:453:GLU:OE2	1:C:483:LEU:HD22	2.20	0.41
1:D:216:ARG:HA	1:D:226:LEU:O	2.21	0.41
1:D:235:GLY:C	1:D:236:LEU:HD12	2.46	0.41
1:A:450:ASP:O	1:A:454:VAL:HG23	2.21	0.40
1:B:547:ARG:HD2	1:B:547:ARG:HA	1.76	0.40
1:D:437:LEU:HB2	6:D:601[B]:I07:C16	2.50	0.40
1:B:362:LEU:HD11	1:B:444:ALA:HB1	2.03	0.40
1:C:272:LEU:O	1:C:276:VAL:HG23	2.20	0.40
1:C:525:TRP:HZ2	2:C:601:FBP:H12	1.86	0.40
1:B:366:PRO:HB3	1:B:508:LEU:O	2.22	0.40
1:C:453:GLU:HG2	1:C:483:LEU:HD13	2.03	0.40
1:D:321:LYS:HD2	1:D:321:LYS:HA	1.76	0.40
1:C:290:LYS:CD	1:C:292:SER:CB	2.88	0.40
1:C:525:TRP:NE1	1:C:560:PRO:HG3	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	510/544 (94%)	498 (98%)	10 (2%)	2 (0%)	30	32
1	B	516/544 (95%)	503 (98%)	9 (2%)	4 (1%)	16	16
1	C	497/544 (91%)	484 (97%)	10 (2%)	3 (1%)	21	22
1	D	508/544 (93%)	492 (97%)	14 (3%)	2 (0%)	30	32
All	All	2031/2176 (93%)	1977 (97%)	43 (2%)	11 (0%)	24	26

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	171	PRO
1	C	233	PRO
1	B	168	GLN
1	B	371	THR
1	C	371	THR
1	D	175	VAL
1	D	371	THR
1	A	566	ASN
1	B	170	GLY
1	A	371	THR
1	C	566	ASN

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	413/436 (95%)	413 (100%)	0	100	100
1	B	415/436 (95%)	415 (100%)	0	100	100
1	C	404/436 (93%)	404 (100%)	0	100	100
1	D	411/436 (94%)	411 (100%)	0	100	100
All	All	1643/1744 (94%)	1643 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	HIS
1	A	206	ASN
1	A	278	HIS
1	A	434	HIS
1	B	59	GLN
1	C	57	GLN
1	C	182	GLN
1	C	421	GLN
1	D	239	GLN
1	D	434	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 8 are monoatomic - leaving 12 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	I07	D	601[A]	-	26,29,29	2.83	12 (46%)	29,43,43	1.62	6 (20%)
2	FBP	B	602	-	18,20,20	0.91	1 (5%)	21,32,32	0.98	0
3	PYR	D	603	4	5,5,5	2.80	3 (60%)	3,6,6	2.62	1 (33%)
2	FBP	C	601	-	18,20,20	1.02	1 (5%)	21,32,32	1.01	1 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FBP	D	602	-	18,20,20	0.85	1 (5%)	21,32,32	0.92	1 (4%)
3	PYR	A	602	4	5,5,5	2.83	3 (60%)	3,6,6	3.70	3 (100%)
3	PYR	C	602	4	5,5,5	2.92	3 (60%)	3,6,6	3.15	2 (66%)
6	I07	D	601[B]	-	26,29,29	2.90	11 (42%)	29,43,43	1.92	9 (31%)
2	FBP	A	601	-	18,20,20	0.89	1 (5%)	21,32,32	1.13	1 (4%)
6	I07	B	601[B]	-	26,29,29	2.84	12 (46%)	29,43,43	1.87	6 (20%)
6	I07	B	601[A]	-	26,29,29	3.00	13 (50%)	29,43,43	1.79	8 (27%)
3	PYR	B	603	4	5,5,5	2.89	3 (60%)	3,6,6	1.82	2 (66%)

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	I07	D	601[A]	-	-	0/6/10/10	0/4/4/4
2	FBP	B	602	-	-	3/13/32/32	0/1/1/1
3	PYR	D	603	4	-	0/4/4/4	-
2	FBP	C	601	-	-	8/13/32/32	0/1/1/1
2	FBP	D	602	-	-	2/13/32/32	0/1/1/1
3	PYR	A	602	4	-	0/4/4/4	-
3	PYR	C	602	4	-	0/4/4/4	-
6	I07	D	601[B]	-	-	0/6/10/10	0/4/4/4
2	FBP	A	601	-	-	5/13/32/32	0/1/1/1
6	I07	B	601[B]	-	-	0/6/10/10	0/4/4/4
6	I07	B	601[A]	-	-	0/6/10/10	0/4/4/4
3	PYR	B	603	4	-	0/4/4/4	-

All (64) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	601[B]	I07	C7-N2	8.44	1.39	1.29
6	B	601[B]	I07	C7-N2	8.21	1.39	1.29
6	B	601[A]	I07	C7-N2	8.17	1.39	1.29
6	D	601[A]	I07	C7-N2	7.65	1.38	1.29
6	B	601[A]	I07	C8-C10	-6.51	1.39	1.51
6	D	601[A]	I07	C8-C10	-5.79	1.41	1.51
6	D	601[B]	I07	C8-C10	-5.54	1.41	1.51
6	B	601[B]	I07	C8-C10	-5.31	1.41	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	602	PYR	O3-CA	4.36	1.33	1.23
6	D	601[A]	I07	C5-C3	-4.28	1.32	1.44
6	B	601[A]	I07	C3-S	-4.19	1.65	1.74
3	B	603	PYR	CA-C	-4.18	1.40	1.54
6	B	601[A]	I07	C5-C3	-4.14	1.32	1.44
6	B	601[B]	I07	C-S	-4.11	1.57	1.70
6	D	601[B]	I07	C5-C3	-4.09	1.32	1.44
6	B	601[B]	I07	C5-C3	-4.06	1.32	1.44
3	D	603	PYR	O-C	3.90	1.32	1.22
3	A	602	PYR	O-C	3.85	1.32	1.22
3	D	603	PYR	CA-C	-3.82	1.41	1.54
6	D	601[A]	I07	C7-C5	-3.78	1.33	1.41
3	A	602	PYR	O3-CA	3.76	1.31	1.23
6	B	601[A]	I07	C7-C5	-3.75	1.33	1.41
6	B	601[B]	I07	C7-C5	-3.71	1.33	1.41
3	B	603	PYR	O3-CA	3.68	1.31	1.23
6	D	601[A]	I07	C4-C6	-3.63	1.33	1.43
6	D	601[B]	I07	C3-S	-3.62	1.66	1.74
6	D	601[B]	I07	C7-C5	-3.61	1.33	1.41
6	D	601[A]	I07	C4-N	-3.58	1.33	1.39
6	D	601[B]	I07	C4-C6	-3.53	1.34	1.43
3	C	602	PYR	O-C	3.50	1.31	1.22
6	D	601[A]	I07	C1-C2	-3.48	1.34	1.42
6	B	601[A]	I07	C4-C6	-3.40	1.34	1.43
6	B	601[A]	I07	C-S	-3.39	1.59	1.70
3	C	602	PYR	CA-C	-3.37	1.43	1.54
6	B	601[B]	I07	C2-C3	-3.36	1.34	1.39
6	D	601[B]	I07	C4-N	-3.35	1.33	1.39
6	D	601[B]	I07	C2-C3	-3.34	1.34	1.39
6	D	601[B]	I07	C1-C2	-3.29	1.34	1.42
3	B	603	PYR	O-C	3.27	1.30	1.22
3	A	602	PYR	CA-C	-3.24	1.43	1.54
6	B	601[A]	I07	C2-C3	-3.24	1.34	1.39
6	B	601[B]	I07	C4-C6	-3.15	1.35	1.43
6	D	601[A]	I07	C2-C3	-3.13	1.34	1.39
6	D	601[B]	I07	C5-C4	-3.06	1.33	1.43
6	D	601[A]	I07	C5-C4	-3.01	1.33	1.43
6	B	601[A]	I07	C5-C4	-2.99	1.33	1.43
3	D	603	PYR	O3-CA	2.98	1.30	1.23
6	B	601[A]	I07	C1-C2	-2.96	1.35	1.42
6	B	601[A]	I07	C4-N	-2.94	1.34	1.39
6	B	601[B]	I07	C5-C4	-2.90	1.33	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	FBP	O2-C2	2.84	1.45	1.40
6	B	601[B]	I07	C1-C2	-2.82	1.35	1.42
6	B	601[B]	I07	C4-N	-2.80	1.34	1.39
6	D	601[B]	I07	C2-N	-2.69	1.32	1.39
2	C	601	FBP	O2-C2	2.67	1.45	1.40
6	D	601[A]	I07	C3-S	-2.65	1.68	1.74
2	D	602	FBP	O2-C2	2.48	1.45	1.40
2	B	602	FBP	O2-C2	2.44	1.44	1.40
6	D	601[A]	I07	C2-N	-2.43	1.33	1.39
6	B	601[B]	I07	C3-S	-2.37	1.69	1.74
6	B	601[A]	I07	C2-N	-2.16	1.34	1.39
6	D	601[A]	I07	C-S	-2.13	1.63	1.70
6	B	601[A]	I07	C9-N	2.13	1.50	1.47
6	B	601[B]	I07	C2-N	-2.06	1.34	1.39

