



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:40 AM UTC

PDB ID : 7FS7 / pdb_00007fs7
Title : Structure of liver pyruvate kinase in complex with allosteric modulator 20
Authors : Lulla, A.; Nilsson, O.; Brear, P.; Nain-Perez, A.; Grotli, M.; Hyvonen, M.
Deposited on : 2022-12-18
Resolution : 2.77 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

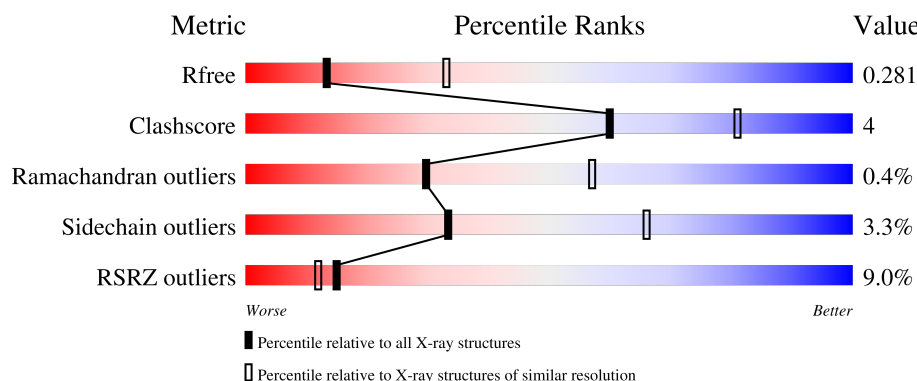
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	447	18% 84% 10% 6%
1	B	447	8% 85% 12% ..
1	C	447	6% 79% 15% • 5%
1	D	447	2% 81% 13% • 5%
1	E	447	19% 82% 11% • 6%

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Mol	Chain	Length	Quality of chain
1	F	447	
1	G	447	
1	H	447	

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	OXL	A	602	-	X	-	-
3	OXL	B	602	-	X	-	-
3	OXL	C	602	-	X	-	-
3	OXL	D	602	-	X	-	-
3	OXL	E	602	-	X	-	-
3	OXL	F	602	-	X	-	-
3	OXL	G	602	-	X	-	-
3	OXL	H	602	-	X	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 26801 atoms, of which 63 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	422	Total	C	N	O	S	0	6	0
			3236	2034	585	597	20			
1	B	436	Total	C	N	O	S	0	4	0
			3329	2090	604	615	20			
1	C	425	Total	C	N	O	S	0	4	0
			3247	2040	585	603	19			
1	D	425	Total	C	N	O	S	0	6	0
			3252	2042	590	601	19			
1	E	419	Total	C	N	O	S	0	5	0
			3210	2018	579	593	20			
1	F	432	Total	C	N	O	S	0	7	0
			3321	2090	597	614	20			
1	G	421	Total	C	N	O	S	0	6	0
			3231	2031	581	600	19			
1	H	425	Total	C	N	O	S	0	4	0
			3251	2040	594	598	19			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P30613
A	0	SER	-	expression tag	UNP P30613
A	1	MET	-	expression tag	UNP P30613
A	2	GLU	-	expression tag	UNP P30613
A	12	ASP	SER	conflict	UNP P30613
A	130	GLY	-	linker	UNP P30613
A	131	SER	-	linker	UNP P30613
A	132	GLY	-	linker	UNP P30613
B	-1	GLY	-	expression tag	UNP P30613
B	0	SER	-	expression tag	UNP P30613
B	1	MET	-	expression tag	UNP P30613
B	2	GLU	-	expression tag	UNP P30613
B	12	ASP	SER	conflict	UNP P30613

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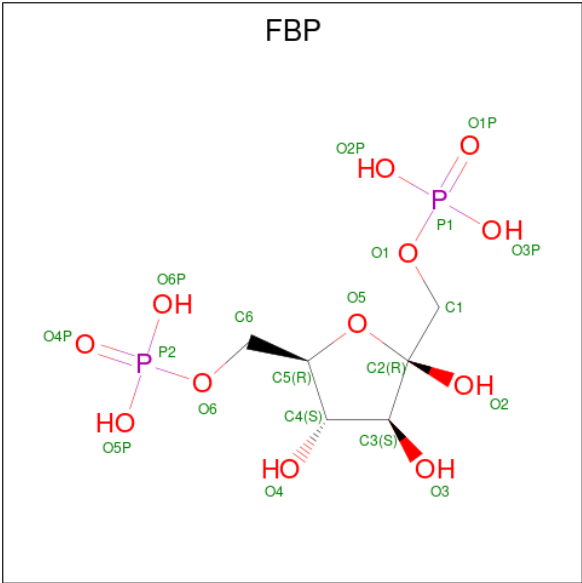
Chain	Residue	Modelled	Actual	Comment	Reference
B	130	GLY	-	linker	UNP P30613
B	131	SER	-	linker	UNP P30613
B	132	GLY	-	linker	UNP P30613
C	-1	GLY	-	expression tag	UNP P30613
C	0	SER	-	expression tag	UNP P30613
C	1	MET	-	expression tag	UNP P30613
C	2	GLU	-	expression tag	UNP P30613
C	12	ASP	SER	conflict	UNP P30613
C	130	GLY	-	linker	UNP P30613
C	131	SER	-	linker	UNP P30613
C	132	GLY	-	linker	UNP P30613
D	-1	GLY	-	expression tag	UNP P30613
D	0	SER	-	expression tag	UNP P30613
D	1	MET	-	expression tag	UNP P30613
D	2	GLU	-	expression tag	UNP P30613
D	12	ASP	SER	conflict	UNP P30613
D	130	GLY	-	linker	UNP P30613
D	131	SER	-	linker	UNP P30613
D	132	GLY	-	linker	UNP P30613
E	-1	GLY	-	expression tag	UNP P30613
E	0	SER	-	expression tag	UNP P30613
E	1	MET	-	expression tag	UNP P30613
E	2	GLU	-	expression tag	UNP P30613
E	12	ASP	SER	conflict	UNP P30613
E	228	GLY	-	linker	UNP P30613
E	229	SER	-	linker	UNP P30613
E	230	GLY	-	linker	UNP P30613
F	-1	GLY	-	expression tag	UNP P30613
F	0	SER	-	expression tag	UNP P30613
F	1	MET	-	expression tag	UNP P30613
F	2	GLU	-	expression tag	UNP P30613
F	12	ASP	SER	conflict	UNP P30613
F	228	GLY	-	linker	UNP P30613
F	229	SER	-	linker	UNP P30613
F	230	GLY	-	linker	UNP P30613
G	-1	GLY	-	expression tag	UNP P30613
G	0	SER	-	expression tag	UNP P30613
G	1	MET	-	expression tag	UNP P30613
G	2	GLU	-	expression tag	UNP P30613
G	12	ASP	SER	conflict	UNP P30613
G	130	GLY	-	linker	UNP P30613
G	131	SER	-	linker	UNP P30613

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Chain	Residue	Modelled	Actual	Comment	Reference
G	132	GLY	-	linker	UNP P30613
H	-1	GLY	-	expression tag	UNP P30613
H	0	SER	-	expression tag	UNP P30613
H	1	MET	-	expression tag	UNP P30613
H	2	GLU	-	expression tag	UNP P30613
H	12	ASP	SER	conflict	UNP P30613
H	130	GLY	-	linker	UNP P30613
H	131	SER	-	linker	UNP P30613
H	132	GLY	-	linker	UNP P30613

- Molecule 2 is 1,6-di-O-phosphono-beta-D-fructofuranose (CCD ID: FBP) (formula: C₆H₁₄O₁₂P₂).



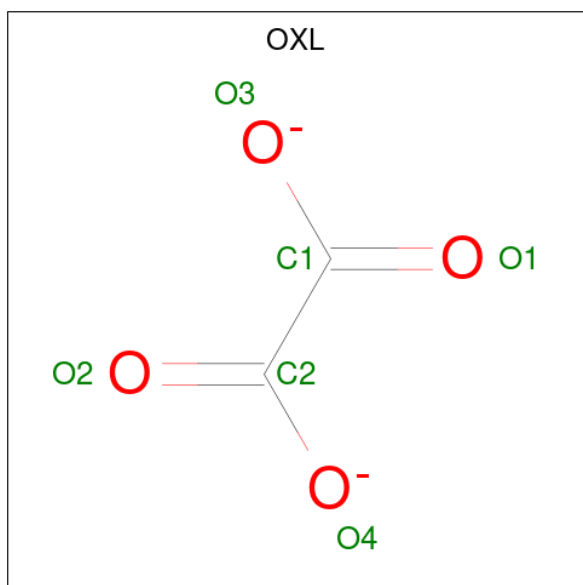
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			20	6	12	2		
2	B	1	Total	C	O	P	0	0
			20	6	12	2		
2	C	1	Total	C	O	P	0	0
			20	6	12	2		
2	D	1	Total	C	O	P	0	0
			20	6	12	2		
2	E	1	Total	C	O	P	0	0
			20	6	12	2		
2	F	1	Total	C	O	P	0	0
			20	6	12	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	G	1	Total	C	O	P	0	0
			20	6	12	2		
2	H	1	Total	C	O	P	0	0
			20	6	12	2		

- Molecule 3 is OXALATE ION (CCD ID: OXL) (formula: C_2O_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	2	4		
3	B	1	Total	C	O	0	0
			6	2	4		
3	C	1	Total	C	O	0	0
			6	2	4		
3	D	1	Total	C	O	0	0
			6	2	4		
3	E	1	Total	C	O	0	0
			6	2	4		
3	F	1	Total	C	O	0	0
			6	2	4		
3	G	1	Total	C	O	0	0
			6	2	4		
3	H	1	Total	C	O	0	0
			6	2	4		

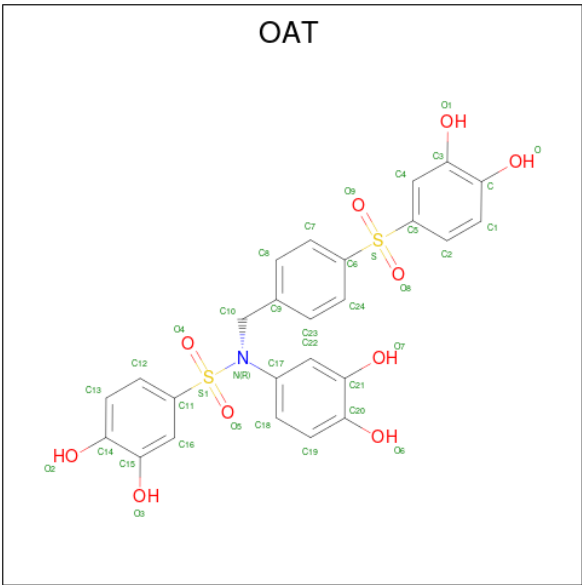
- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mg 1 1	0	0
4	B	1	Total Mg 1 1	0	0
4	C	1	Total Mg 1 1	0	0
4	D	1	Total Mg 1 1	0	0
4	E	1	Total Mg 1 1	0	0
4	F	1	Total Mg 1 1	0	0
4	G	1	Total Mg 1 1	0	0
4	H	1	Total Mg 1 1	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total K 1 1	0	0
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
5	E	1	Total K 1 1	0	0
5	F	1	Total K 1 1	0	0
5	G	1	Total K 1 1	0	0
5	H	1	Total K 1 1	0	0

- Molecule 6 is N-{[4-(3,4-dihydroxybenzene-1-sulfonyl)phenyl]methyl}-N-(3,4-dihydroxyphenyl)-3,4-dihydroxybenzene-1-sulfonamide (CCD ID: OAT) (formula: C₂₅H₂₁NO₁₀S₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	B	1	Total	C	H	N	O	S	21	0
			59	25	21	1	10	2		
6	G	1	Total	C	H	N	O	S	21	0
			59	25	21	1	10	2		
6	H	1	Total	C	H	N	O	S	21	0
			59	25	21	1	10	2		

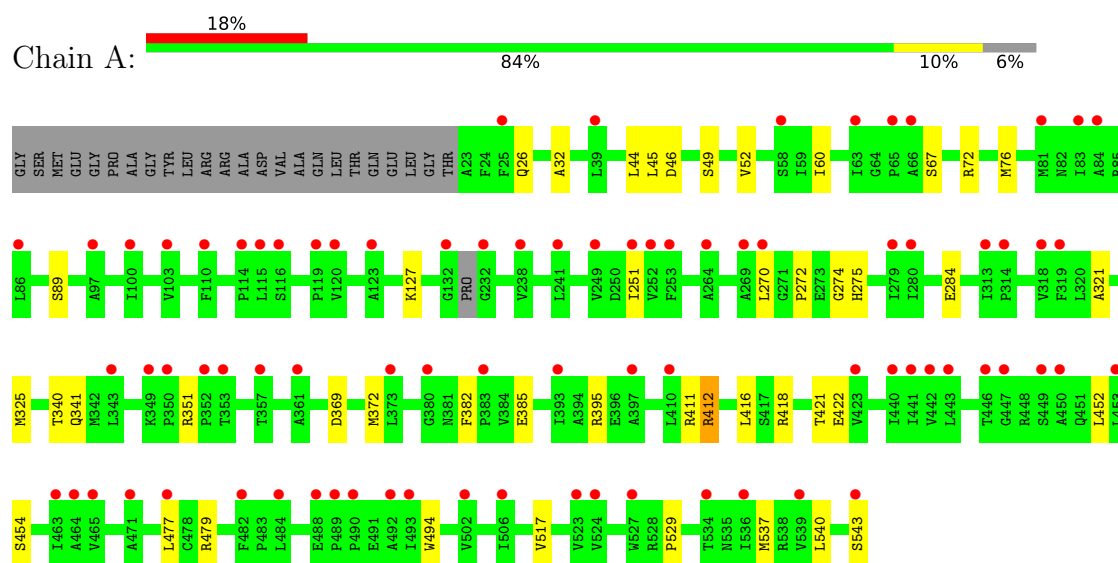
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	5	Total	O	0	0
			5	5		
7	B	15	Total	O	0	0
			15	15		
7	C	45	Total	O	0	0
			45	45		
7	D	79	Total	O	0	0
			79	79		
7	E	9	Total	O	0	0
			9	9		
7	F	35	Total	O	0	0
			35	35		
7	G	48	Total	O	0	0
			48	48		
7	H	87	Total	O	0	0
			87	87		

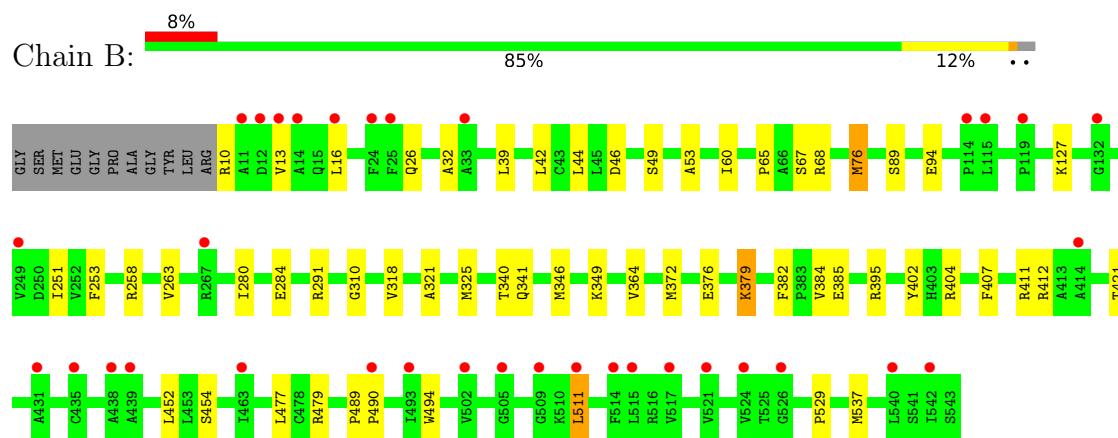
3 Residue-property plots [i](#)