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	601[B]	I07	C3-C2-N	-4.72	106.59	110.19
6	B	601[A]	I07	C3-C2-N	-4.60	106.69	110.19
6	B	601[B]	I07	C8-N1-N2	4.51	119.76	114.38
3	A	602	PYR	OXT-C-O	-4.45	113.31	123.90
6	B	601[B]	I07	C3-C2-N	-4.43	106.82	110.19
6	D	601[B]	I07	C8-N1-N2	4.35	119.56	114.38
6	B	601[B]	I07	C2-C3-S	-3.98	108.17	111.16
3	C	602	PYR	OXT-C-O	-3.90	114.61	123.90
6	D	601[A]	I07	C3-C2-N	-3.68	107.39	110.19
3	D	603	PYR	OXT-C-CA	3.66	123.75	113.59
6	D	601[A]	I07	C10-C8-N1	-3.58	107.21	112.76
6	D	601[B]	I07	C10-C8-N1	-3.56	107.25	112.76
3	A	602	PYR	OXT-C-CA	3.55	123.43	113.59
2	A	601	FBP	O6-P2-O4P	3.45	115.77	106.44
6	B	601[B]	I07	C10-C8-N1	-3.45	107.43	112.76
3	C	602	PYR	OXT-C-CA	3.44	123.13	113.59
6	D	601[A]	I07	C8-N1-N2	3.33	118.35	114.38
6	D	601[B]	I07	C2-C3-S	-3.15	108.80	111.16
6	B	601[A]	I07	C2-C3-S	-3.12	108.82	111.16
6	B	601[A]	I07	C8-N1-N2	3.09	118.06	114.38
3	A	602	PYR	O3-CA-CB	-2.96	113.15	119.77
6	B	601[A]	I07	C17-O2-C12	-2.90	111.27	117.50
6	D	601[A]	I07	O-C6-C4	-2.86	120.52	126.38
6	B	601[A]	I07	O-C6-C4	-2.76	120.73	126.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	601[B]	I07	O-C6-C4	-2.71	120.83	126.38
6	D	601[B]	I07	C17-O2-C12	-2.56	112.02	117.50
6	D	601[B]	I07	C9-N-C4	2.53	132.39	125.40
6	B	601[B]	I07	C5-C4-N	-2.53	106.92	109.41
6	D	601[B]	I07	C5-C4-N	-2.41	107.04	109.41
6	B	601[A]	I07	C10-C8-N1	-2.36	109.10	112.76
6	B	601[A]	I07	C8-N1-C6	2.33	121.40	119.22
6	B	601[B]	I07	O-C6-N1	-2.28	117.17	121.17
3	B	603	PYR	OXT-C-CA	2.27	119.88	113.59
2	C	601	FBP	O6P-P2-O6	2.27	112.58	106.67
6	D	601[B]	I07	C9-N-C2	-2.26	118.20	124.46
2	D	602	FBP	O2P-P1-O1	2.26	112.56	106.67
3	B	603	PYR	OXT-C-O	-2.19	118.69	123.90
6	B	601[A]	I07	C5-C4-N	-2.15	107.29	109.41
6	D	601[A]	I07	C3-S-C	-2.14	92.55	95.52
6	D	601[A]	I07	C8-N1-C6	2.12	121.20	119.22

There are no chirality outliers.

All (18) torsion outliers are listed below:

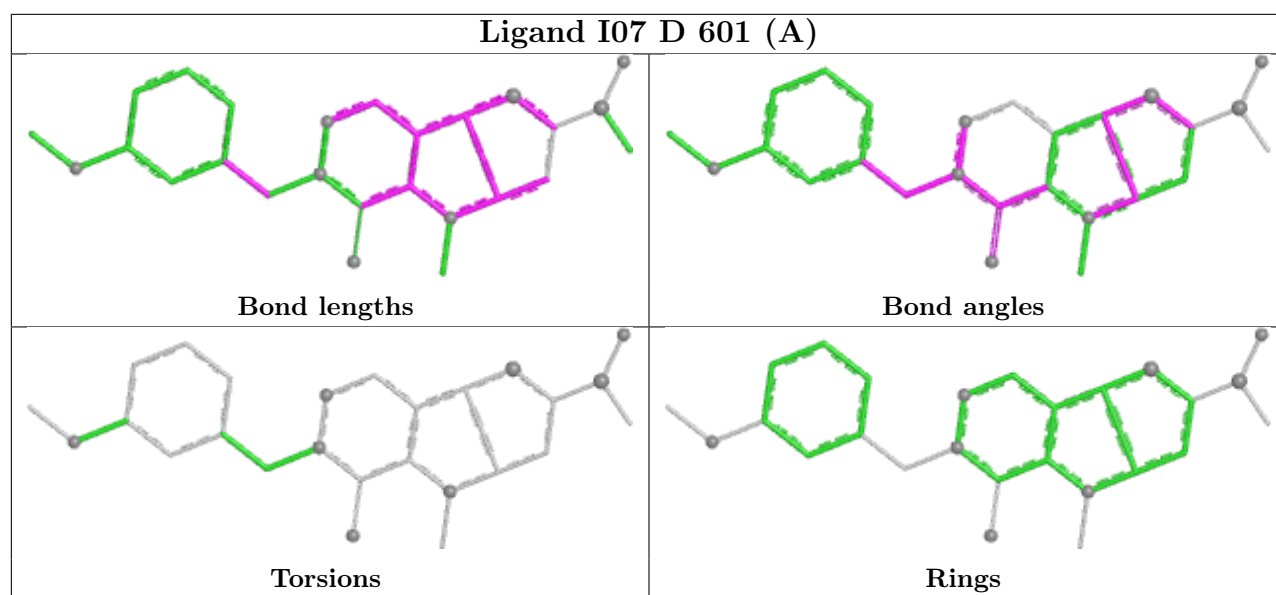
Mol	Chain	Res	Type	Atoms
2	A	601	FBP	C1-O1-P1-O2P
2	A	601	FBP	C1-O1-P1-O3P
2	B	602	FBP	C1-O1-P1-O3P
2	C	601	FBP	C1-O1-P1-O1P
2	C	601	FBP	C1-O1-P1-O3P
2	C	601	FBP	C6-O6-P2-O4P
2	C	601	FBP	C6-O6-P2-O5P
2	C	601	FBP	C6-O6-P2-O6P
2	D	602	FBP	C4-C5-C6-O6
2	A	601	FBP	C4-C5-C6-O6
2	B	602	FBP	C4-C5-C6-O6
2	D	602	FBP	O5-C5-C6-O6
2	C	601	FBP	O5-C5-C6-O6
2	A	601	FBP	O5-C5-C6-O6
2	B	602	FBP	O5-C5-C6-O6
2	C	601	FBP	C4-C5-C6-O6
2	A	601	FBP	C1-O1-P1-O1P
2	C	601	FBP	C1-O1-P1-O2P