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

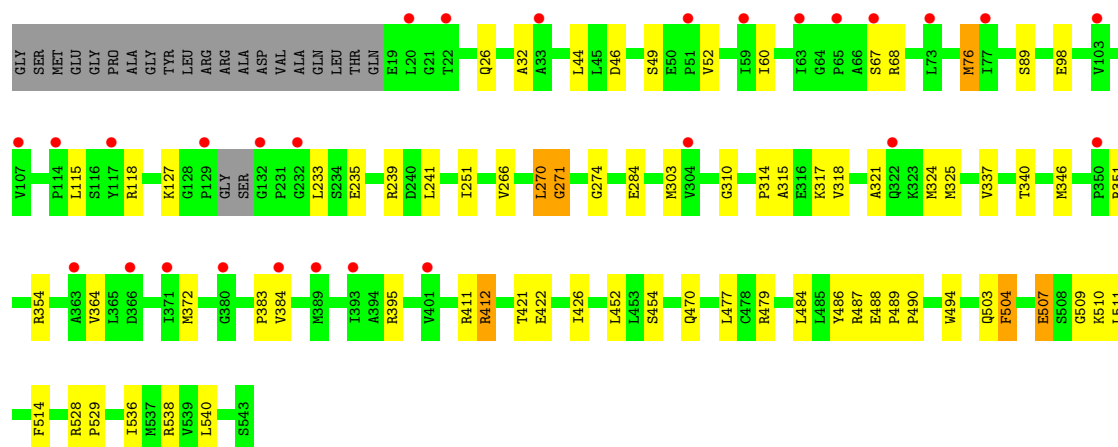


- Molecule 1: Pyruvate kinase PKLR

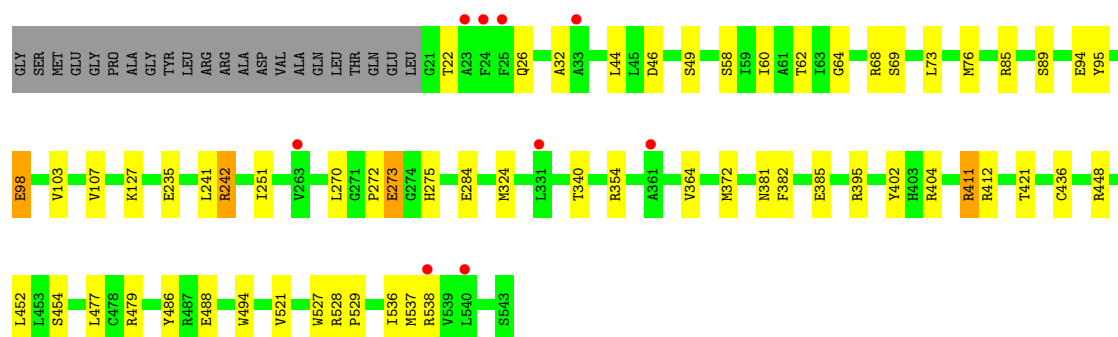
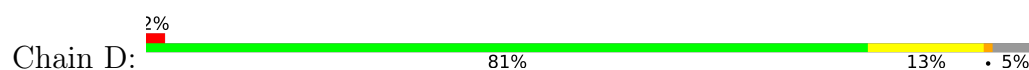


- Molecule 1: Pyruvate kinase PKLR

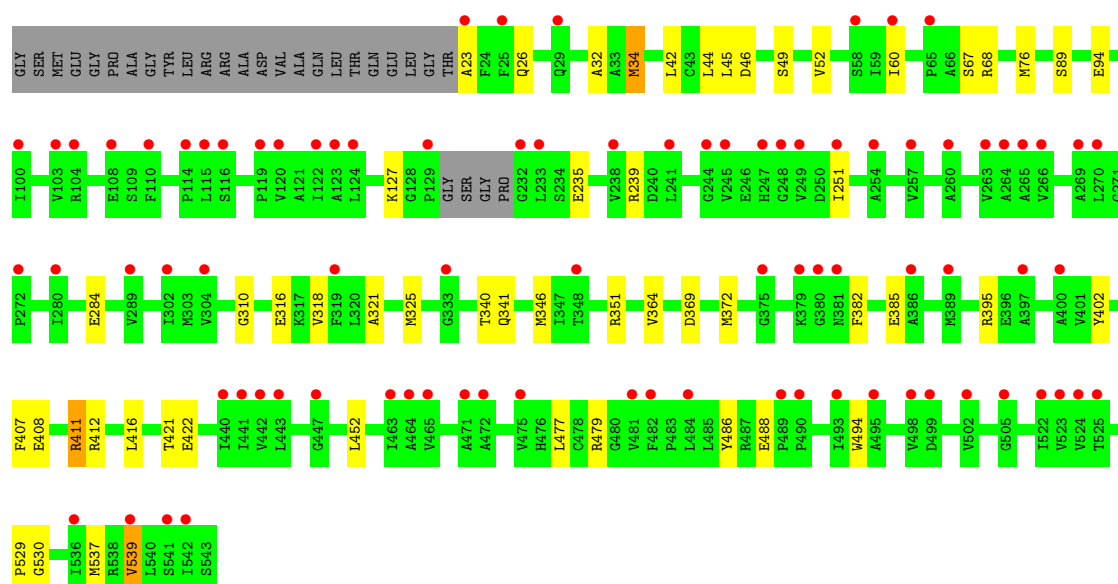
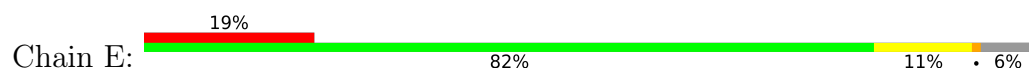




• Molecule 1: Pyruvate kinase PKLR



• Molecule 1: Pyruvate kinase PKLR



• Molecule 1: Pyruvate kinase PKLR

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	206.02Å 113.18Å 187.30Å 90.00° 92.90° 90.00°	Depositor
Resolution (Å)	187.06 – 2.77 187.06 – 2.77	Depositor EDS
% Data completeness (in resolution range)	69.5 (187.06-2.77) 69.5 (187.06-2.77)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 2.77Å)	Xtriage
Refinement program	BUSTER 2.10.4 (16-JUL-2021)	Depositor
R, R_{free}	0.251 , 0.301 0.243 , 0.281	Depositor DCC
R_{free} test set	3781 reflections (3.47%)	wwPDB-VP
Wilson B-factor (Å ²)	80.3	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 91.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.012 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	26801	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 91.37 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.5377e-08. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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, K, FBP, OAT, OXL

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	0.67	0/3307	1.08	1/4469 (0.0%)
1	B	0.70	0/3396	1.07	0/4592
1	C	0.79	0/3313	1.12	4/4479 (0.1%)
1	D	0.83	0/3326	1.13	1/4497 (0.0%)
1	E	0.68	0/3278	1.08	2/4430 (0.0%)
1	F	0.71	1/3396 (0.0%)	1.08	3/4591 (0.1%)
1	G	0.78	1/3303 (0.0%)	1.11	2/4465 (0.0%)
1	H	0.82	0/3316	1.11	4/4483 (0.1%)
All	All	0.75	2/26635 (0.0%)	1.10	17/36006 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	116	SER	CA-C	5.05	1.60	1.52
1	F	313	ILE	CA-C	5.01	1.56	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	422	GLU	CB-CG-CD	6.88	124.29	112.60
1	A	422	GLU	CB-CG-CD	6.58	123.79	112.60
1	C	270	LEU	CA-C-N	5.85	131.05	121.87
1	C	270	LEU	C-N-CA	5.85	131.05	121.87
1	G	270	LEU	CA-C-N	5.68	130.80	121.87
1	G	270	LEU	C-N-CA	5.68	130.80	121.87
1	E	539	VAL	N-CA-CB	5.57	117.34	111.00
1	H	385	GLU	CB-CG-CD	5.52	121.98	112.60
1	H	311	ILE	CB-CA-C	5.41	115.82	111.06
1	F	408	GLU	CB-CG-CD	-5.25	103.67	112.60
1	H	411	ARG	CA-C-N	5.13	127.12	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	411	ARG	C-N-CA	5.13	127.12	120.44
1	F	27	GLN	CA-C-N	5.08	129.62	122.46
1	F	27	GLN	C-N-CA	5.08	129.62	122.46
1	C	504	PHE	CA-C-N	5.04	125.43	119.94
1	C	504	PHE	C-N-CA	5.04	125.43	119.94
1	D	381	ASN	CA-CB-CG	5.01	117.61	112.60