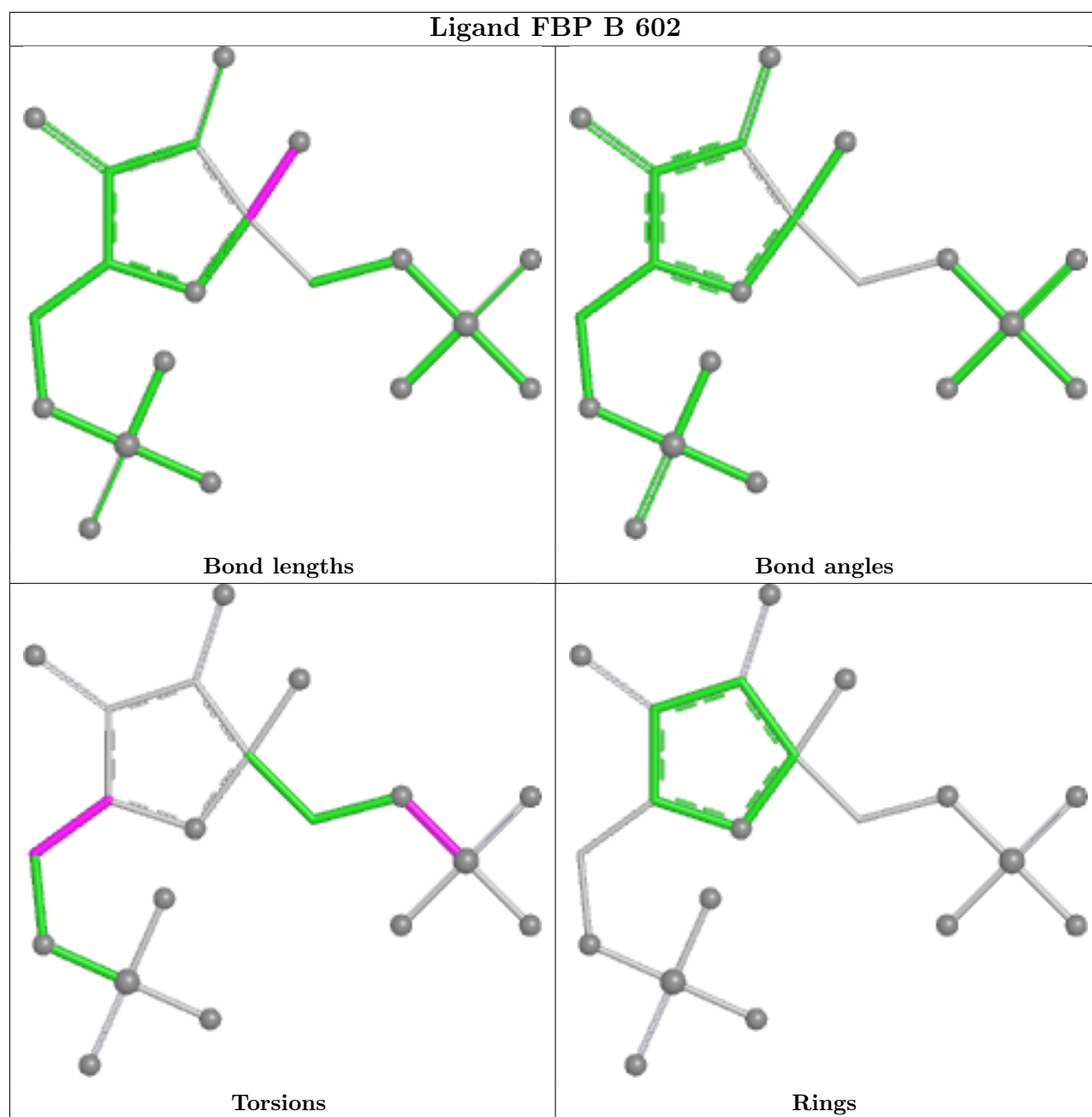
There are no ring outliers.

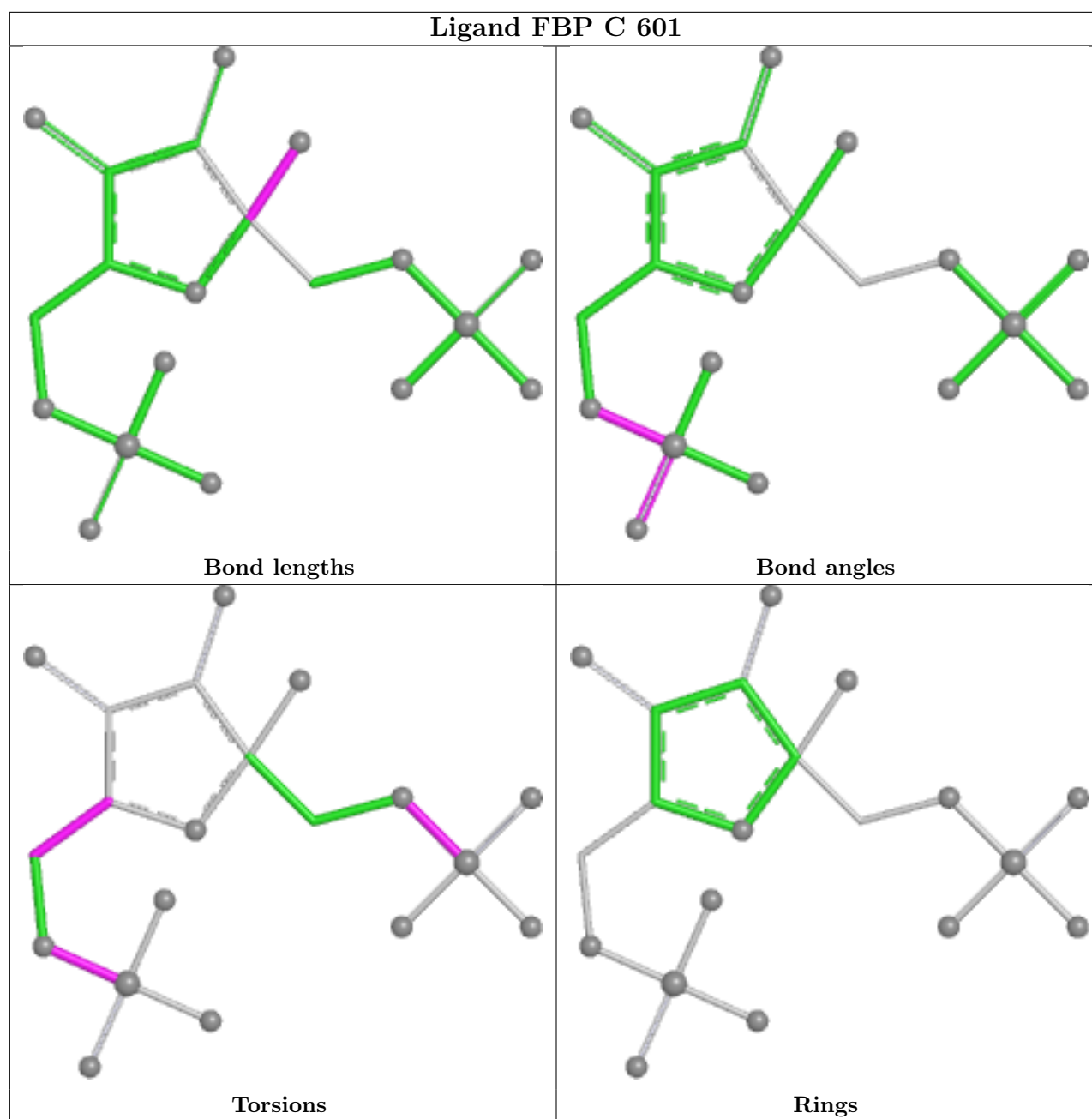
3 monomers are involved in 5 short contacts:

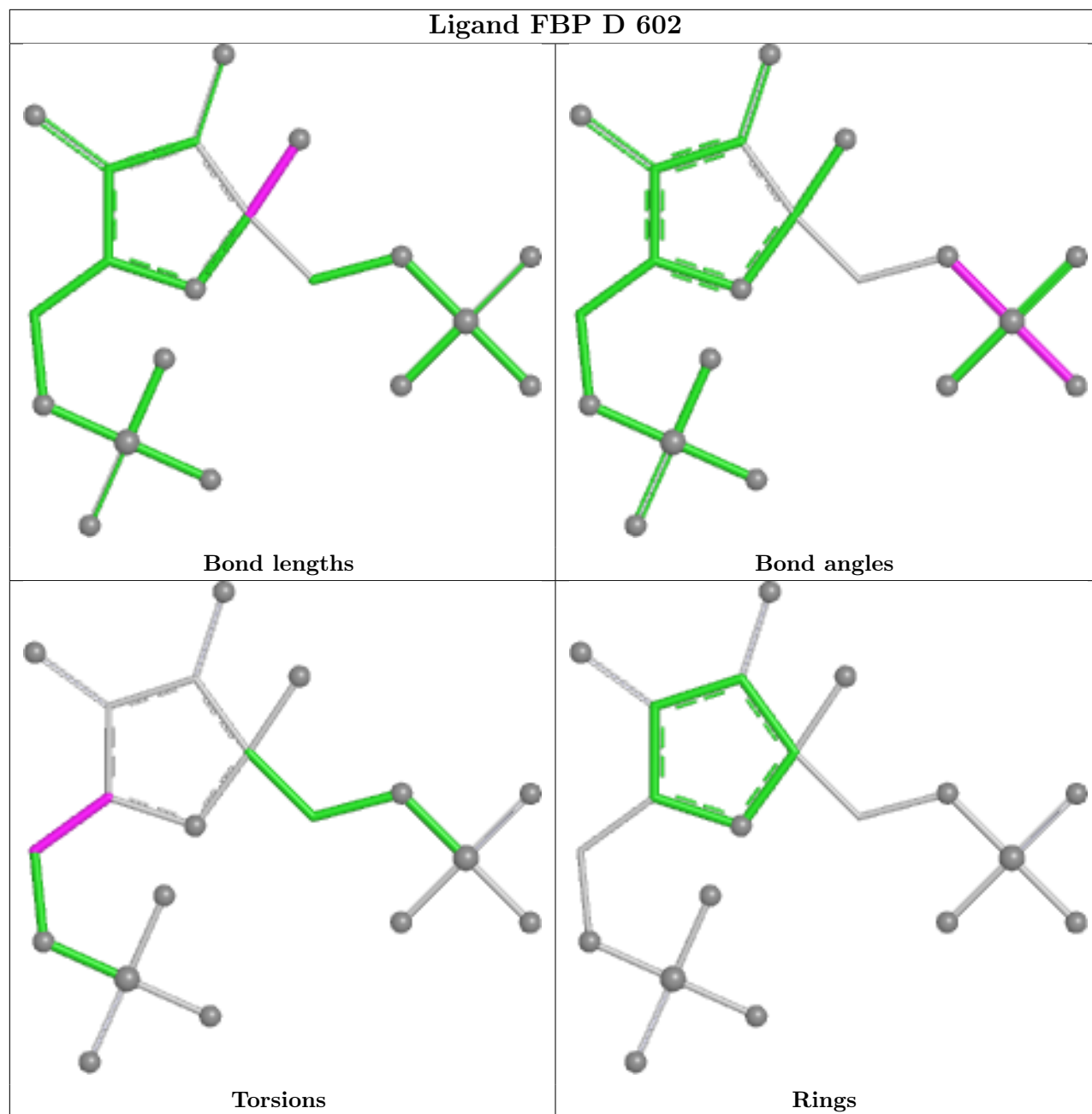
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	603	PYR	2	0
2	C	601	FBP	2	0
6	D	601[B]	I07	1	0

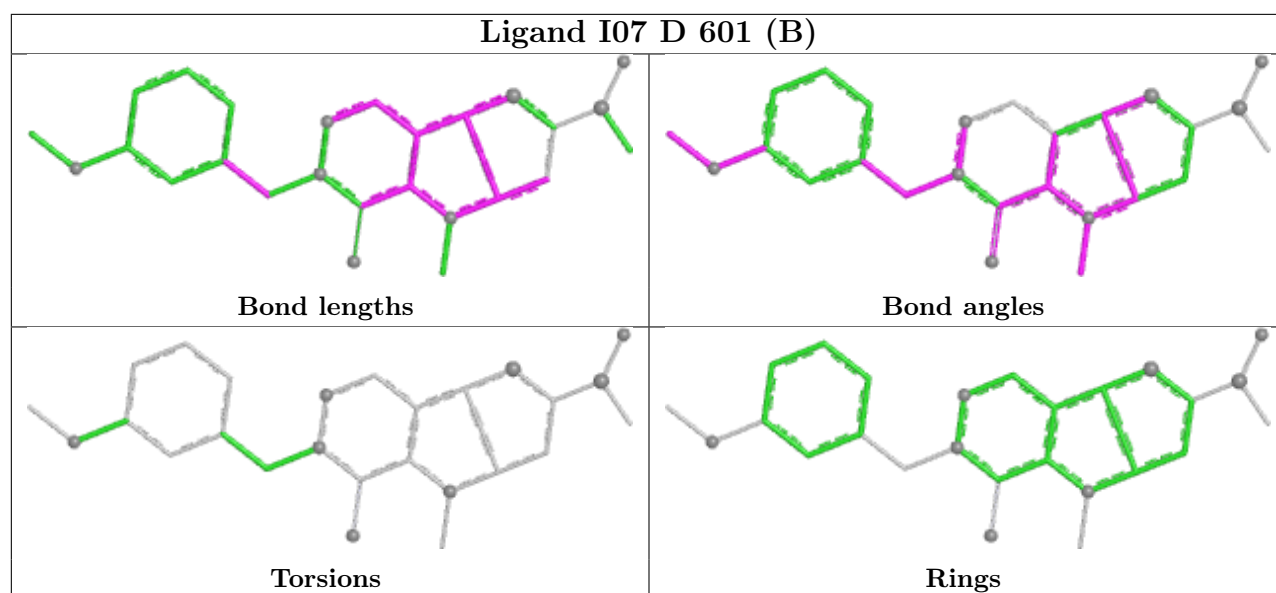
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