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	3236	0	3299	23	0
1	B	3329	0	3394	30	1
1	C	3247	0	3299	45	0
1	D	3252	0	3310	30	0
1	E	3210	0	3270	31	0
1	F	3321	0	3393	30	0
1	G	3231	0	3284	48	0
1	H	3251	0	3306	37	0
2	A	20	0	10	0	0
2	B	20	0	10	0	0
2	C	20	0	10	0	0
2	D	20	0	10	0	0
2	E	20	0	10	1	0
2	F	20	0	10	0	0
2	G	20	0	10	0	0
2	H	20	0	10	0	0
3	A	6	0	0	0	0
3	B	6	0	0	0	0
3	C	6	0	0	0	0
3	D	6	0	0	0	0
3	E	6	0	0	0	0
3	F	6	0	0	0	0
3	G	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	6	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	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
5	E	1	0	0	0	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
5	H	1	0	0	0	0
6	B	38	21	0	4	0
6	G	38	21	0	3	0
6	H	38	21	0	0	0
7	A	5	0	0	0	0
7	B	15	0	0	0	0
7	C	45	0	0	0	0
7	D	79	0	0	1	0
7	E	9	0	0	0	0
7	F	35	0	0	0	0
7	G	48	0	0	1	0
7	H	87	0	0	4	0
All	All	26738	63	26635	228	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (228) 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:H:449:SER:HB2	7:H:737:HOH:O	1.65	0.95
1:C:411:ARG:HG3	1:C:426:ILE:HD11	1.57	0.86
1:H:442:VAL:HG22	7:H:737:HOH:O	1.76	0.86
1:H:444:THR:HB	7:H:737:HOH:O	1.75	0.85
1:C:528:ARG:HH22	1:G:233:LEU:HB3	1.48	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:402:TYR:HB3	6:G:605:OAT:O5	1.84	0.78
1:E:235:GLU:O	1:E:239:ARG:HD3	1.87	0.73
1:D:242[A]:ARG:NH2	1:D:273:GLU:OE1	2.22	0.72
1:A:517:VAL:HG22	1:A:543:SER:HB3	1.73	0.70
1:C:351:ARG:HH22	1:C:354:ARG:HH22	1.38	0.69
1:D:94:GLU:O	1:D:98:GLU:HG3	1.93	0.69
1:G:411:ARG:HH22	1:H:411:ARG:NH2	1.89	0.69
6:B:605:OAT:O5	1:D:402:TYR:HB3	1.92	0.68
1:H:94:GLU:O	1:H:98:GLU:HG3	1.94	0.67
1:H:272:PRO:HA	1:H:275[A]:HIS:CE1	2.30	0.67
1:D:89:SER:HA	1:D:127:LYS:HG3	1.78	0.66
1:D:272:PRO:HA	1:D:275[A]:HIS:CE1	2.31	0.65
1:C:270:LEU:O	1:C:271:GLY:O	2.14	0.65
1:A:89:SER:HA	1:A:127:LYS:HG3	1.78	0.64
1:H:89:SER:HA	1:H:127:LYS:HG3	1.79	0.64
1:G:270:LEU:O	1:G:271:GLY:O	2.14	0.64
1:E:89:SER:HA	1:E:127:LYS:HG3	1.78	0.64
1:C:89:SER:HA	1:C:127:LYS:HG3	1.79	0.64
1:D:64:GLY:O	1:D:68:ARG:HG3	1.98	0.63
1:F:89:SER:HA	1:F:127:LYS:HG3	1.79	0.63
1:C:235:GLU:CD	1:G:494:TRP:HB2	2.24	0.62
1:C:411:ARG:HH12	1:D:411:ARG:NH2	1.97	0.62
1:B:89:SER:HA	1:B:127:LYS:HG3	1.80	0.62
1:G:89:SER:HA	1:G:127:LYS:HG3	1.80	0.62
1:E:382:PHE:HB3	1:E:385:GLU:HB2	1.82	0.61
1:G:422[A]:GLU:HG3	1:G:452:LEU:HD13	1.82	0.61
1:C:422[A]:GLU:HG3	1:C:452:LEU:HD13	1.82	0.61
1:D:251:ILE:HG12	1:D:477:LEU:HD11	1.83	0.61
1:F:382:PHE:HB3	1:F:385:GLU:HB2	1.81	0.61
1:A:382:PHE:HB3	1:A:385:GLU:HB2	1.84	0.59
1:A:412:ARG:NH1	1:B:404:ARG:HH11	2.00	0.59
1:C:528:ARG:NH2	1:G:233:LEU:HB3	2.17	0.59
1:B:382:PHE:HB3	1:B:385:GLU:HB2	1.84	0.59
1:A:270:LEU:HB3	1:A:274:GLY:HA3	1.85	0.58
1:E:34:MET:HE2	1:E:34:MET:HA	1.85	0.58
1:G:241:LEU:HD22	1:G:270:LEU:HD21	1.86	0.58
1:A:411:ARG:NH2	1:B:411:ARG:HH22	2.01	0.57
1:C:251:ILE:HG12	1:C:477:LEU:HD11	1.87	0.56
1:D:103:VAL:O	1:D:107:VAL:HG23	2.06	0.56
1:F:408:GLU:HG2	1:F:411:ARG:NH2	2.20	0.56
1:A:411:ARG:HH21	1:B:411:ARG:HH22	1.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:407:PHE:CE2	1:B:411:ARG:NH1	2.75	0.55
1:G:411:ARG:NH2	1:H:411:ARG:NH2	2.54	0.54
1:E:351:ARG:HH21	1:G:315:ALA:HB2	1.72	0.54
1:D:382:PHE:HB3	1:D:385:GLU:HB2	1.90	0.54
1:C:503:GLN:O	1:C:507:GLU:HG2	2.09	0.53
1:F:407:PHE:CE2	1:F:411:ARG:NH1	2.76	0.53
1:C:412:ARG:NH1	1:D:404:ARG:HD3	2.24	0.53
1:H:251:ILE:HG12	1:H:477:LEU:HD11	1.91	0.52
1:C:421:THR:HG22	1:C:452:LEU:HD12	1.90	0.52
1:D:46:ASP:HB3	1:D:49:SER:HB2	1.92	0.52
1:B:341:GLN:OE1	1:D:354:ARG:HG2	2.10	0.52
1:C:239:ARG:HH21	1:G:494:TRP:HD1	1.57	0.52
1:C:115:LEU:O	1:C:118:ARG:NH1	2.43	0.52
1:D:241:LEU:HD22	1:D:270:LEU:HD21	1.92	0.51
1:H:267[B]:ARG:NH1	1:H:274:GLY:O	2.42	0.51
1:H:382:PHE:HB3	1:H:385:GLU:HB2	1.91	0.51
1:C:536:ILE:HG12	1:D:538:ARG:HG3	1.93	0.51
1:H:46:ASP:HB3	1:H:49:SER:HB2	1.93	0.51
1:A:272:PRO:HA	1:A:275:HIS:NE2	2.26	0.51
1:G:251:ILE:HG12	1:G:477:LEU:HD11	1.92	0.50
1:H:411:ARG:HG2	1:H:426:ILE:HD11	1.91	0.50
1:B:258:ARG:O	1:B:291:ARG:HD3	2.11	0.50
1:E:67:SER:HB2	1:E:76[A]:MET:SD	2.52	0.50
1:F:26:GLN:O	1:F:49:SER:OG	2.28	0.50
1:G:421:THR:HG22	1:G:452:LEU:HD12	1.94	0.50
1:H:259:LYS:NZ	7:H:702:HOH:O	2.40	0.50
1:B:26:GLN:O	1:B:49:SER:OG	2.27	0.50
1:C:46:ASP:HB3	1:C:49:SER:HB2	1.94	0.50
1:A:26:GLN:O	1:A:49:SER:OG	2.30	0.50
1:A:341:GLN:OE1	1:C:354:ARG:HG2	2.11	0.50
1:E:251:ILE:HG12	1:E:477:LEU:HD11	1.93	0.49
1:A:67:SER:HB2	1:A:76[A]:MET:SD	2.52	0.49
1:C:233:LEU:O	1:G:528:ARG:NH2	2.32	0.49
1:G:46:ASP:HB3	1:G:49:SER:HB2	1.93	0.49
1:B:251:ILE:HG12	1:B:477:LEU:HD11	1.94	0.49
1:D:26:GLN:O	1:D:49:SER:OG	2.30	0.49
1:E:421:THR:HG22	1:E:452:LEU:HD12	1.94	0.49
1:B:65:PRO:HG2	1:B:379:LYS:HG2	1.94	0.49
1:A:251:ILE:HG12	1:A:477:LEU:HD11	1.94	0.49
1:A:421:THR:HG22	1:A:452:LEU:HD12	1.94	0.49
1:A:418[A]:ARG:HG3	1:B:16:LEU:HD11	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:ASP:HB3	1:B:49:SER:HB2	1.95	0.48
1:D:68:ARG:NH2	1:D:95:TYR:O	2.46	0.48
1:F:251:ILE:HG12	1:F:477:LEU:HD11	1.94	0.48
1:F:46:ASP:HB3	1:F:49:SER:HB2	1.95	0.48
1:G:538:ARG:HG2	1:H:536:ILE:HG12	1.95	0.48
1:G:60:ILE:HB	1:G:372:MET:HG3	1.96	0.48
1:G:67:SER:HB2	1:G:76[B]:MET:SD	2.53	0.48
1:G:537:MET:HE3	1:H:537:MET:HG2	1.96	0.48
1:E:23:ALA:HB3	1:E:26:GLN:HG2	1.95	0.48
1:F:408:GLU:HG2	1:F:411:ARG:HH22	1.79	0.48
1:H:421:THR:HG22	1:H:452:LEU:HD12	1.96	0.48
1:C:241:LEU:HD22	1:C:270:LEU:HD21	1.95	0.48
1:G:26:GLN:O	1:G:49:SER:OG	2.28	0.48
1:G:538:ARG:CG	1:H:536:ILE:HG12	2.43	0.48
1:C:60:ILE:HB	1:C:372:MET:HG3	1.96	0.47
1:E:408:GLU:CD	1:F:411:ARG:HH21	2.21	0.47
1:G:33:ALA:HB2	7:G:703:HOH:O	2.13	0.47
1:F:67:SER:HB2	1:F:76[B]:MET:SD	2.55	0.47
1:A:46:ASP:HB3	1:A:49:SER:HB2	1.97	0.47
1:B:67:SER:HB2	1:B:76[B]:MET:SD	2.55	0.47
1:C:67:SER:HB2	1:C:76[B]:MET:SD	2.55	0.47
1:E:408:GLU:OE1	1:F:411:ARG:NH2	2.44	0.47
1:E:412:ARG:HD3	1:F:404:ARG:NH1	2.30	0.47
1:G:411:ARG:HG3	1:G:426:ILE:HD11	1.97	0.47
1:B:402:TYR:HB3	6:B:605:OAT:O9	2.14	0.46
1:G:527:TRP:CD2	1:G:528:ARG:HG2	2.50	0.46
1:F:421:THR:HG22	1:F:452:LEU:HD12	1.97	0.46
1:E:46:ASP:HB3	1:E:49:SER:HB2	1.96	0.46
1:D:421:THR:HG22	1:D:452:LEU:HD12	1.97	0.46
1:F:32:ALA:HA	1:F:44:LEU:HB2	1.98	0.46
1:B:421:THR:HG22	1:B:452:LEU:HD12	1.97	0.46
1:C:239:ARG:HH21	1:G:494:TRP:CD1	2.34	0.46
1:A:494:TRP:CD1	1:A:529:PRO:HG3	2.51	0.45
1:C:486:TYR:CZ	1:C:488:GLU:HB2	2.51	0.45
1:A:60:ILE:HB	1:A:372:MET:HG3	1.98	0.45
1:C:68:ARG:NH2	1:C:98:GLU:HB3	2.30	0.45
1:B:60:ILE:HB	1:B:372:MET:HG3	1.98	0.45
1:G:536:ILE:HG23	1:H:538[A]:ARG:HG2	1.98	0.45
1:B:346:MET:HA	1:B:349:LYS:O	2.16	0.45
1:C:529:PRO:HB2	1:G:238:VAL:HG21	1.97	0.45
1:D:60:ILE:HB	1:D:372:MET:HG3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:448:ARG:HD3	7:D:732:HOH:O	2.16	0.45
1:C:32:ALA:HA	1:C:44:LEU:HB2	1.98	0.45
1:H:68:ARG:NH1	1:H:98:GLU:OE1	2.45	0.45
1:G:39:LEU:HB2	6:G:605:OAT:O6	2.17	0.45
1:C:239:ARG:NH2	1:G:494:TRP:CD1	2.85	0.45
1:D:494:TRP:CD1	1:D:529:PRO:HG3	2.53	0.44
1:E:494:TRP:CD1	1:E:529:PRO:HG3	2.52	0.44
1:A:32:ALA:HA	1:A:44:LEU:HB2	1.99	0.44
1:C:364:VAL:O	1:C:479:ARG:NH1	2.48	0.44
1:C:494:TRP:HB2	1:G:235:GLU:CD	2.43	0.44
1:F:241:LEU:HD22	1:F:270:LEU:HD21	2.00	0.44
1:G:32:ALA:HA	1:G:44:LEU:HB2	1.99	0.44
1:B:32:ALA:HA	1:B:44:LEU:HB2	1.99	0.44
1:C:509:GLY:HA2	1:C:514:PHE:HD2	1.82	0.44
1:D:62:THR:HA	1:D:85:ARG:HB3	1.99	0.44
1:A:351:ARG:HH21	1:C:315:ALA:HB2	1.82	0.44
1:H:60:ILE:HB	1:H:372:MET:HG3	1.98	0.44
1:E:407:PHE:O	1:E:411:ARG:HB2	2.18	0.44
1:H:494:TRP:CD1	1:H:529:PRO:HG3	2.52	0.44
1:B:76[A]:MET:HG2	1:B:384:VAL:HG22	1.99	0.44
1:F:494:TRP:CD1	1:F:529:PRO:HG3	2.53	0.44
1:G:321:ALA:O	1:G:325:MET:HG3	2.18	0.44
1:G:76[A]:MET:HG2	1:G:384:VAL:HG22	1.99	0.43
1:G:272:PRO:HD2	1:G:273:GLU:OE1	2.18	0.43
1:C:321:ALA:O	1:C:325:MET:HG3	2.17	0.43
1:E:341:GLN:OE1	1:G:354:ARG:HG2	2.19	0.43
1:E:416:LEU:HB2	1:F:16:LEU:HD22	2.00	0.43
1:F:364:VAL:O	1:F:479:ARG:NH1	2.48	0.43
1:G:411:ARG:NH2	1:H:411:ARG:CZ	2.81	0.43
1:C:76[A]:MET:HG2	1:C:384:VAL:HG22	1.99	0.43
1:F:76[A]:MET:HG2	1:F:384:VAL:HG22	2.00	0.43
1:H:26:GLN:O	1:H:49:SER:OG	2.29	0.43
1:B:494:TRP:CD1	1:B:529:PRO:HG3	2.53	0.43
1:C:484:LEU:HB3	1:C:504:PHE:CE2	2.53	0.43
1:H:258:ARG:O	1:H:291:ARG:HD3	2.19	0.43
1:G:412:ARG:NH2	1:H:404:ARG:HD3	2.34	0.43
1:D:527:TRP:CD2	1:D:528:ARG:HG2	2.54	0.43
1:F:341:GLN:OE1	1:H:354:ARG:HG2	2.19	0.43
1:G:266:VAL:O	1:G:270:LEU:HG	2.19	0.43
1:F:42:LEU:O	1:H:324:MET:HA	2.19	0.43
1:E:26:GLN:O	1:E:49:SER:OG	2.29	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:32:ALA:HA	1:E:44:LEU:HB2	1.99	0.43
1:C:314:PRO:HG2	1:C:317:LYS:HD2	2.01	0.42
1:D:364:VAL:O	1:D:479:ARG:NH1	2.49	0.42
1:E:45:LEU:HB2	1:G:324:MET:HB2	2.01	0.42
1:E:364:VAL:O	1:E:479:ARG:NH1	2.47	0.42
1:F:55:ARG:HB2	1:F:395:ARG:HG3	2.02	0.42
1:F:60:ILE:HB	1:F:372:MET:HG3	1.99	0.42
1:F:267[B]:ARG:HH21	1:F:270:LEU:HD12	1.83	0.42
1:C:538:ARG:CG	1:D:536:ILE:HG12	2.50	0.42
1:E:60:ILE:HB	1:E:372:MET:HG3	2.00	0.42
1:C:351:ARG:HH22	1:C:354:ARG:NH2	2.12	0.42
1:G:412:ARG:CZ	1:H:404:ARG:HD3	2.50	0.42
1:H:310:GLY:HA2	1:H:318:VAL:HG21	2.01	0.42
1:B:53:ALA:O	1:B:395:ARG:NH2	2.52	0.42
1:B:364:VAL:O	1:B:479:ARG:NH1	2.51	0.42
1:F:310:GLY:HA2	1:F:318:VAL:HG21	2.02	0.42
1:H:67:SER:HB2	1:H:76[A]:MET:SD	2.60	0.42
1:D:73:LEU:HA	1:D:76[B]:MET:HG3	2.01	0.41
1:A:321:ALA:O	1:A:325:MET:HG3	2.20	0.41
1:G:411:ARG:HH22	1:H:411:ARG:HH21	1.65	0.41
1:E:310:GLY:HA2	1:E:318:VAL:HG21	2.03	0.41
1:E:530:GLY:O	2:E:601:FBP:O4	2.29	0.41
1:H:32:ALA:HA	1:H:44:LEU:HB2	2.02	0.41
1:A:369:ASP:HA	1:A:479:ARG:HB2	2.03	0.41
1:B:310:GLY:HA2	1:B:318:VAL:HG21	2.02	0.41
1:F:346:MET:HA	1:F:349:LYS:O	2.20	0.41
1:E:316:GLU:HB2	1:G:393:ILE:HA	2.02	0.41
1:E:321:ALA:O	1:E:325:MET:HG3	2.20	0.41
1:C:489:PRO:HA	1:C:490:PRO:HD3	1.92	0.41
1:E:486:TYR:CE2	1:E:488:GLU:HB2	2.56	0.41
1:F:321:ALA:O	1:F:325:MET:HG3	2.20	0.41
1:H:407:PHE:CE2	1:H:411:ARG:HD3	2.56	0.41
1:B:39:LEU:HB2	6:B:605:OAT:O6	2.20	0.41
1:B:321:ALA:O	1:B:325:MET:HG3	2.21	0.41
1:C:266:VAL:O	1:C:270:LEU:HG	2.21	0.41
1:C:310:GLY:HA2	1:C:318:VAL:HG21	2.03	0.41
1:D:486:TYR:CE2	1:D:488:GLU:HB2	2.56	0.41
1:E:369:ASP:HA	1:E:479:ARG:HB2	2.03	0.41
1:G:402:TYR:HB3	6:G:605:OAT:O8	2.21	0.41
1:A:45:LEU:HB2	1:C:324:MET:HB2	2.02	0.41
6:B:605:OAT:C22	6:B:605:OAT:O4	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:364:VAL:O	1:G:479:ARG:NH1	2.50	0.41
1:G:422[B]:GLU:HG3	1:H:434:LYS:HE2	2.03	0.41
1:H:346:MET:HE2	1:H:346:MET:HB3	1.96	0.41
1:B:489:PRO:HA	1:B:490:PRO:HD3	1.98	0.41
1:F:354:ARG:HG2	1:H:341:GLN:OE1	2.21	0.41
1:G:252:VAL:HG23	1:G:277:ILE:HG21	2.03	0.41
1:G:369:ASP:HA	1:G:479:ARG:HB2	2.03	0.41
1:B:42:LEU:O	1:D:324:MET:HA	2.21	0.40
1:B:253:PHE:HD1	1:B:280:ILE:HB	1.86	0.40
1:A:416:LEU:CB	1:B:16:LEU:HD22	2.52	0.40
1:C:303:MET:HG3	1:C:337:VAL:HB	2.04	0.40
1:D:68:ARG:HD2	1:D:95:TYR:OH	2.21	0.40
1:F:489:PRO:HA	1:F:490:PRO:HD3	1.99	0.40
1:C:26:GLN:O	1:C:49:SER:OG	2.28	0.40
1:D:32:ALA:HA	1:D:44:LEU:HB2	2.02	0.40
1:E:42:LEU:O	1:G:324:MET:HA	2.21	0.40
1:E:416:LEU:CB	1:F:16:LEU:HD22	2.52	0.40
1:C:346:MET:HE2	1:C:346:MET:HB3	2.00	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:511:LEU:O	1:B:511:LEU:O[2_555]	2.07	0.13

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	424/447 (95%)	415 (98%)	8 (2%)	1 (0%)	43 70
1	B	438/447 (98%)	427 (98%)	10 (2%)	1 (0%)	43 70

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	425/447 (95%)	412 (97%)	9 (2%)	4 (1%)	14	37
1	D	429/447 (96%)	421 (98%)	7 (2%)	1 (0%)	43	70
1	E	420/447 (94%)	412 (98%)	7 (2%)	1 (0%)	43	70
1	F	435/447 (97%)	424 (98%)	10 (2%)	1 (0%)	43	70
1	G	423/447 (95%)	412 (97%)	8 (2%)	3 (1%)	18	44
1	H	427/447 (96%)	416 (97%)	10 (2%)	1 (0%)	43	70
All	All	3421/3576 (96%)	3339 (98%)	69 (2%)	13 (0%)	30	57

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	271	GLY
1	G	271	GLY
1	A	340	THR
1	D	340	THR
1	B	340	THR
1	C	274	GLY
1	C	340	THR
1	E	340	THR
1	F	340	THR
1	G	340	THR
1	H	340	THR
1	C	383	PRO
1	G	383	PRO

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	340/352 (97%)	331 (97%)	9 (3%)	40	72
1	B	349/352 (99%)	334 (96%)	15 (4%)	26	57
1	C	341/352 (97%)	328 (96%)	13 (4%)	29	61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	342/352 (97%)	325 (95%)	17 (5%)	22	51
1	E	338/352 (96%)	327 (97%)	11 (3%)	33	66
1	F	350/352 (99%)	337 (96%)	13 (4%)	30	62
1	G	340/352 (97%)	328 (96%)	12 (4%)	32	64
1	H	340/352 (97%)	330 (97%)	10 (3%)	37	70
All	All	2740/2816 (97%)	2640 (96%)	100 (4%)	33	63

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	VAL
1	A	72	ARG
1	A	284	GLU
1	A	395	ARG
1	A	412	ARG
1	A	454	SER
1	A	537[A]	MET
1	A	537[B]	MET
1	A	540	LEU
1	B	10	ARG
1	B	13	VAL
1	B	68	ARG
1	B	76[A]	MET
1	B	76[B]	MET
1	B	94	GLU
1	B	263	VAL
1	B	284	GLU
1	B	376	GLU
1	B	379	LYS
1	B	412	ARG
1	B	454	SER
1	B	511	LEU
1	B	537[A]	MET
1	B	537[B]	MET
1	C	52	VAL
1	C	76[A]	MET
1	C	76[B]	MET
1	C	284	GLU
1	C	395	ARG
1	C	412	ARG

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Mol	Chain	Res	Type
1	C	454	SER
1	C	470	GLN
1	C	487	ARG
1	C	507	GLU
1	C	510	LYS
1	C	511	LEU
1	C	540	LEU
1	D	22	THR
1	D	58[A]	SER
1	D	58[B]	SER
1	D	69	SER
1	D	98	GLU
1	D	235	GLU
1	D	242[A]	ARG
1	D	242[B]	ARG
1	D	273	GLU
1	D	284	GLU
1	D	395	ARG
1	D	411	ARG
1	D	412	ARG
1	D	436	CYS
1	D	454	SER
1	D	521	VAL
1	D	537	MET
1	E	34	MET
1	E	52	VAL
1	E	68	ARG
1	E	94	GLU
1	E	284	GLU
1	E	346	MET
1	E	395	ARG
1	E	411	ARG
1	E	537[A]	MET
1	E	537[B]	MET
1	E	539	VAL
1	F	13	VAL
1	F	16	LEU
1	F	68	ARG
1	F	76[A]	MET
1	F	76[B]	MET
1	F	118	ARG
1	F	242	ARG

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Mol	Chain	Res	Type
1	F	284	GLU
1	F	412	ARG
1	F	454	SER
1	F	511	LEU
1	F	537[A]	MET
1	F	537[B]	MET
1	G	52	VAL
1	G	68	ARG
1	G	76[A]	MET
1	G	76[B]	MET
1	G	113	SER
1	G	273	GLU
1	G	284	GLU
1	G	331	LEU
1	G	395	ARG
1	G	412	ARG
1	G	454	SER
1	G	540	LEU
1	H	22	THR
1	H	98	GLU
1	H	235	GLU
1	H	273	GLU
1	H	284	GLU
1	H	395	ARG
1	H	411	ARG
1	H	412	ARG
1	H	454	SER
1	H	537	MET

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

Mol	Chain	Res	Type
1	A	236	GLN
1	C	470	GLN
1	D	236	GLN
1	F	15	GLN
1	H	236	GLN

5.3.3 RNA ⓘ

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 35 ligands modelled in this entry, 16 are monoatomic - leaving 19 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	OXL	B	602	4	5,5,5	2.11	2 (40%)	6,6,6	1.64	2 (33%)
3	OXL	A	602	4	5,5,5	2.11	2 (40%)	6,6,6	1.82	1 (16%)
3	OXL	H	602	4	5,5,5	2.21	2 (40%)	6,6,6	1.71	1 (16%)
2	FBP	E	601	-	18,20,20	0.51	0	21,32,32	0.75	0
6	OAT	G	605	-	41,41,41	0.35	0	60,62,62	0.73	2 (3%)
6	OAT	H	605	-	41,41,41	0.32	0	60,62,62	0.63	2 (3%)
2	FBP	F	601	-	18,20,20	0.58	0	21,32,32	0.62	0
3	OXL	G	602	4	5,5,5	2.13	2 (40%)	6,6,6	1.79	2 (33%)
2	FBP	D	601	-	18,20,20	0.79	1 (5%)	21,32,32	0.99	1 (4%)
3	OXL	E	602	4	5,5,5	2.30	3 (60%)	6,6,6	1.83	2 (33%)
3	OXL	C	602	4	5,5,5	2.03	2 (40%)	6,6,6	1.92	2 (33%)
2	FBP	B	601	-	18,20,20	0.73	1 (5%)	21,32,32	0.79	0
2	FBP	A	601	-	18,20,20	0.61	0	21,32,32	0.66	0
6	OAT	B	605	-	41,41,41	0.39	0	60,62,62	0.67	2 (3%)
2	FBP	C	601	-	18,20,20	0.86	1 (5%)	21,32,32	1.23	1 (4%)
3	OXL	D	602	4	5,5,5	2.27	2 (40%)	6,6,6	1.80	2 (33%)
2	FBP	G	601	-	18,20,20	0.66	1 (5%)	21,32,32	1.16	1 (4%)
2	FBP	H	601	-	18,20,20	0.56	0	21,32,32	1.37	2 (9%)
3	OXL	F	602	4	5,5,5	2.83	4 (80%)	6,6,6	2.18	3 (50%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	OXL	B	602	4	-	4/4/4/4	-
3	OXL	A	602	4	-	4/4/4/4	-
3	OXL	H	602	4	-	4/4/4/4	-
2	FBP	E	601	-	-	3/13/32/32	0/1/1/1
6	OAT	G	605	-	-	11/32/32/32	0/4/4/4
6	OAT	H	605	-	-	3/32/32/32	0/4/4/4
2	FBP	F	601	-	-	2/13/32/32	0/1/1/1
3	OXL	G	602	4	-	4/4/4/4	-
2	FBP	D	601	-	-	2/13/32/32	0/1/1/1
3	OXL	E	602	4	-	4/4/4/4	-
3	OXL	C	602	4	-	4/4/4/4	-
2	FBP	B	601	-	-	2/13/32/32	0/1/1/1
2	FBP	A	601	-	-	3/13/32/32	0/1/1/1
6	OAT	B	605	-	-	9/32/32/32	0/4/4/4
2	FBP	C	601	-	-	7/13/32/32	0/1/1/1
3	OXL	D	602	4	-	4/4/4/4	-
2	FBP	G	601	-	-	2/13/32/32	0/1/1/1
2	FBP	H	601	-	-	3/13/32/32	0/1/1/1
3	OXL	F	602	4	-	4/4/4/4	-

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	602	OXL	O2-C2	3.92	1.32	1.22
3	E	602	OXL	O1-C1	3.83	1.32	1.22
3	H	602	OXL	O2-C2	3.81	1.32	1.22
3	B	602	OXL	O2-C2	3.73	1.31	1.22
3	A	602	OXL	O2-C2	3.60	1.31	1.22
3	F	602	OXL	O2-C2	3.52	1.31	1.22
3	G	602	OXL	O1-C1	3.31	1.30	1.22
3	C	602	OXL	O1-C1	3.30	1.30	1.22
3	F	602	OXL	O3-C1	-3.29	1.21	1.30
3	F	602	OXL	O1-C1	3.21	1.30	1.22
3	G	602	OXL	O3-C1	-3.19	1.21	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	FBP	P1-O3P	-2.95	1.43	1.54
3	C	602	OXL	O3-C1	-2.88	1.22	1.30
2	D	601	FBP	P2-O5P	-2.81	1.44	1.54
3	D	602	OXL	O4-C2	-2.78	1.23	1.30
3	B	602	OXL	O4-C2	-2.77	1.23	1.30
3	E	602	OXL	O3-C1	-2.76	1.23	1.30
3	A	602	OXL	O4-C2	-2.63	1.23	1.30
3	F	602	OXL	O4-C2	-2.58	1.23	1.30
3	H	602	OXL	O4-C2	-2.51	1.23	1.30
2	B	601	FBP	P2-O5P	-2.10	1.47	1.54
2	G	601	FBP	P1-O3P	-2.03	1.47	1.54
3	E	602	OXL	C2-C1	2.02	1.58	1.54

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	601	FBP	O5P-P2-O6	4.39	118.11	106.67
2	C	601	FBP	O6-P2-O4P	3.96	117.15	106.44
3	F	602	OXL	O4-C2-C1	3.65	119.92	112.83
3	A	602	OXL	O4-C2-C1	3.63	119.87	112.83
3	C	602	OXL	O3-C1-C2	3.60	119.82	112.83
6	G	605	OAT	C9-C10-N	3.41	118.80	112.02
3	H	602	OXL	O4-C2-C1	3.40	119.42	112.83
3	G	602	OXL	O3-C1-C2	3.34	119.30	112.83
3	E	602	OXL	O3-C1-C2	3.25	119.14	112.83
3	D	602	OXL	O4-C2-C1	3.10	118.85	112.83
3	B	602	OXL	O4-C2-C1	3.06	118.77	112.83
3	F	602	OXL	O3-C1-C2	3.04	118.72	112.83
2	G	601	FBP	O6P-P2-O6	2.72	113.77	106.67
6	B	605	OAT	C9-C10-N	2.70	117.38	112.02
3	D	602	OXL	O3-C1-C2	2.68	118.03	112.83
3	E	602	OXL	O4-C2-C1	2.64	117.95	112.83
2	H	601	FBP	O6P-P2-O6	-2.61	99.88	106.67
6	G	605	OAT	C17-N-S1	-2.51	112.68	117.72
3	C	602	OXL	O4-C2-C1	2.36	117.41	112.83
3	G	602	OXL	O4-C2-C1	2.32	117.33	112.83
6	H	605	OAT	C9-C10-N	2.18	116.36	112.02
6	B	605	OAT	C10-N-S1	2.17	122.54	117.95
3	B	602	OXL	O3-C1-C2	2.16	117.02	112.83
3	F	602	OXL	O2-C2-C1	-2.10	116.26	120.63
2	D	601	FBP	O1-P1-O1P	2.08	112.07	106.44
6	H	605	OAT	C17-N-S1	-2.01	113.68	117.72

There are no chirality outliers.

All (79) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	601	FBP	C1-O1-P1-O2P
2	C	601	FBP	C1-O1-P1-O3P
2	C	601	FBP	O1-C1-C2-O2
2	C	601	FBP	O1-C1-C2-C3
2	C	601	FBP	O1-C1-C2-O5
2	D	601	FBP	C4-C5-C6-O6
6	G	605	OAT	C16-C11-S1-N
6	G	605	OAT	C17-N-S1-O4
6	G	605	OAT	C12-C11-S1-N
2	H	601	FBP	C4-C5-C6-O6
6	H	605	OAT	C10-N-S1-O5
6	B	605	OAT	C17-N-S1-O4
6	B	605	OAT	C17-N-S1-O5
6	G	605	OAT	C17-N-S1-O5
6	G	605	OAT	C17-N-S1-C11
6	G	605	OAT	C12-C11-S1-O4
2	A	601	FBP	C4-C5-C6-O6
2	B	601	FBP	C4-C5-C6-O6
2	D	601	FBP	O5-C5-C6-O6
2	E	601	FBP	C4-C5-C6-O6
2	F	601	FBP	C4-C5-C6-O6
2	G	601	FBP	C4-C5-C6-O6
6	G	605	OAT	C16-C11-S1-O4
6	B	605	OAT	C22-C17-N-S1
6	H	605	OAT	C10-N-S1-C11
6	B	605	OAT	C17-N-S1-C11
6	B	605	OAT	C12-C11-S1-O4
2	C	601	FBP	C1-O1-P1-O1P
2	E	601	FBP	C6-O6-P2-O4P
6	G	605	OAT	C10-N-S1-O4
6	B	605	OAT	C16-C11-S1-O4
6	B	605	OAT	C12-C11-S1-N
6	B	605	OAT	C16-C11-S1-N
3	A	602	OXL	O3-C1-C2-O4
3	C	602	OXL	O3-C1-C2-O4
3	D	602	OXL	O3-C1-C2-O4
3	E	602	OXL	O3-C1-C2-O4
3	F	602	OXL	O3-C1-C2-O4
3	G	602	OXL	O3-C1-C2-O4
3	B	602	OXL	O3-C1-C2-O4

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Mol	Chain	Res	Type	Atoms
3	H	602	OXL	O3-C1-C2-O4
2	H	601	FBP	O5-C5-C6-O6
6	G	605	OAT	C9-C10-N-C17
3	A	602	OXL	O1-C1-C2-O2
3	B	602	OXL	O1-C1-C2-O2
3	D	602	OXL	O1-C1-C2-O2
3	E	602	OXL	O1-C1-C2-O2
6	G	605	OAT	C10-N-S1-O5
3	C	602	OXL	O1-C1-C2-O2
3	F	602	OXL	O1-C1-C2-O2
3	G	602	OXL	O1-C1-C2-O2
3	H	602	OXL	O1-C1-C2-O2
6	H	605	OAT	C10-N-S1-O4
2	C	601	FBP	C6-O6-P2-O4P
2	A	601	FBP	O5-C5-C6-O6
2	E	601	FBP	O5-C5-C6-O6
2	F	601	FBP	O5-C5-C6-O6
2	G	601	FBP	O5-C5-C6-O6
6	G	605	OAT	C10-N-S1-C11
3	D	602	OXL	O3-C1-C2-O2
3	E	602	OXL	O1-C1-C2-O4
3	E	602	OXL	O3-C1-C2-O2
3	A	602	OXL	O1-C1-C2-O4
3	A	602	OXL	O3-C1-C2-O2
3	B	602	OXL	O1-C1-C2-O4
3	B	602	OXL	O3-C1-C2-O2
3	C	602	OXL	O1-C1-C2-O4
3	C	602	OXL	O3-C1-C2-O2
3	D	602	OXL	O1-C1-C2-O4
3	F	602	OXL	O3-C1-C2-O2
3	G	602	OXL	O1-C1-C2-O4
3	G	602	OXL	O3-C1-C2-O2
3	H	602	OXL	O1-C1-C2-O4
2	B	601	FBP	O5-C5-C6-O6
3	F	602	OXL	O1-C1-C2-O4
3	H	602	OXL	O3-C1-C2-O2
2	A	601	FBP	C6-O6-P2-O5P
2	H	601	FBP	C6-O6-P2-O4P
6	B	605	OAT	C18-C17-N-S1

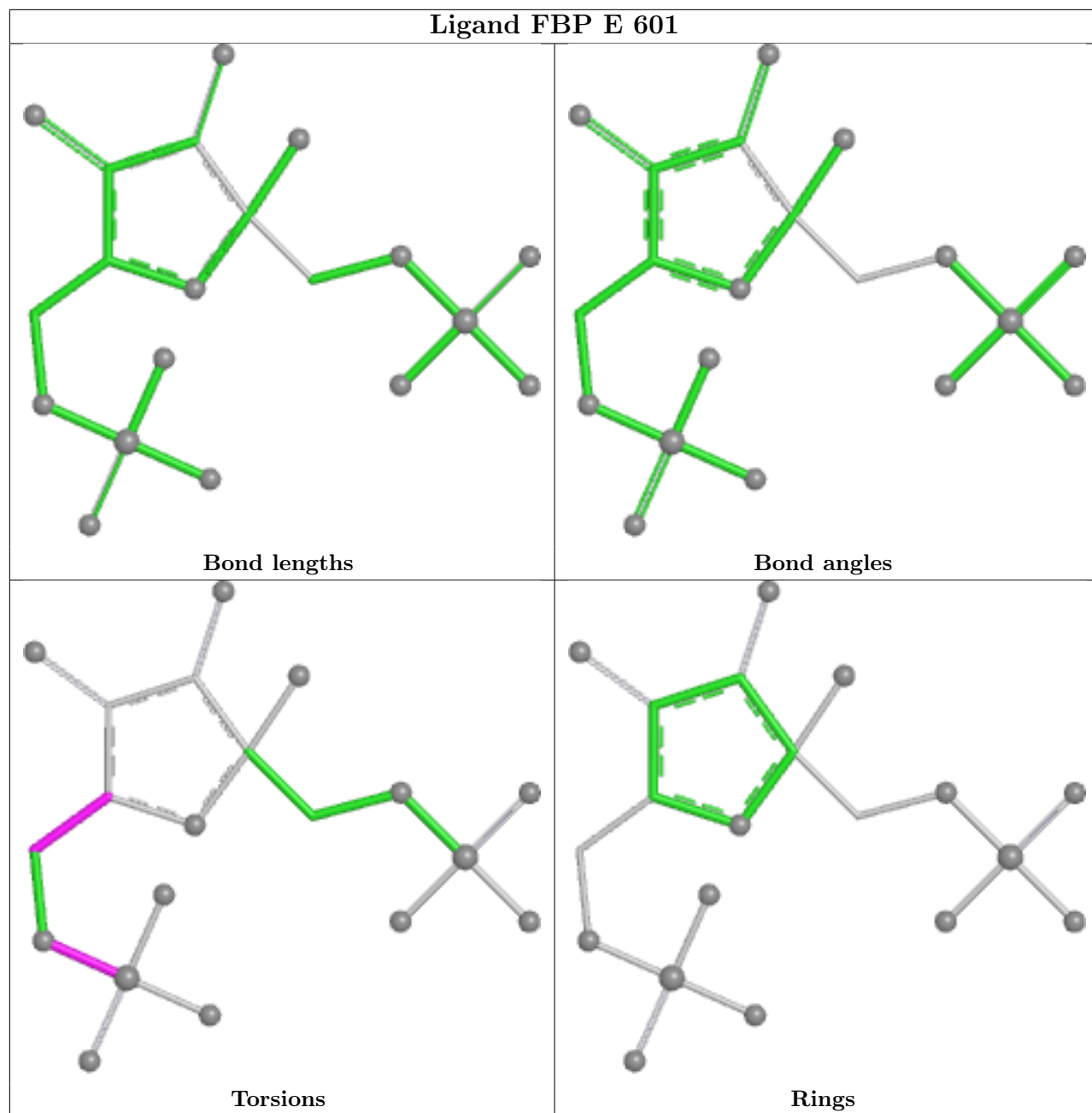
There are no ring outliers.

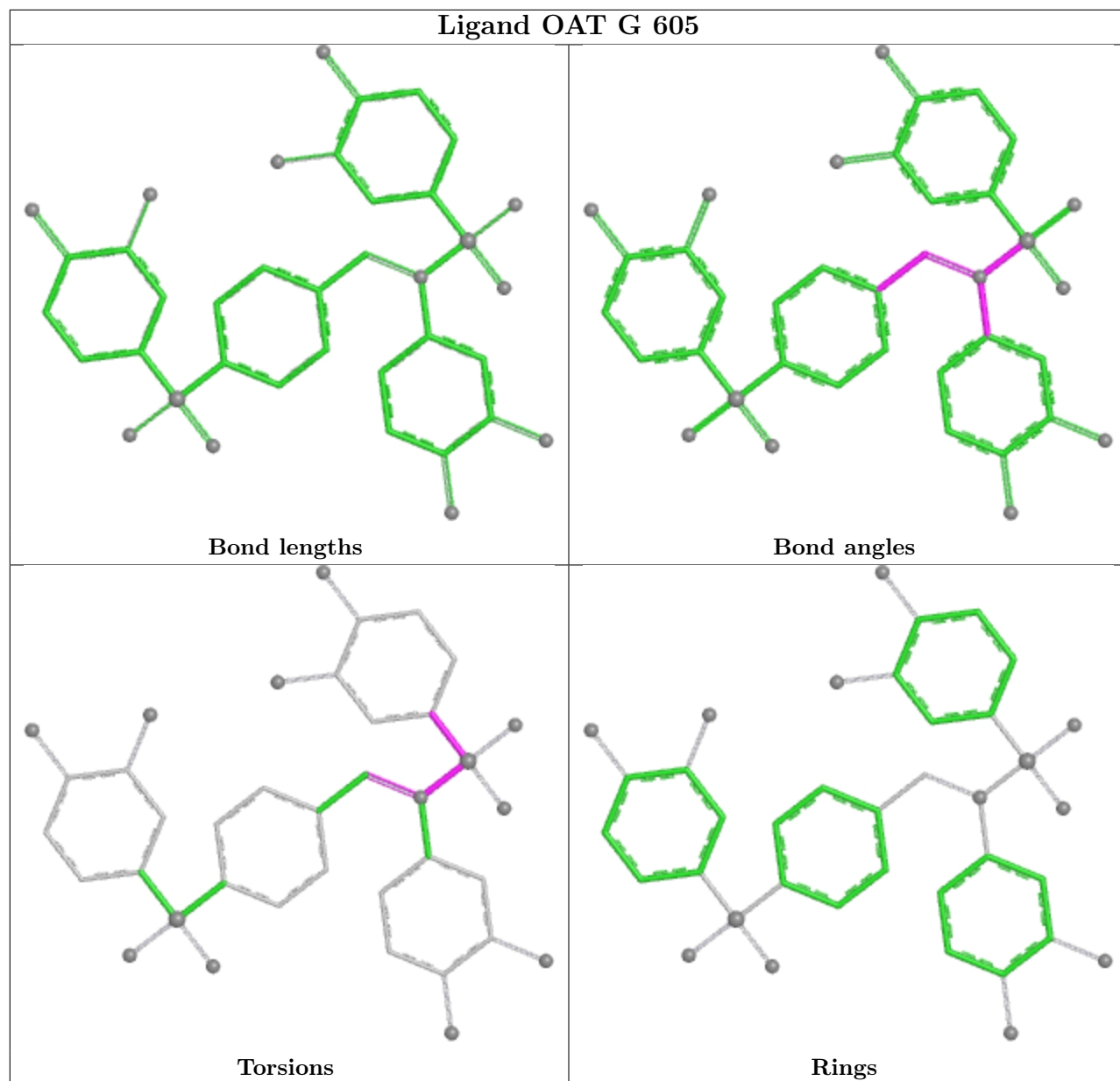
3 monomers are involved in 8 short contacts:

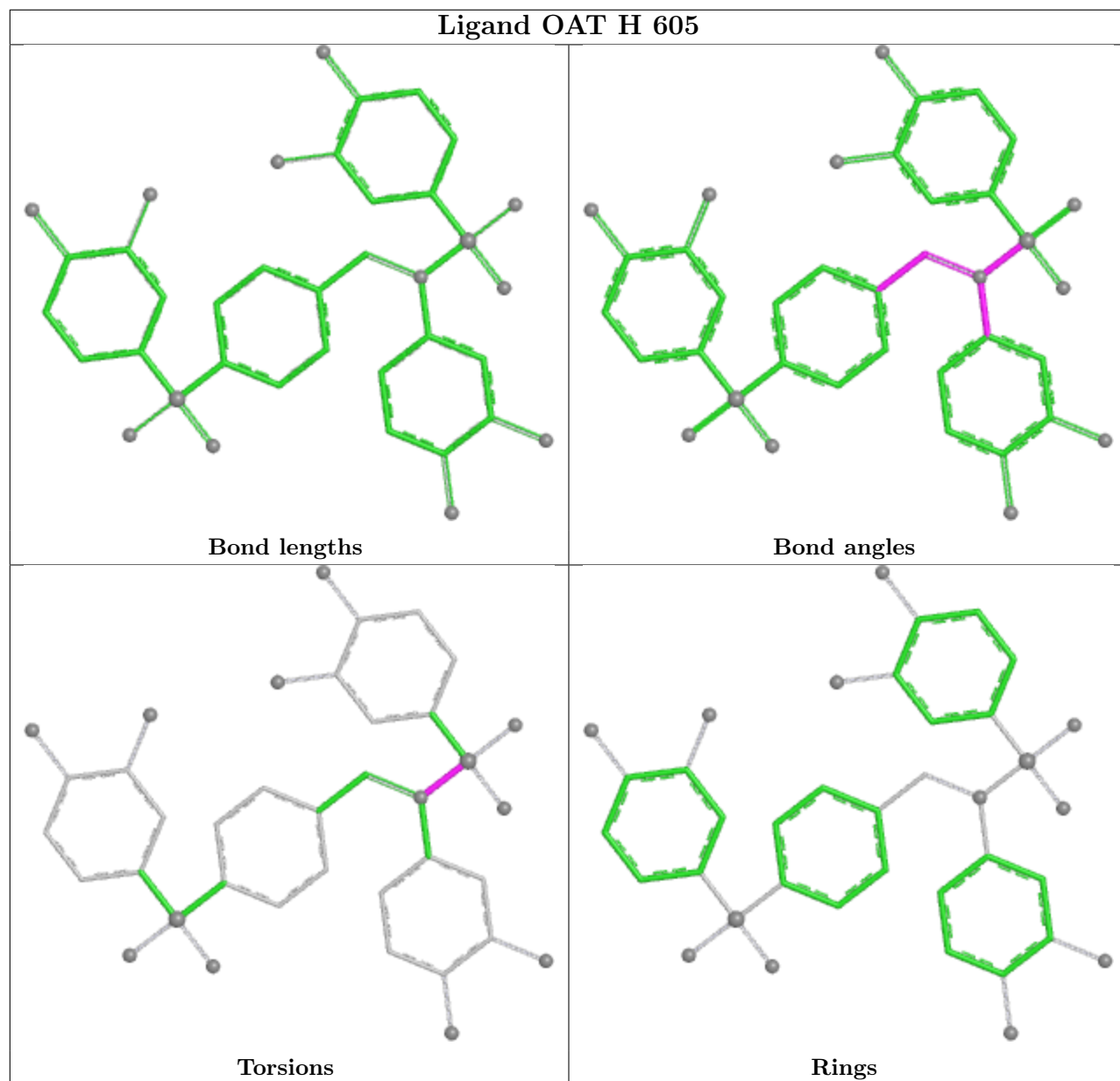
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	601	FBP	1	0
6	G	605	OAT	3	0
6	B	605	OAT	4	0

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.