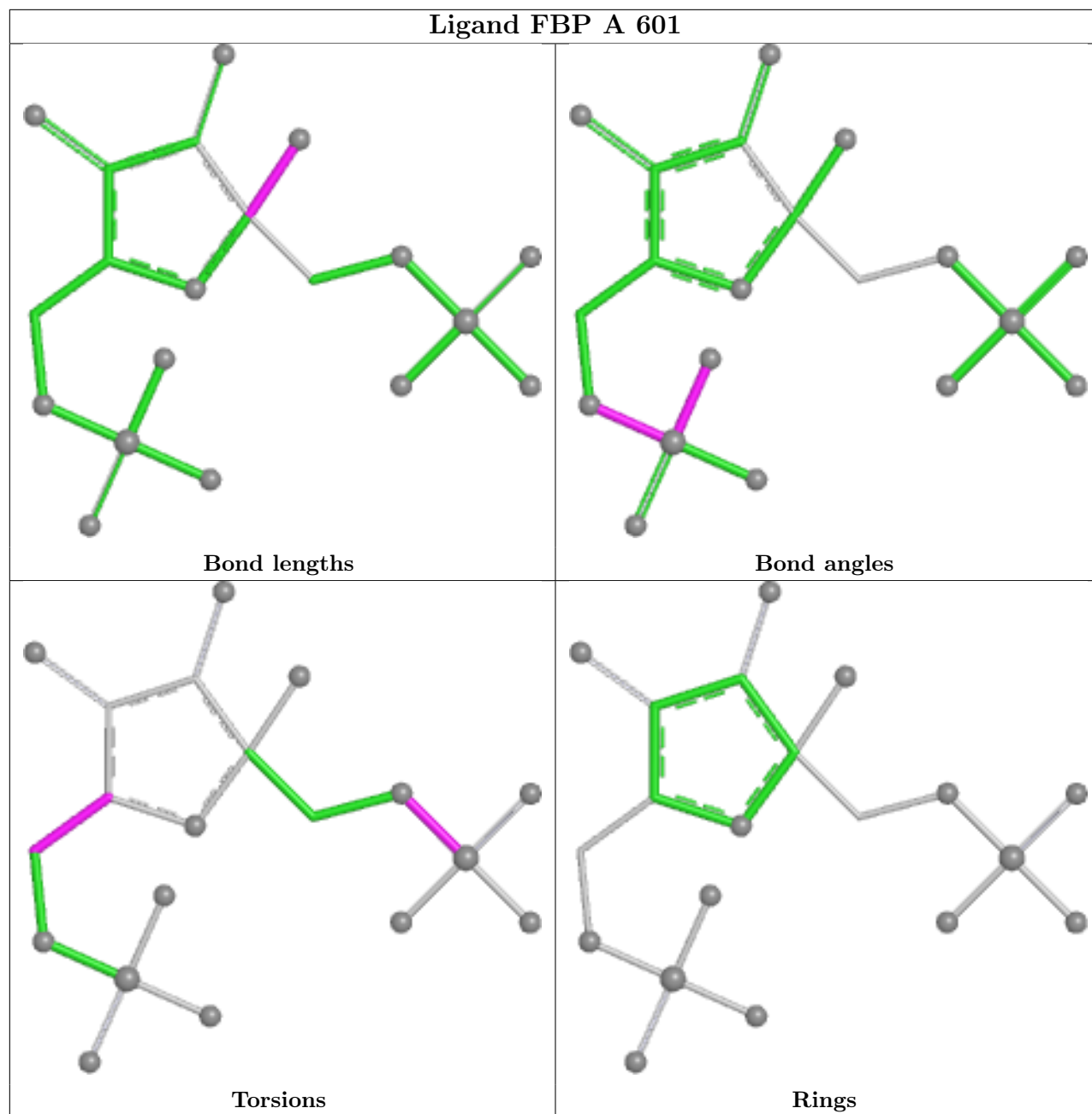


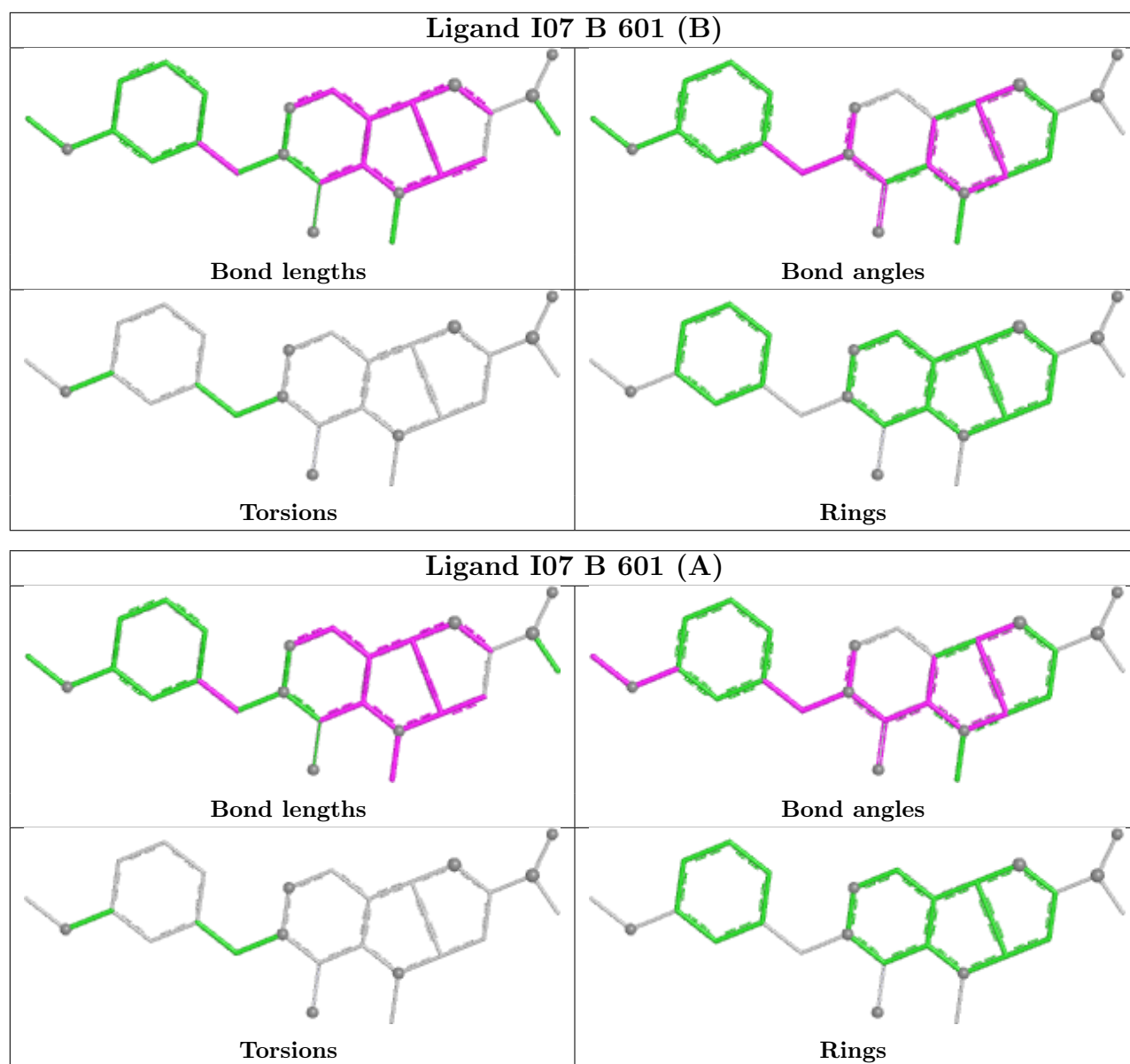






Ligand FBP A 601





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	514/544 (94%)	0.27	27 (5%) 32 37	34, 49, 95, 119	0
1	B	518/544 (95%)	0.26	19 (3%) 45 51	32, 51, 84, 123	0
1	C	503/544 (92%)	0.58	41 (8%) 17 21	35, 55, 97, 141	0
1	D	512/544 (94%)	0.20	33 (6%) 25 30	31, 47, 92, 119	0
All	All	2047/2176 (94%)	0.33	120 (5%) 28 33	31, 51, 93, 141	0

All (120) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	224	ILE	5.2
1	C	231	ILE	4.8
1	D	191	PHE	4.3
1	D	166	ILE	4.2
1	D	229	GLN	3.9
1	B	180	GLY	3.7
1	B	166	ILE	3.7
1	C	144	SER	3.5
1	C	232	GLY	3.5
1	C	260	ASP	3.3
1	D	231	ILE	3.3
1	B	191	PHE	3.2
1	A	229	GLN	3.2
1	C	84	ALA	3.1
1	A	231	ILE	3.1
1	C	544	GLY	3.1
1	D	165	GLY	3.1
1	D	210	VAL	3.0
1	D	233	PRO	3.0
1	A	180	GLY	3.0
1	C	166	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	181	SER	2.9
1	D	434	HIS	2.9
1	D	178	VAL	2.9
1	D	181	SER	2.8
1	D	200	VAL	2.8
1	B	171	PRO	2.8
1	C	296	ALA	2.8
1	A	210	VAL	2.8
1	A	237	VAL	2.8
1	B	181	SER	2.8
1	A	213	VAL	2.8
1	C	306	HIS	2.7
1	D	447	LEU	2.7
1	A	227	VAL	2.7
1	C	185	VAL	2.7
1	C	180	GLY	2.7
1	D	173	SER	2.7
1	D	237	VAL	2.7
1	C	233	PRO	2.7
1	C	261	LEU	2.7
1	A	238	THR	2.6
1	A	243	GLY	2.6
1	C	546	LEU	2.6
1	C	227	VAL	2.6
1	D	199	THR	2.6
1	B	83	VAL	2.6
1	C	165	GLY	2.6
1	D	232	GLY	2.6
1	B	64	ALA	2.6
1	A	240	VAL	2.6
1	A	225	SER	2.6
1	C	173	SER	2.6
1	C	200	VAL	2.5
1	C	213	VAL	2.5
1	D	230	LYS	2.5
1	C	58	GLN	2.5
1	C	208	VAL	2.5
1	A	241	GLU	2.5
1	A	244	GLY	2.5
1	C	202	VAL	2.4
1	D	213	VAL	2.4
1	B	521	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	205	PRO	2.4
1	A	235	GLY	2.4
1	B	167	LEU	2.4
1	D	550	ASP	2.4
1	C	287	PHE	2.4
1	C	538	GLU	2.3
1	B	67	ASP	2.3
1	B	170	GLY	2.3
1	C	162	ILE	2.3
1	C	259	VAL	2.3
1	C	145	PRO	2.2
1	D	574	SER	2.2
1	A	224	ILE	2.2
1	D	180	GLY	2.2
1	C	181	SER	2.2
1	C	245	VAL	2.2
1	A	206	ASN	2.2
1	C	303	PRO	2.2
1	D	186	THR	2.2
1	D	247	GLY	2.2
1	A	173	SER	2.2
1	B	186	THR	2.2
1	A	166	ILE	2.1
1	B	65	MET	2.1
1	B	178	VAL	2.1
1	A	233	PRO	2.1
1	B	84	ALA	2.1
1	B	190	ALA	2.1
1	C	65	MET	2.1
1	D	184	LEU	2.1
1	A	81	GLU	2.1
1	A	260	ASP	2.1
1	C	196	ASN	2.1
1	C	228	VAL	2.1
1	D	183	VAL	2.1
1	A	182	GLN	2.1
1	B	306	HIS	2.1
1	D	236	LEU	2.1
1	D	188	ASP	2.1
1	C	240	VAL	2.1
1	D	245	VAL	2.1
1	C	246	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	242	ASN	2.1
1	B	245	VAL	2.1
1	D	187	VAL	2.1
1	D	228	VAL	2.1
1	C	214	GLY	2.1
1	C	266	GLU	2.0
1	D	197	ALA	2.0
1	B	57	GLN	2.0
1	C	242	ASN	2.0
1	A	228	VAL	2.0
1	D	185	VAL	2.0
1	D	240	VAL	2.0
1	C	234	GLU	2.0
1	A	239	GLN	2.0
1	A	246	LEU	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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

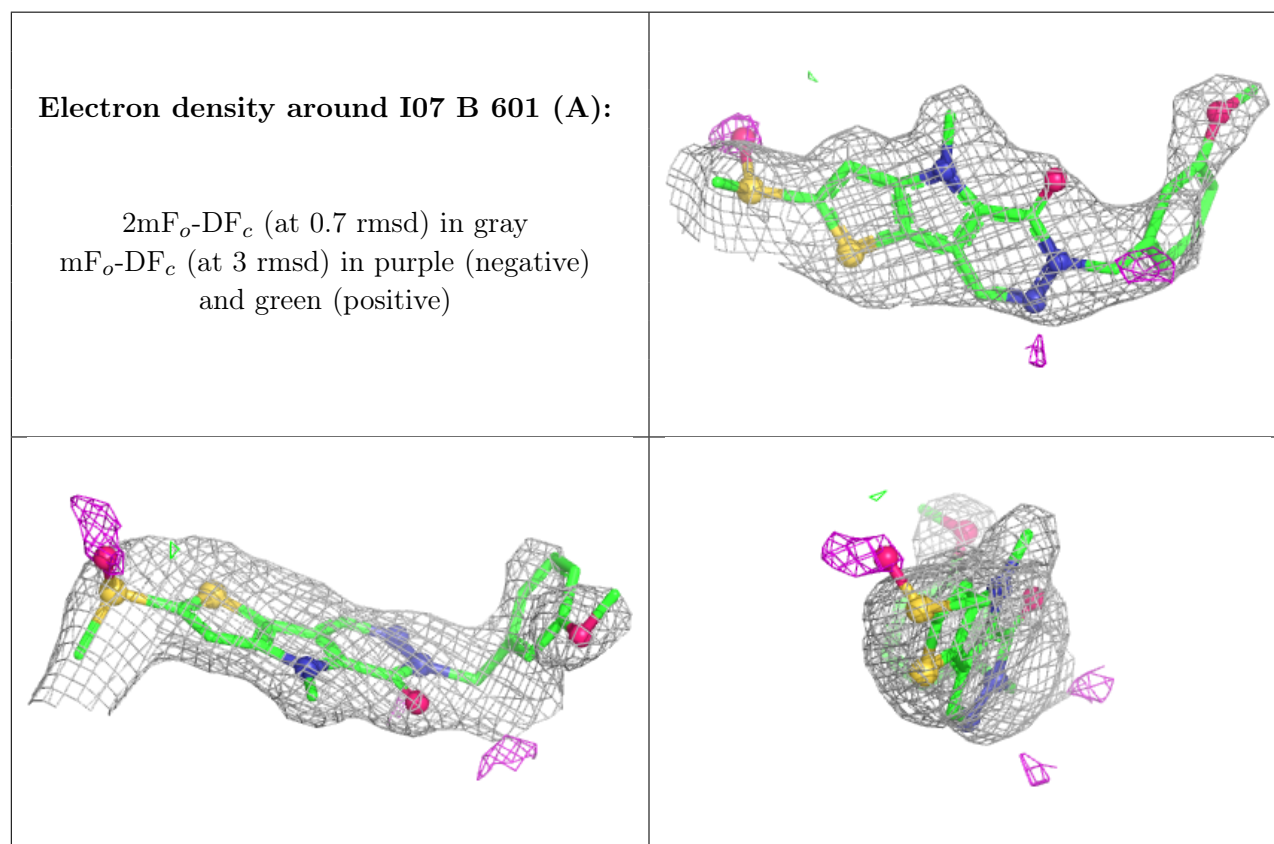
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	K	C	604	1/1	0.80	0.19	73,73,73,73	0
3	PYR	B	603	6/6	0.86	0.17	51,65,85,120	0
3	PYR	D	603	6/6	0.89	0.12	36,48,54,72	0
3	PYR	C	602	6/6	0.89	0.12	51,53,65,66	0
4	MN	B	604	1/1	0.90	0.28	163,163,163,163	0
6	I07	B	601[A]	26/26	0.90	0.14	39,45,47,51	26
6	I07	B	601[B]	26/26	0.90	0.14	42,45,47,49	26
3	PYR	A	602	6/6	0.92	0.12	44,48,51,58	0