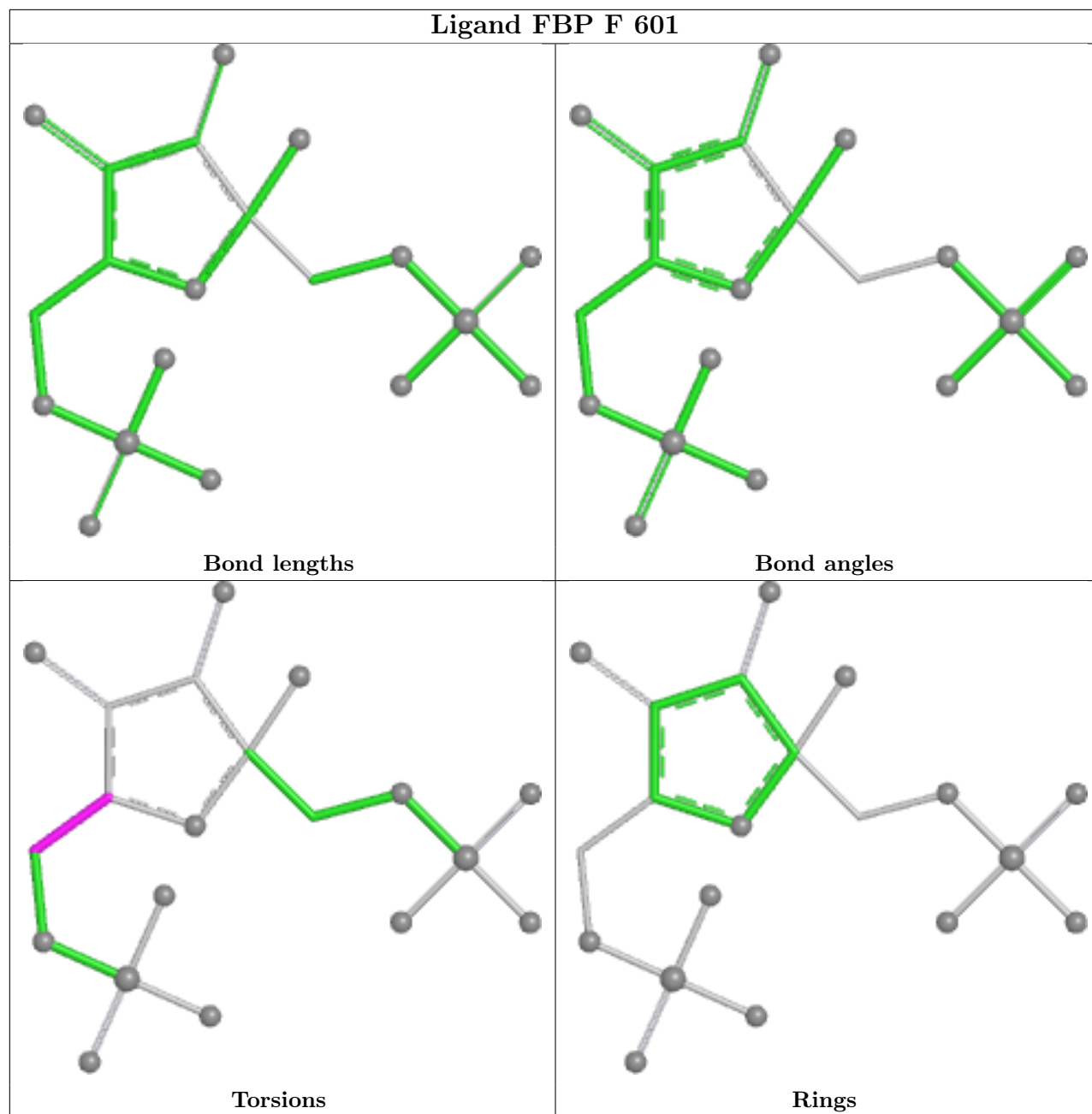
Ligand FBP E 601

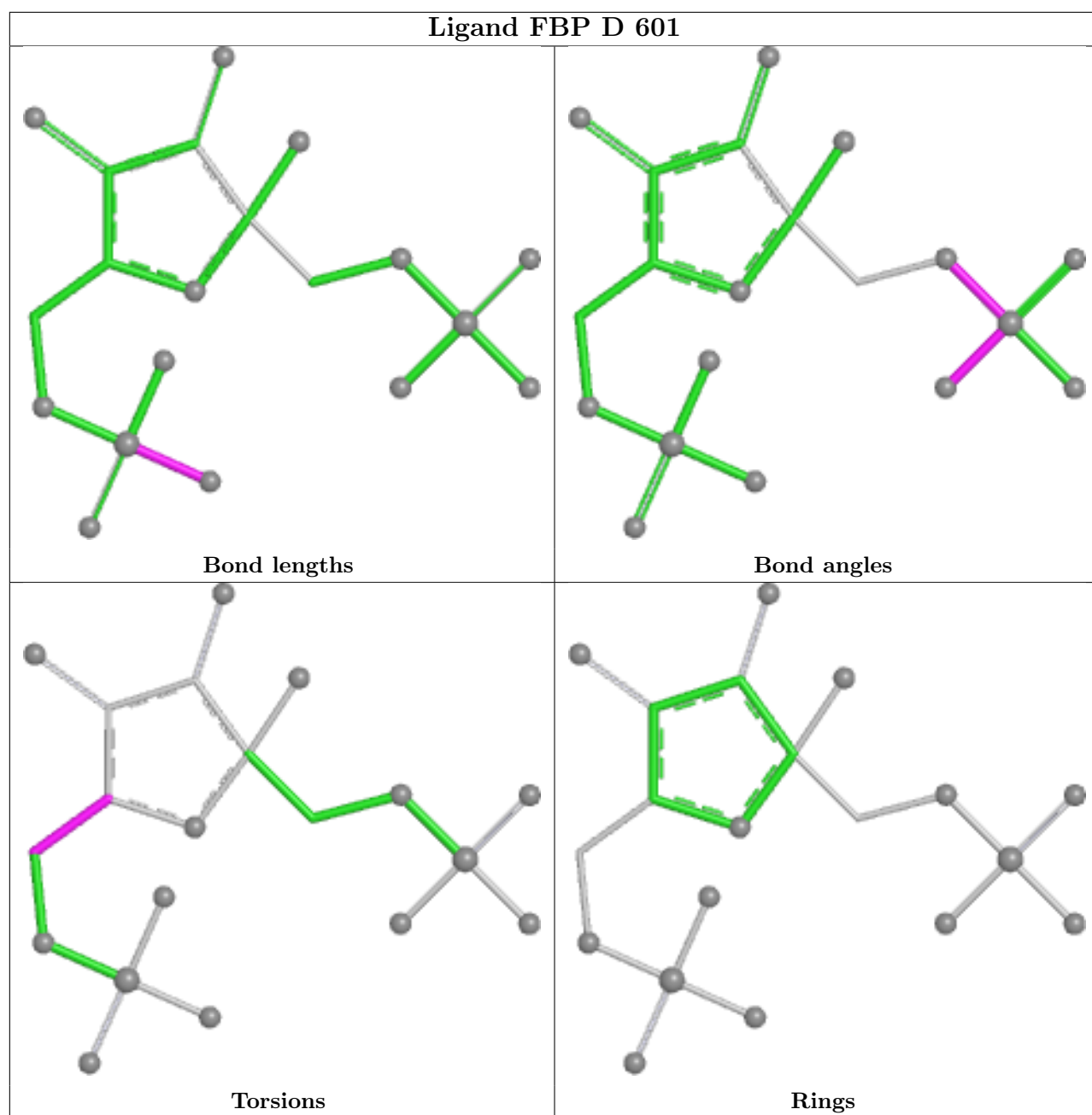


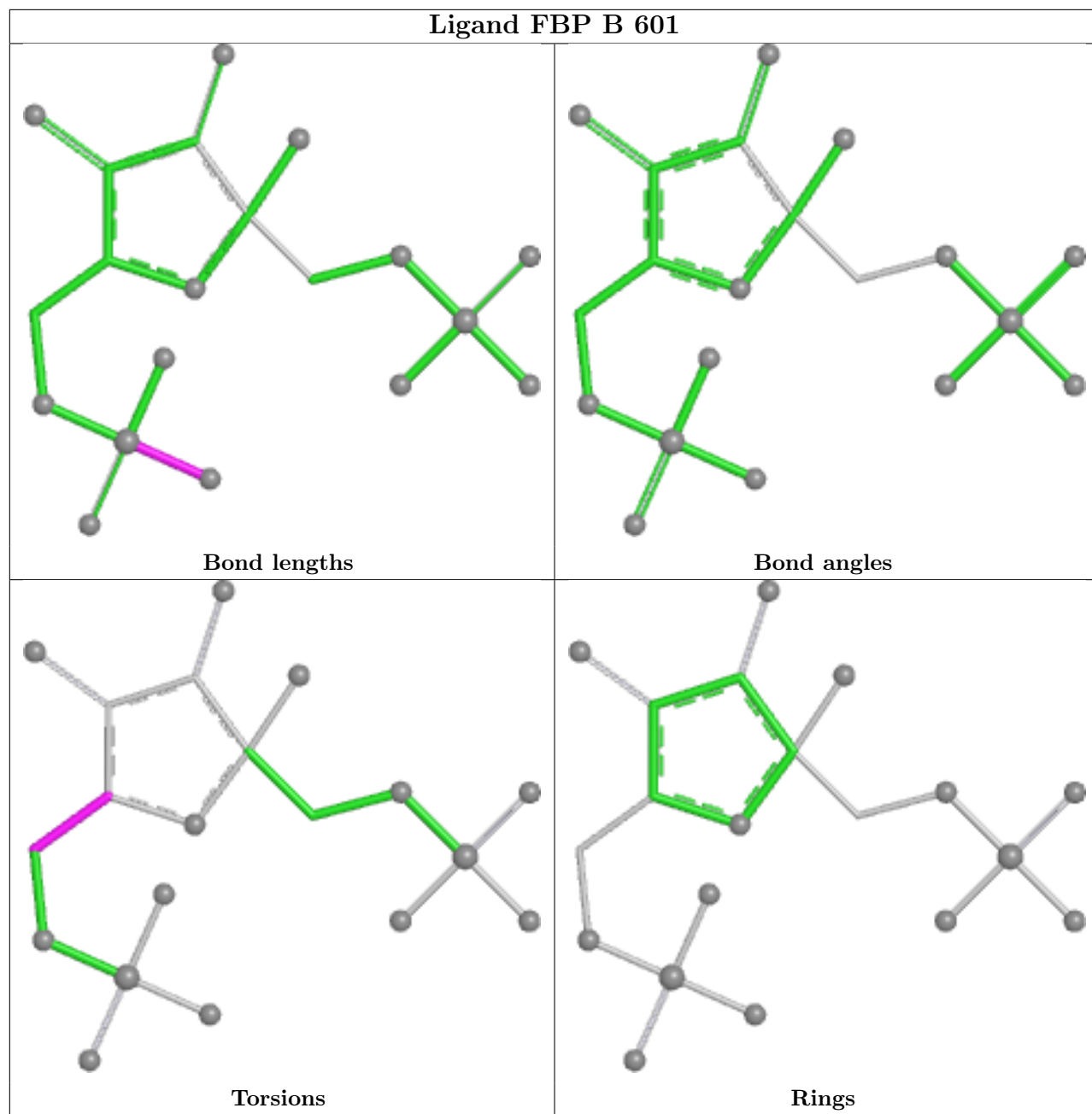


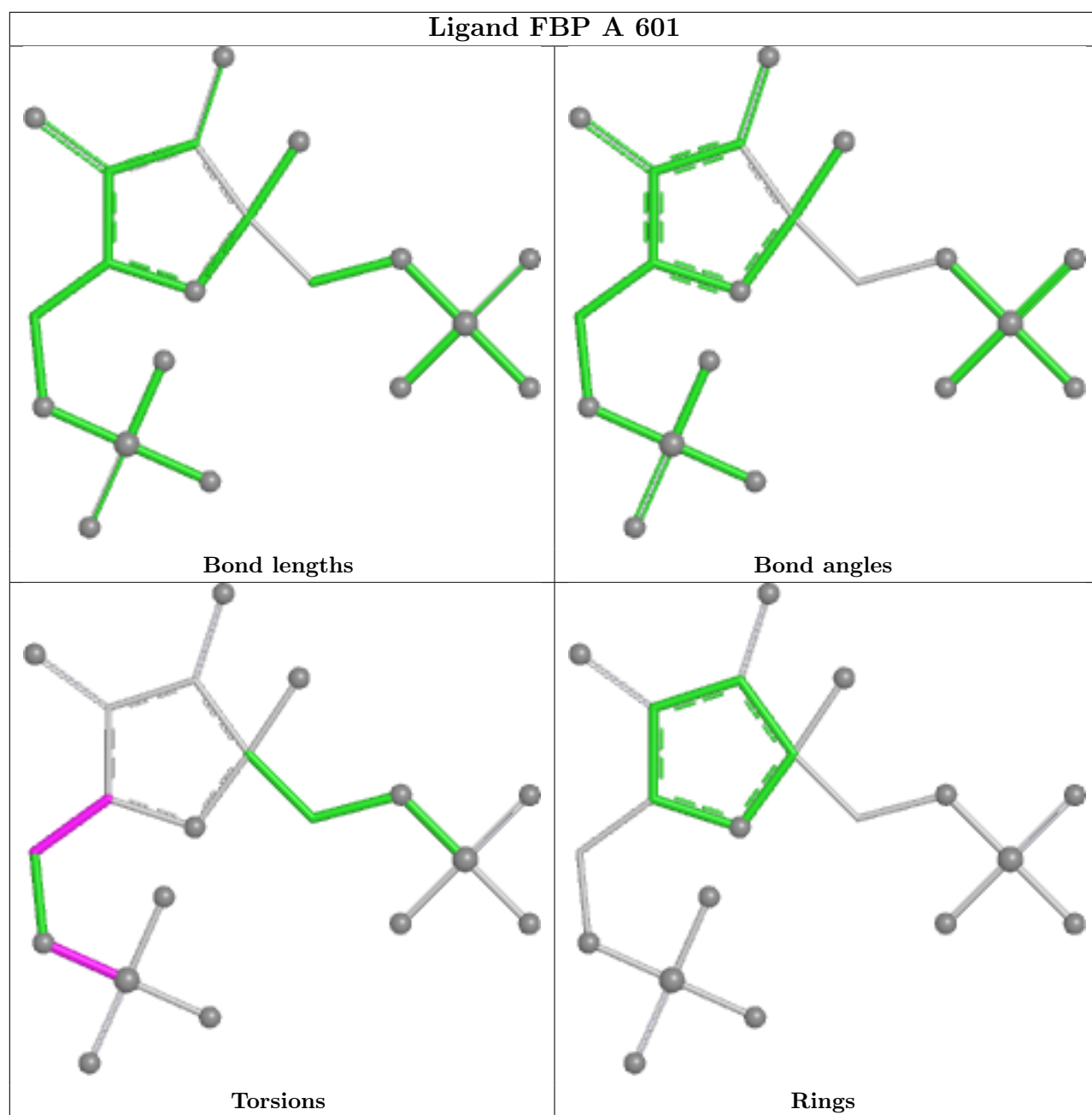


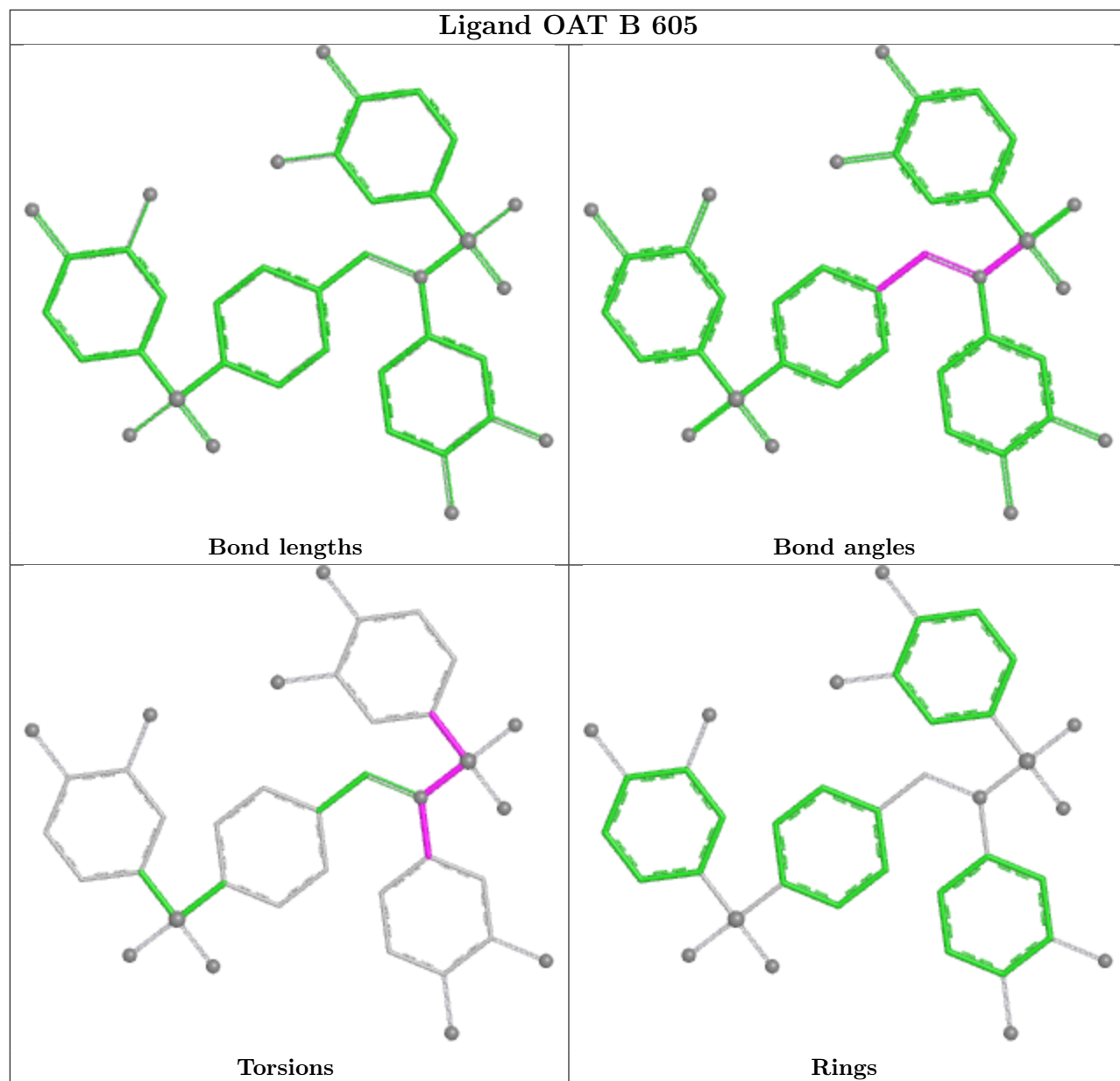
Ligand FBP F 601

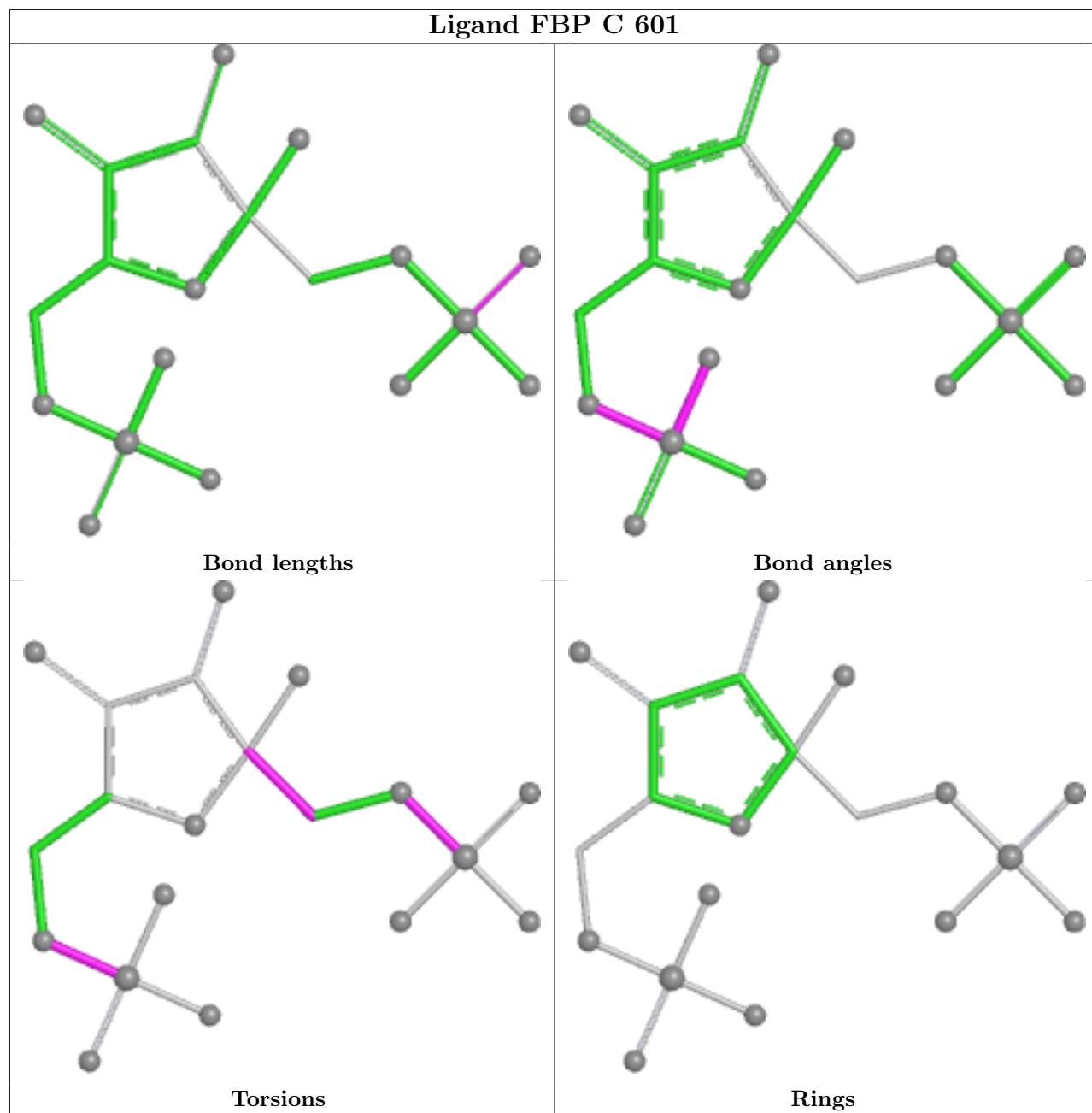


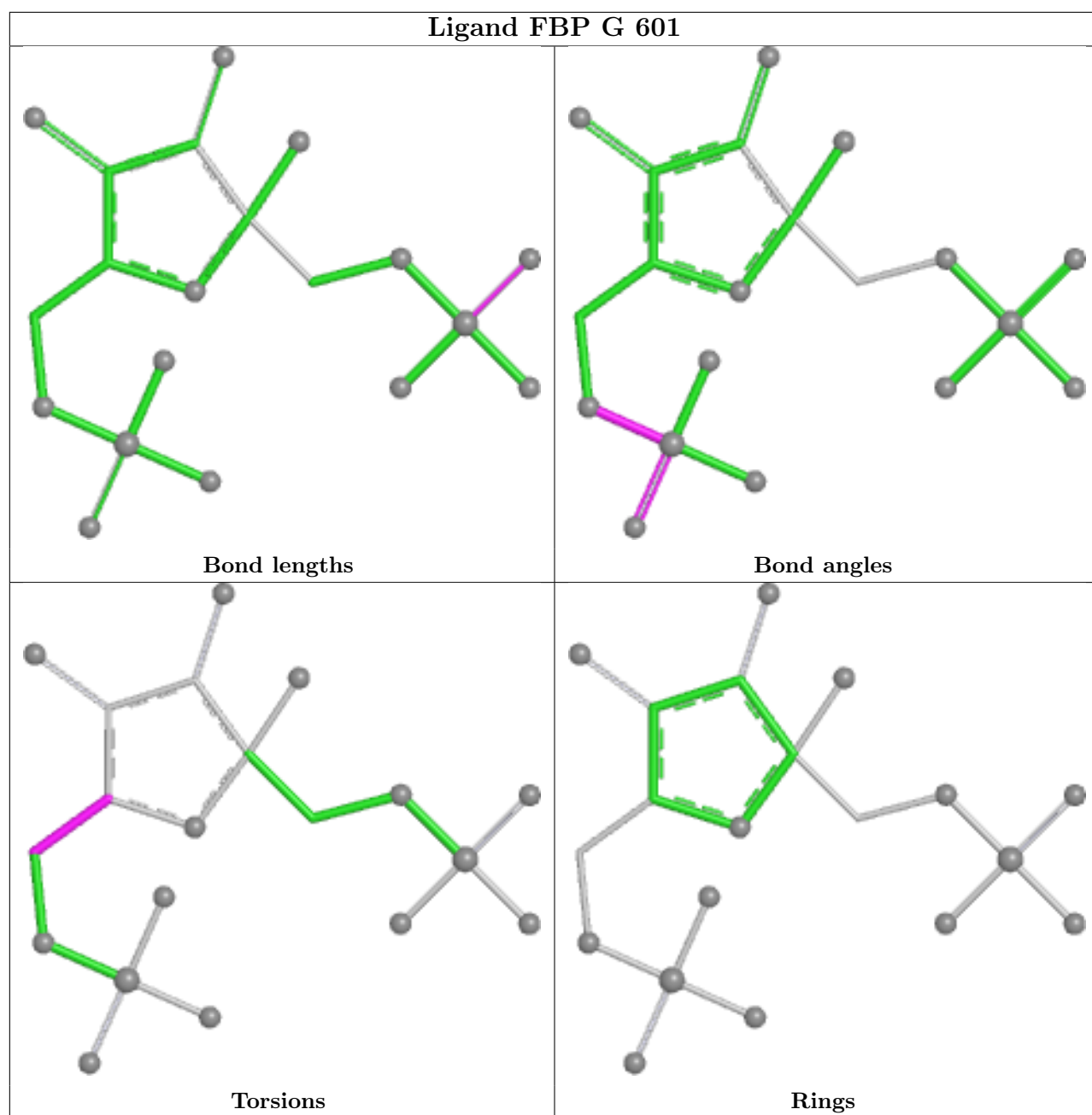


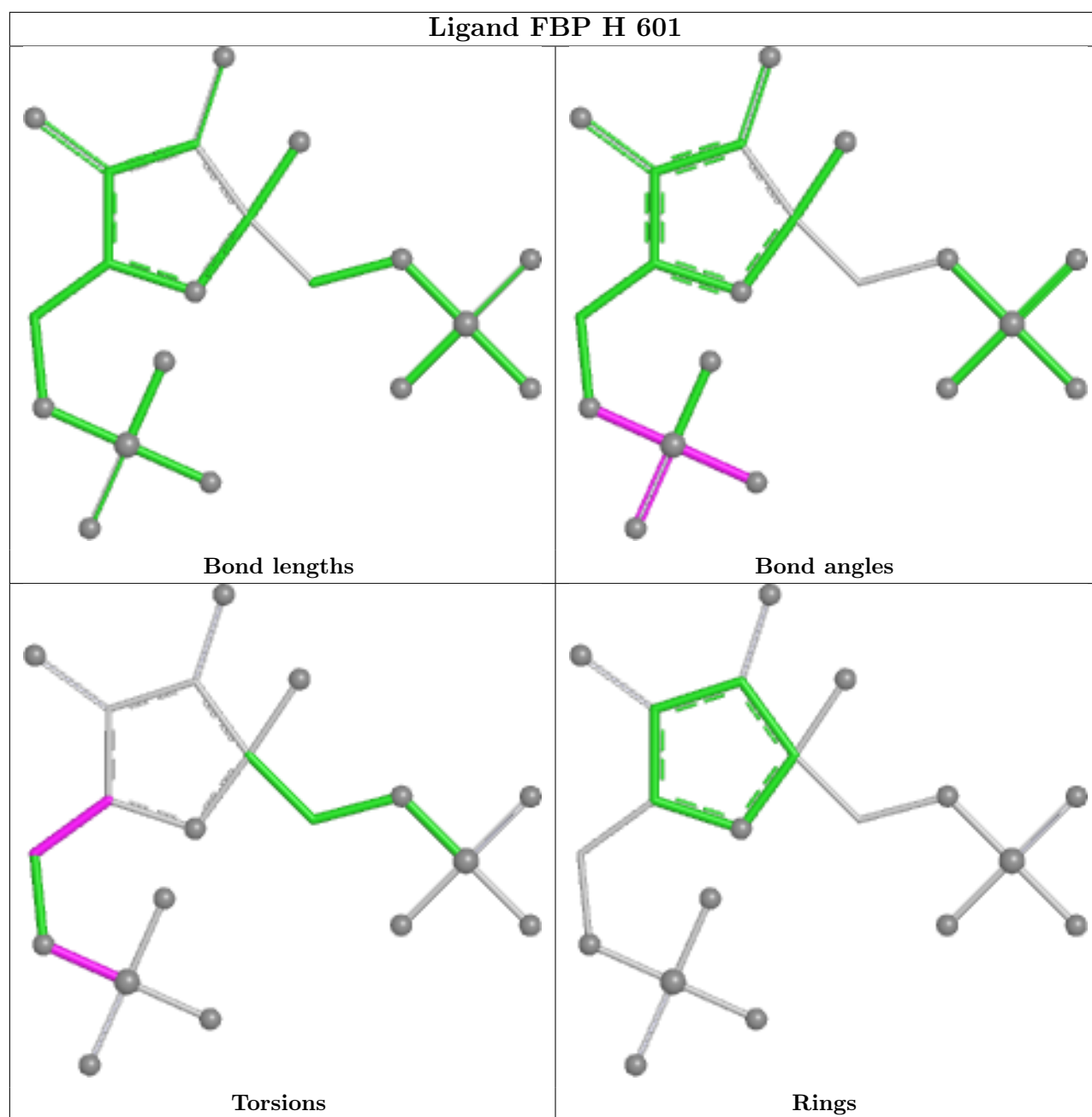












5.7 Other polymers [i](#)

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	422/447 (94%)	1.26	81 (19%) 3 2	71, 141, 183, 196	6 (1%)
1	B	436/447 (97%)	0.69	34 (7%) 19 15	66, 107, 146, 160	4 (0%)
1	C	425/447 (95%)	0.69	28 (6%) 24 19	40, 97, 143, 171	4 (0%)
1	D	425/447 (95%)	0.12	9 (2%) 63 56	29, 64, 104, 157	6 (1%)
1	E	419/447 (93%)	1.22	85 (20%) 3 2	62, 133, 172, 183	5 (1%)
1	F	432/447 (96%)	0.77	43 (9%) 12 10	60, 103, 141, 156	7 (1%)
1	G	421/447 (94%)	0.49	17 (4%) 42 36	38, 86, 121, 146	7 (1%)
1	H	425/447 (95%)	0.09	9 (2%) 63 56	28, 63, 101, 155	4 (0%)
All	All	3405/3576 (95%)	0.66	306 (8%) 15 12	28, 100, 163, 196	43 (1%)

All (306) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	442	VAL	7.6
1	A	463	ILE	6.3
1	B	16	LEU	5.1
1	E	463	ILE	5.1
1	E	447	GLY	5.0
1	G	116	SER	5.0
1	A	482	PHE	4.9
1	B	13	VAL	4.8
1	E	319	PHE	4.7
1	A	465	VAL	4.7
1	A	241	LEU	4.6
1	A	453	LEU	4.6
1	E	120	VAL	4.6
1	E	397	ALA	4.5
1	G	115	LEU	4.5
1	A	63	ILE	4.4

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Mol	Chain	Res	Type	RSRZ
1	G	114	PRO	4.3
1	G	363	ALA	4.2
1	G	322	GLN	4.1
1	F	115	LEU	4.1
1	E	238	VAL	4.1
1	A	238	VAL	4.0
1	F	33	ALA	3.9
1	A	484	LEU	3.9
1	A	543	SER	3.9
1	B	12	ASP	3.9
1	A	313	ILE	3.7
1	E	249	VAL	3.7
1	A	232	GLY	3.7
1	B	438	ALA	3.7
1	E	244	GLY	3.7
1	B	517	VAL	3.7
1	A	492	ALA	3.6
1	A	489	PRO	3.6
1	E	129	PRO	3.6
1	B	540[A]	LEU	3.6
1	E	502	VAL	3.6
1	C	22	THR	3.6
1	A	319	PHE	3.5
1	A	450	ALA	3.5
1	E	505	GLY	3.5
1	F	116	SER	3.5
1	F	540[A]	LEU	3.5
1	E	103	VAL	3.4
1	F	514	PHE	3.4
1	A	361	ALA	3.4
1	C	20	LEU	3.4
1	H	25	PHE	3.4
1	F	459	ARG	3.3
1	A	357	THR	3.3
1	A	350	PRO	3.3
1	C	114	PRO	3.3
1	F	61	ALA	3.3
1	A	443	LEU	3.3
1	B	115	LEU	3.3
1	C	33	ALA	3.3
1	A	490	PRO	3.3
1	E	348	THR	3.3

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Mol	Chain	Res	Type	RSRZ
1	E	248	GLY	3.2
1	E	400	ALA	3.2
1	A	441	ILE	3.2
1	E	484	LEU	3.2
1	E	254	ALA	3.2
1	E	375	GLY	3.2
1	G	63	ILE	3.2
1	E	264	ALA	3.1
1	B	521	VAL	3.1
1	A	353	THR	3.1
1	E	304	VAL	3.1
1	E	115	LEU	3.1
1	B	11	ALA	3.1
1	E	260	ALA	3.1
1	F	113	SER	3.1
1	F	511	LEU	3.0
1	F	513	GLY	3.0
1	A	264	ALA	3.0
1	E	123	ALA	3.0
1	A	115	LEU	3.0
1	C	51	PRO	3.0
1	E	23	ALA	3.0
1	A	123	ALA	3.0
1	A	442	VAL	3.0
1	F	241	LEU	2.9
1	A	471	ALA	2.9
1	D	25	PHE	2.9
1	F	527	TRP	2.9
1	B	132	GLY	2.9
1	E	465	VAL	2.9
1	F	502	VAL	2.9
1	F	39	LEU	2.9
1	G	110	PHE	2.9
1	A	100	ILE	2.9
1	F	463	ILE	2.9
1	A	86	LEU	2.9
1	A	116	SER	2.9
1	A	423	VAL	2.9
1	A	114	PRO	2.8
1	E	122	ILE	2.8
1	B	33	ALA	2.8
1	F	521	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	65	PRO	2.8
1	E	251	ILE	2.8
1	A	84	ALA	2.8
1	E	119	PRO	2.8
1	F	490	PRO	2.8
1	C	304	VAL	2.8
1	D	33	ALA	2.8
1	E	114	PRO	2.7
1	F	485	LEU	2.7
1	A	269	ALA	2.7
1	A	397	ALA	2.7
1	E	257	VAL	2.7
1	F	524	VAL	2.7
1	E	266	VAL	2.7
1	E	495	ALA	2.7
1	A	25	PHE	2.7
1	B	24	PHE	2.7
1	B	490	PRO	2.7
1	C	371	ILE	2.7
1	E	443	LEU	2.7
1	E	263	VAL	2.7
1	E	482	PHE	2.7
1	B	25	PHE	2.7
1	E	498	VAL	2.7
1	B	119	PRO	2.7
1	C	63	ILE	2.6
1	D	331	LEU	2.6
1	G	64	GLY	2.6
1	A	318	VAL	2.6
1	C	384	VAL	2.6
1	G	254	ALA	2.6
1	B	114	PRO	2.6
1	G	118	ARG	2.6
1	A	132	GLY	2.6
1	A	539	VAL	2.6
1	A	251	ILE	2.6
1	A	506	ILE	2.6
1	E	536	ILE	2.6
1	E	232	GLY	2.6
1	A	81	MET	2.6
1	C	77	ILE	2.6
1	F	512	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	E	489	PRO	2.6
1	A	493	ILE	2.5
1	E	241	LEU	2.5
1	F	516	ARG	2.5
1	F	52	VAL	2.5
1	E	29	GLN	2.5
1	H	23	ALA	2.5
1	A	279	ILE	2.5
1	B	542	ILE	2.5
1	H	21	GLY	2.5
1	B	249	VAL	2.5
1	C	103	VAL	2.5
1	F	11	ALA	2.5
1	C	65	PRO	2.5
1	F	119	PRO	2.5
1	A	120	VAL	2.5
1	E	245	VAL	2.5
1	F	103	VAL	2.5
1	A	536	ILE	2.4
1	H	331	LEU	2.4
1	A	352	PRO	2.4
1	A	380	GLY	2.4