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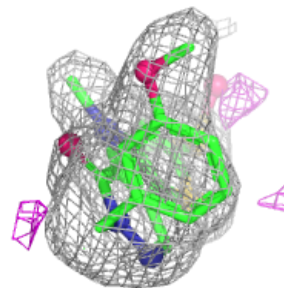
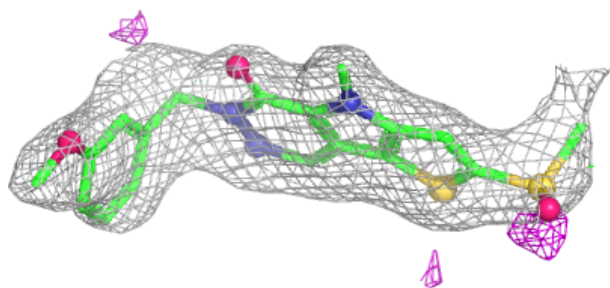
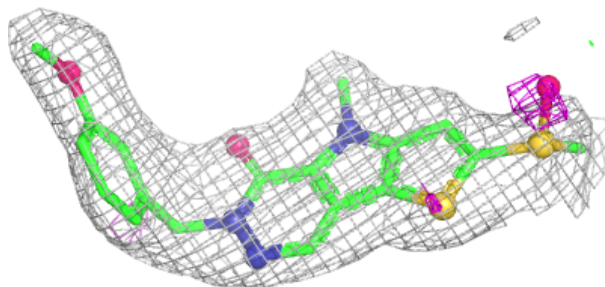
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	I07	D	601[A]	26/26	0.92	0.11	43,47,50,52	26
6	I07	D	601[B]	26/26	0.92	0.11	44,47,49,54	26
2	FBP	C	601	20/20	0.93	0.11	43,47,54,58	0
2	FBP	B	602	20/20	0.94	0.09	43,47,55,56	0
4	MN	A	603	1/1	0.94	0.20	98,98,98,98	0
4	MN	D	604	1/1	0.95	0.14	72,72,72,72	0
2	FBP	A	601	20/20	0.95	0.08	44,50,60,61	0
5	K	B	605	1/1	0.96	0.12	55,55,55,55	0
5	K	D	605	1/1	0.96	0.12	51,51,51,51	0
2	FBP	D	602	20/20	0.97	0.07	37,46,48,50	0
5	K	A	604	1/1	0.99	0.13	54,54,54,54	0
4	MN	C	603	1/1	0.99	0.03	57,57,57,57	0

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

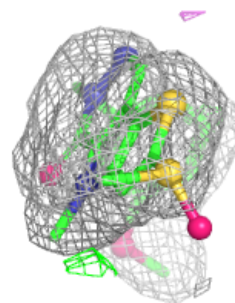
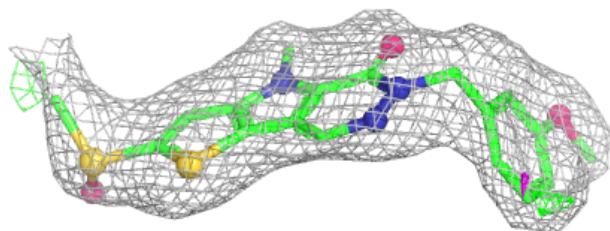
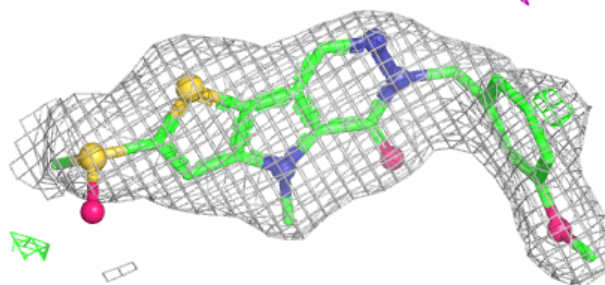


Electron density around I07 B 601 (B):

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

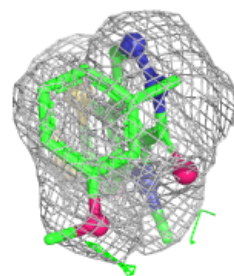
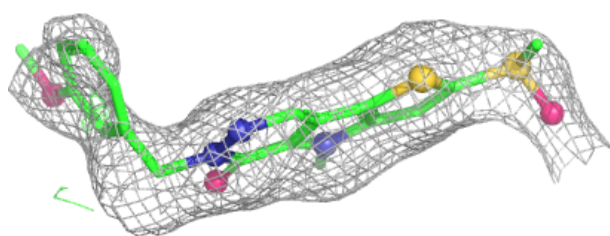
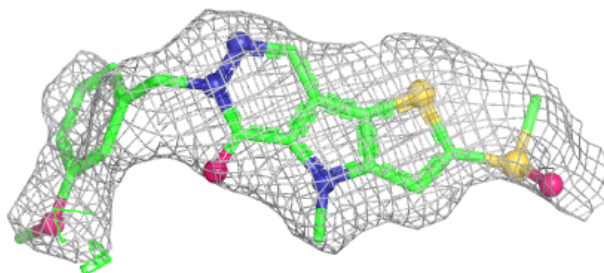
**Electron density around I07 D 601 (A):**

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

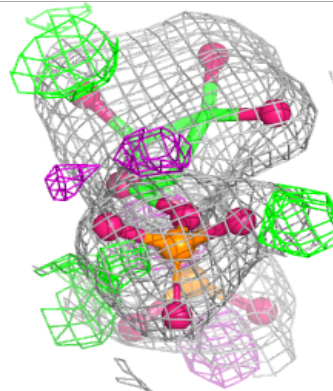
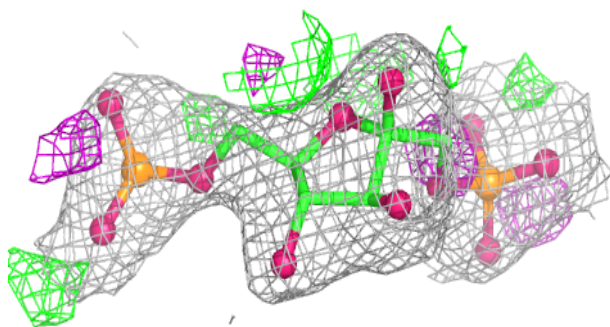
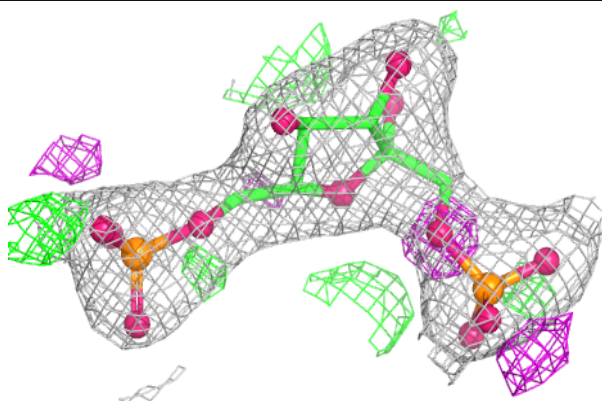


Electron density around I07 D 601 (B):

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

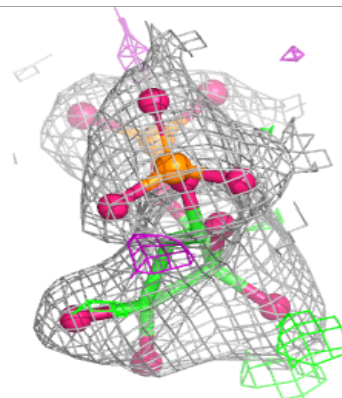
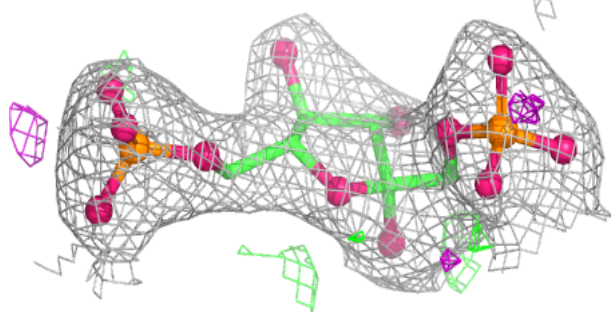
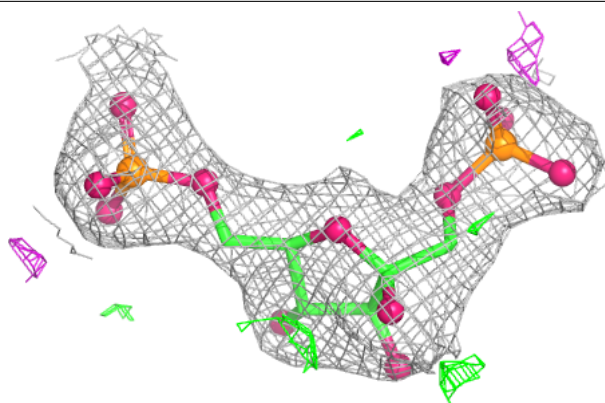
**Electron density around FBP C 601:**

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

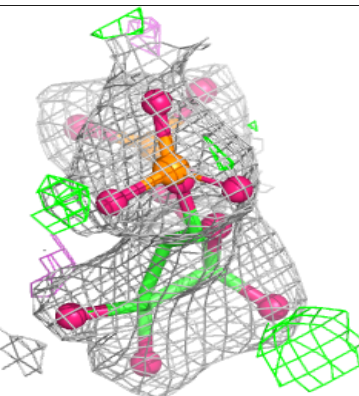
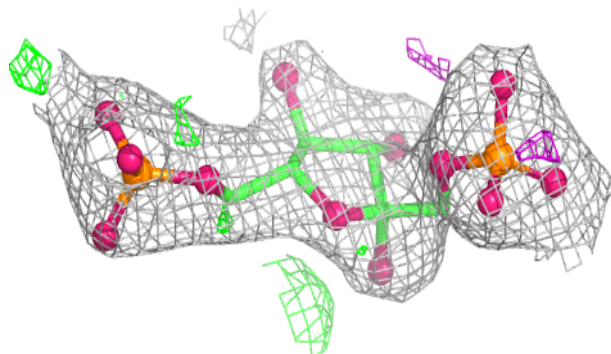
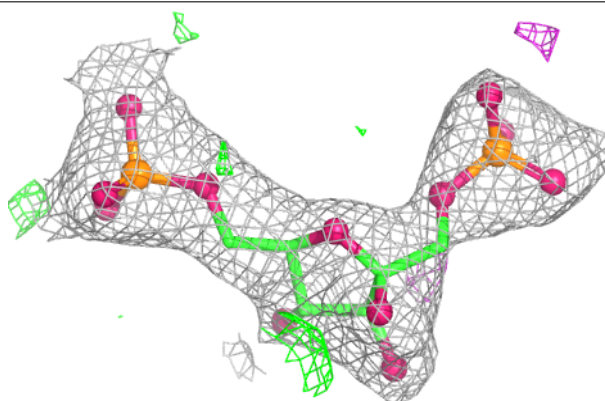


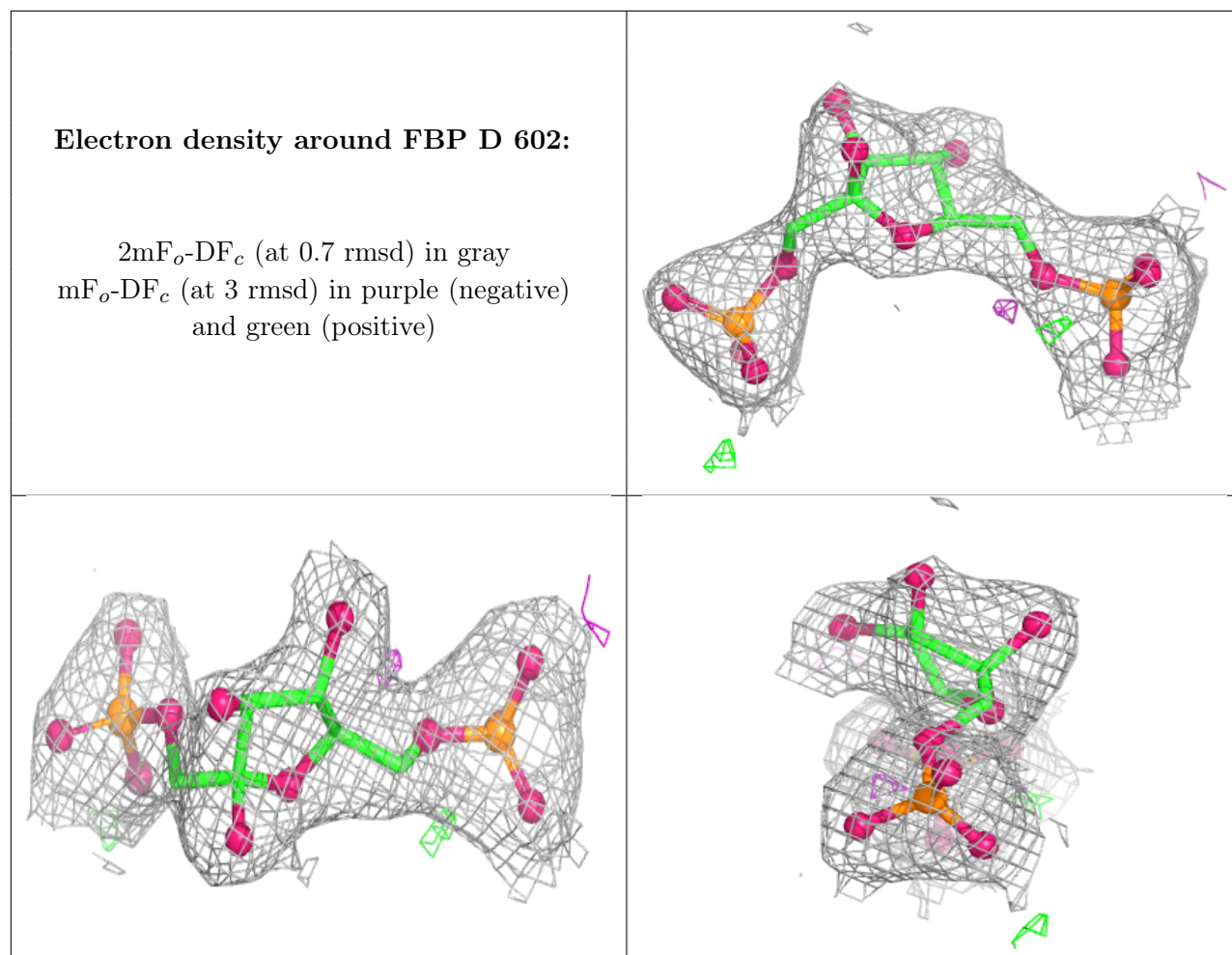
Electron density around FBP B 602:

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

**Electron density around FBP 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)





6.5 Other polymers [i](#)

There are no such residues in this entry.