1	B	502	VAL	2.4
1	A	97	ALA	2.4
1	E	280	ILE	2.4
1	H	30	LEU	2.4
1	G	374[A]	SER	2.4
1	E	523	VAL	2.4
1	A	373	LEU	2.4
1	A	534	THR	2.4
1	E	233	LEU	2.4
1	E	525	THR	2.4
1	E	381	ASN	2.4
1	E	116	SER	2.4
1	C	132	GLY	2.4
1	B	439	ALA	2.4
1	D	24	PHE	2.4
1	E	104	ARG	2.4
1	E	269	ALA	2.4
1	A	314	PRO	2.4
1	C	129	PRO	2.4
1	E	289	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	267[A]	ARG	2.4
1	A	253	PHE	2.4
1	E	25	PHE	2.4
1	H	275[A]	HIS	2.3
1	E	65	PRO	2.3
1	H	242	ARG	2.3
1	D	23	ALA	2.3
1	E	379	LYS	2.3
1	E	490	PRO	2.3
1	C	322	GLN	2.3
1	E	333	GLY	2.3
1	F	28	GLN	2.3
1	G	232	GLY	2.3
1	A	464	ALA	2.3
1	C	73	LEU	2.3
1	C	59	ILE	2.3
1	E	493	ILE	2.3
1	F	12	ASP	2.3
1	G	129	PRO	2.3
1	F	13	VAL	2.3
1	B	493	ILE	2.3
1	D	361	ALA	2.3
1	F	506	ILE	2.3
1	B	505	GLY	2.3
1	C	380	GLY	2.3
1	A	39	LEU	2.3
1	A	502	VAL	2.3
1	E	441	ILE	2.2
1	E	522	ILE	2.2
1	E	247	HIS	2.2
1	F	377	THR	2.2
1	B	267[A]	ARG	2.2
1	E	272	PRO	2.2
1	A	249	VAL	2.2
1	A	343	LEU	2.2
1	B	463	ILE	2.2
1	A	119	PRO	2.2
1	F	114	PRO	2.2
1	E	541[A]	SER	2.2
1	B	511	LEU	2.2
1	D	540	LEU	2.2
1	A	349	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	475	VAL	2.2
1	E	481	VAL	2.2
1	A	393	ILE	2.2
1	C	389	MET	2.2
1	E	464	ALA	2.2
1	E	472	ALA	2.2
1	E	108	GLU	2.2
1	A	383	PRO	2.2
1	F	489	PRO	2.2
1	E	100	ILE	2.2
1	B	509	GLY	2.2
1	B	515	LEU	2.2
1	E	124	LEU	2.2
1	C	366	ASP	2.2
1	A	523	VAL	2.2
1	B	414	ALA	2.1
1	H	33	ALA	2.1
1	B	435	CYS	2.1
1	E	270	LEU	2.1
1	F	453	LEU	2.1
1	A	447	GLY	2.1
1	F	505	GLY	2.1
1	A	449	SER	2.1
1	C	107	VAL	2.1
1	E	58	SER	2.1
1	F	441	ILE	2.1
1	G	543	SER	2.1
1	E	265	ALA	2.1
1	E	386	ALA	2.1
1	A	527	TRP	2.1
1	A	446	THR	2.1
1	B	526	GLY	2.1
1	E	380	GLY	2.1
1	D	263	VAL	2.1
1	E	440	ILE	2.1
1	E	499	ASP	2.1
1	B	514	PHE	2.1
1	G	40[A]	GLU	2.1
1	C	350	PRO	2.1
1	F	483	PRO	2.1
1	C	117	TYR	2.1
1	A	83	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	103	VAL	2.1
1	A	252	VAL	2.1
1	A	524	VAL	2.1
1	C	401	VAL	2.1
1	E	539	VAL	2.1
1	F	24	PHE	2.1
1	B	14	ALA	2.1
1	C	363	ALA	2.1
1	E	389	MET	2.1
1	A	280	ILE	2.1
1	E	60	ILE	2.1
1	B	524	VAL	2.1
1	F	539	VAL	2.1
1	F	25	PHE	2.1
1	G	25	PHE	2.1
1	D	538	ARG	2.0
1	A	66	ALA	2.0
1	C	67	SER	2.0
1	E	471	ALA	2.0
1	A	270	LEU	2.0
1	A	410	LEU	2.0
1	A	477	LEU	2.0
1	A	440	ILE	2.0
1	C	232	GLY	2.0
1	C	393	ILE	2.0
1	E	302	ILE	2.0
1	E	524	VAL	2.0
1	F	475	VAL	2.0
1	F	517	VAL	2.0
1	A	110	PHE	2.0
1	A	488[A]	GLU	2.0
1	E	110	PHE	2.0
1	B	431	ALA	2.0
1	A	58	SER	2.0
1	G	113	SER	2.0
1	H	543	SER	2.0
1	F	515	LEU	2.0
1	E	542	ILE	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.

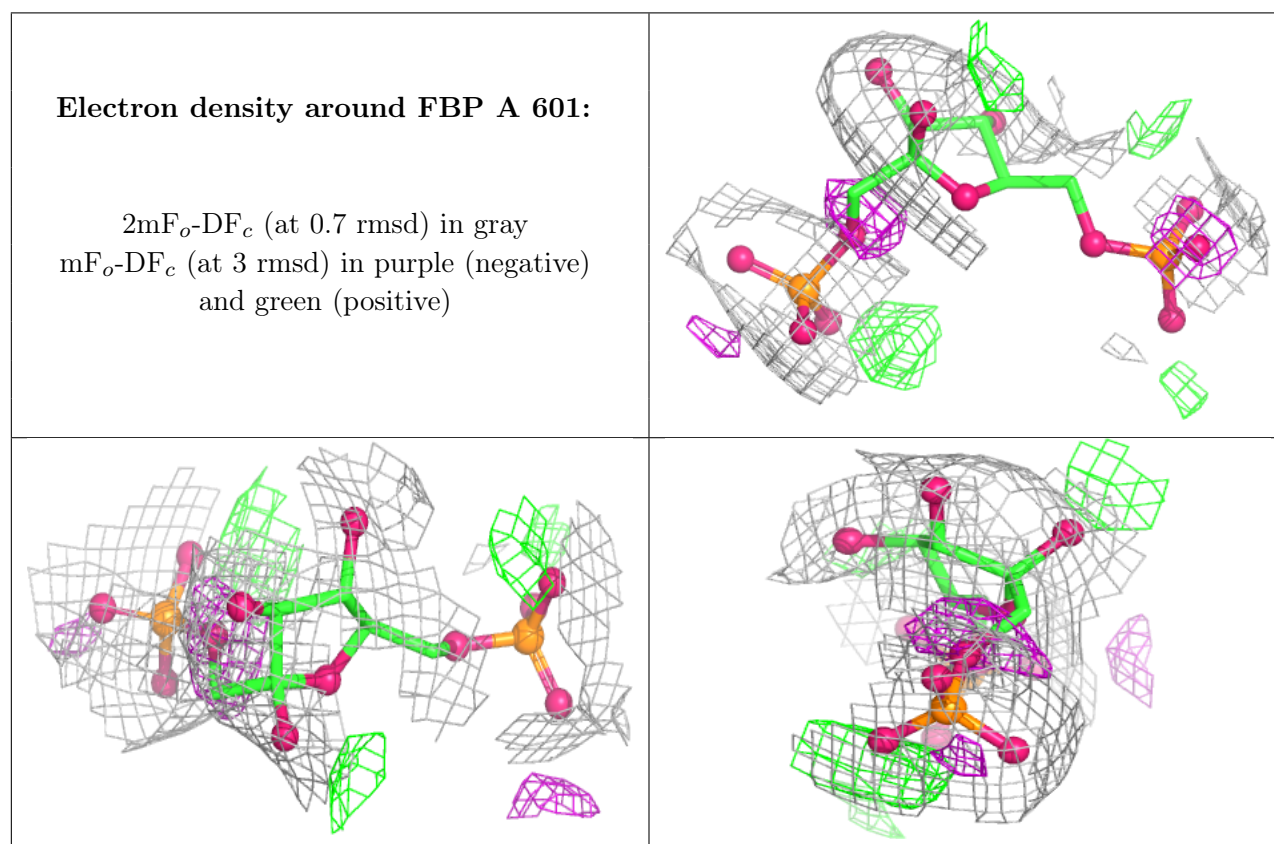
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	K	E	604	1/1	0.20	0.13	228,228,228,228	0
5	K	A	604	1/1	0.63	0.10	201,201,201,201	0
2	FBP	A	601	20/20	0.71	0.14	150,151,151,152	0
3	OXL	A	602	6/6	0.74	0.10	178,178,178,178	0
3	OXL	E	602	6/6	0.77	0.12	167,167,167,167	0
6	OAT	G	605	38/38	0.78	0.17	156,159,164,164	21
2	FBP	E	601	20/20	0.79	0.11	139,140,142,142	0
5	K	C	604	1/1	0.81	0.09	125,125,125,125	0
5	K	F	604	1/1	0.82	0.09	142,142,142,142	0
2	FBP	F	601	20/20	0.82	0.10	127,128,130,131	0
5	K	B	604	1/1	0.83	0.08	138,138,138,138	0
6	OAT	H	605	38/38	0.83	0.19	118,123,125,126	21
4	MG	A	603	1/1	0.85	0.07	141,141,141,141	0
6	OAT	B	605	38/38	0.85	0.16	120,126,130,130	21
3	OXL	C	602	6/6	0.86	0.13	156,156,156,156	0
3	OXL	H	602	6/6	0.89	0.16	126,127,127,127	0
2	FBP	B	601	20/20	0.89	0.09	120,121,125,125	0
3	OXL	F	602	6/6	0.90	0.12	172,172,172,172	0
3	OXL	G	602	6/6	0.91	0.11	123,123,123,124	0
3	OXL	D	602	6/6	0.93	0.15	138,138,138,138	0
4	MG	C	603	1/1	0.93	0.12	88,88,88,88	0
4	MG	E	603	1/1	0.93	0.04	111,111,111,111	0
4	MG	G	603	1/1	0.93	0.07	67,67,67,67	0
4	MG	B	603	1/1	0.95	0.05	79,79,79,79	0
3	OXL	B	602	6/6	0.95	0.08	150,150,151,151	0
5	K	G	604	1/1	0.95	0.04	104,104,104,104	0
5	K	H	604	1/1	0.95	0.07	98,98,98,98	0
2	FBP	G	601	20/20	0.95	0.07	66,69,72,73	0
4	MG	F	603	1/1	0.95	0.07	71,71,71,71	0
5	K	D	604	1/1	0.95	0.06	100,100,100,100	0
2	FBP	C	601	20/20	0.96	0.07	70,71,71,71	0
4	MG	H	603	1/1	0.97	0.07	64,64,64,64	0

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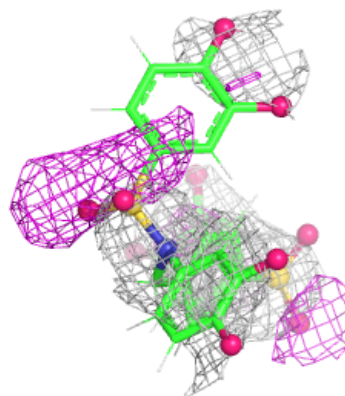
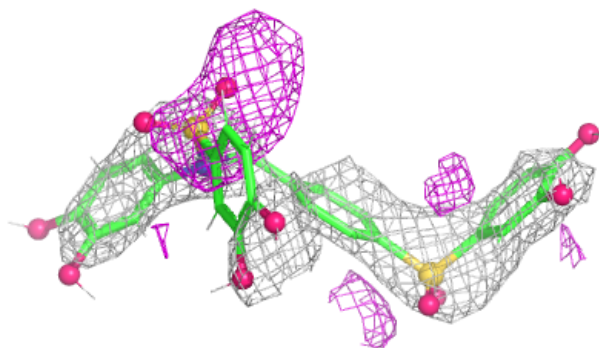
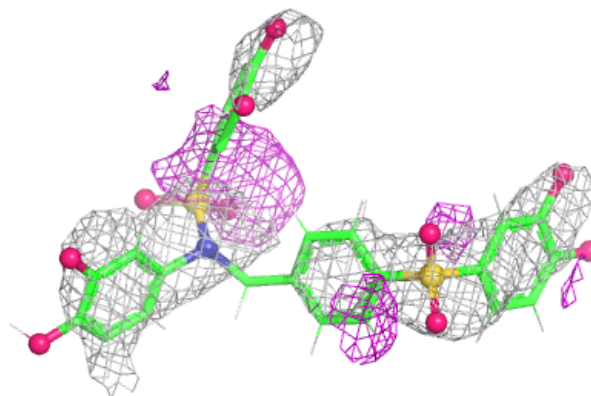
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FBP	H	601	20/20	0.97	0.06	43,52,53,54	0
2	FBP	D	601	20/20	0.98	0.06	49,51,53,53	0
4	MG	D	603	1/1	0.99	0.06	60,60,60,60	0

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.

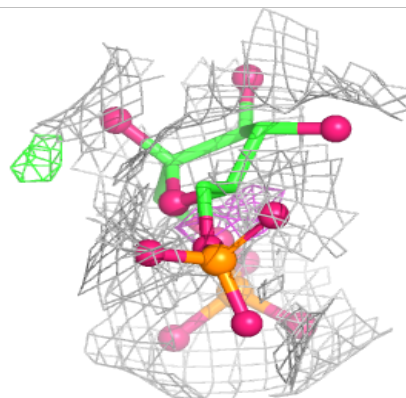
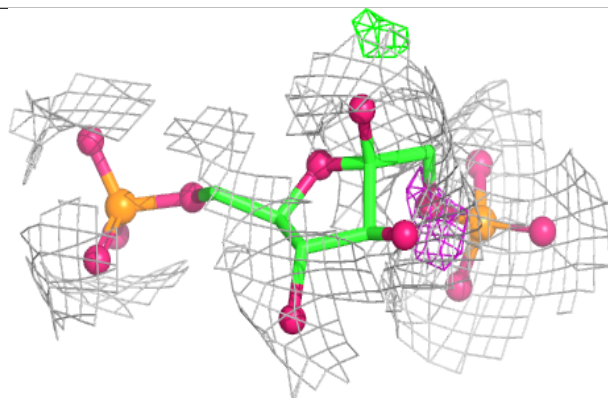
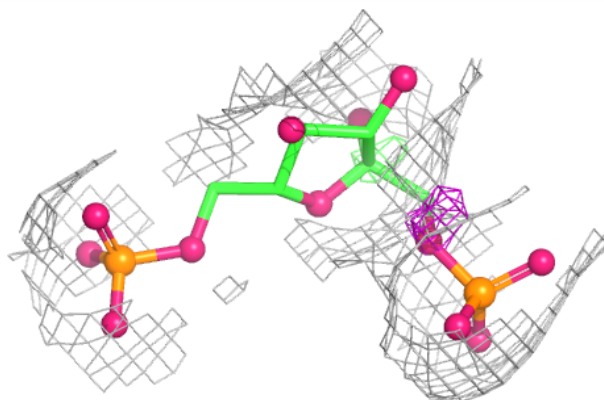


Electron density around OAT G 605:

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

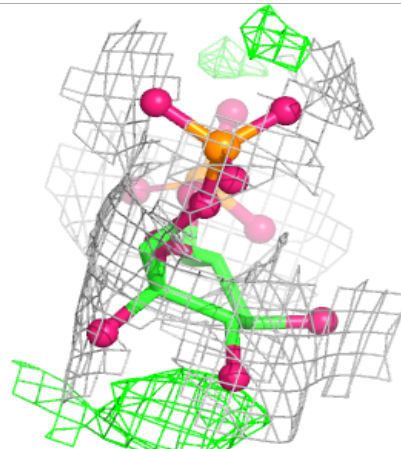
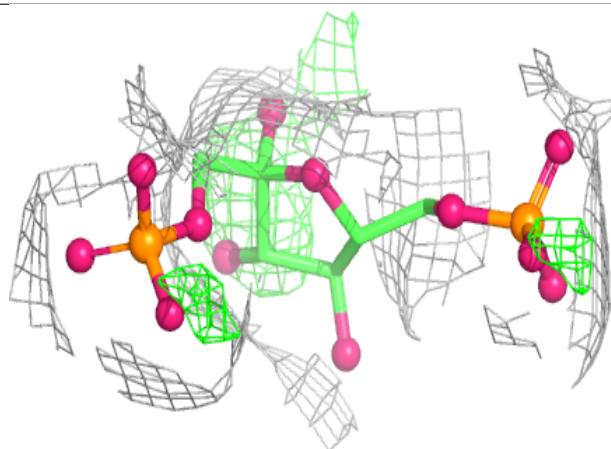
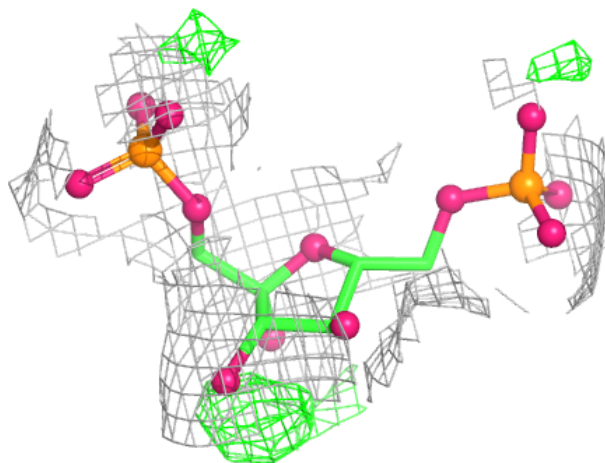
**Electron density around FBP E 601:**

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



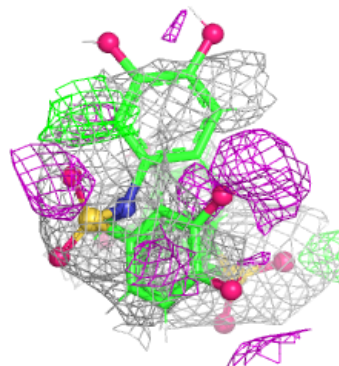
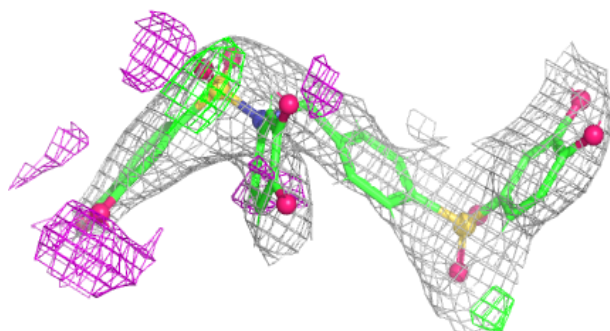
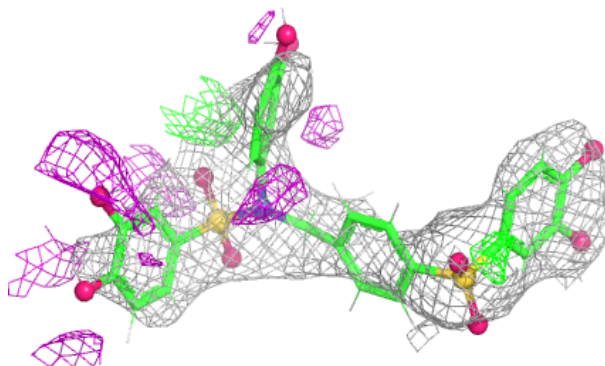
Electron density around FBP F 601:

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

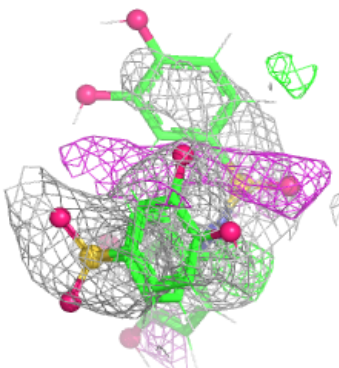
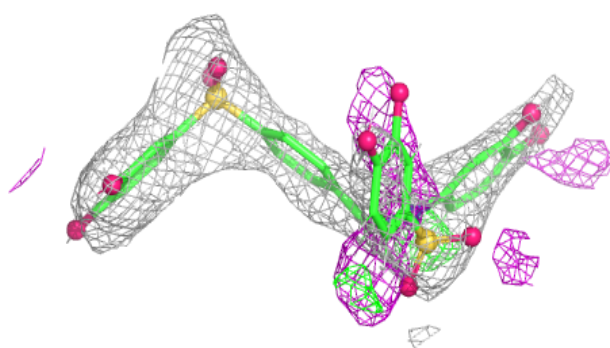
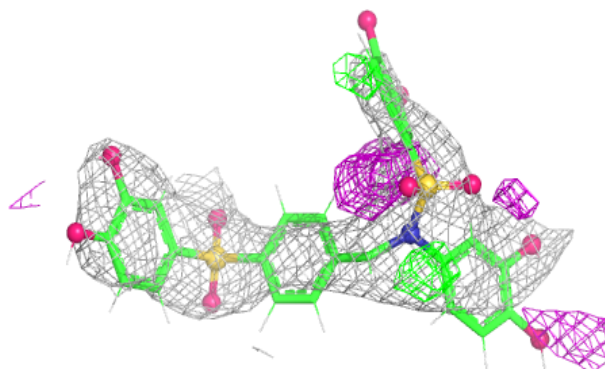


Electron density around OAT H 605:

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

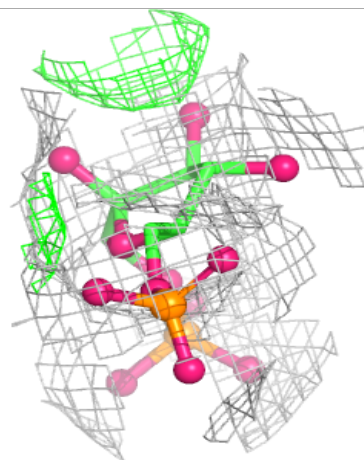
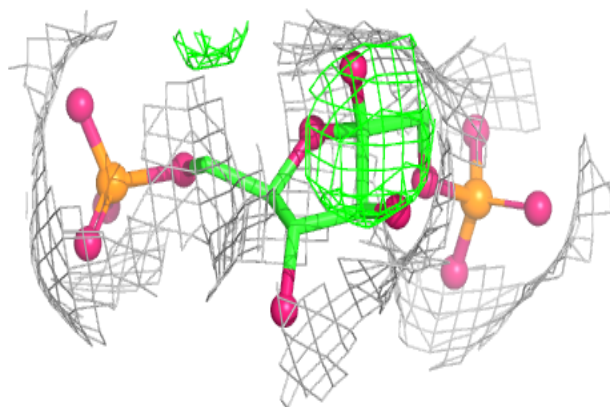
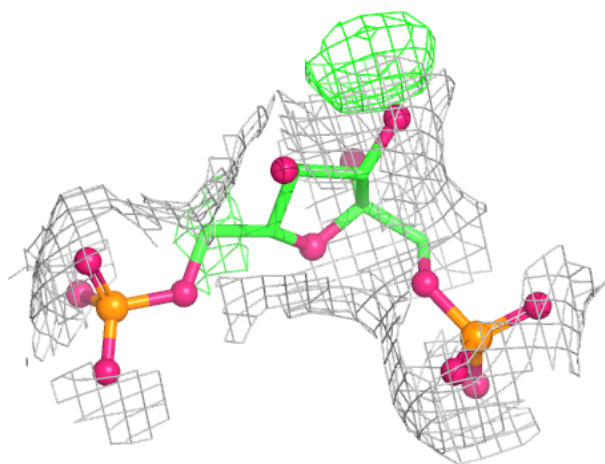
**Electron density around OAT B 605:**

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



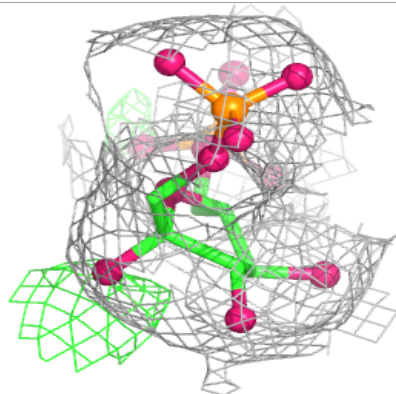
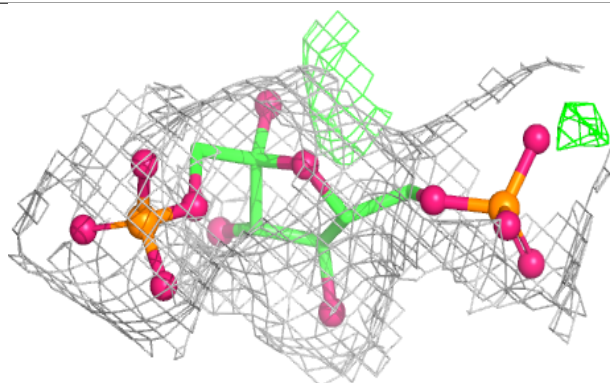
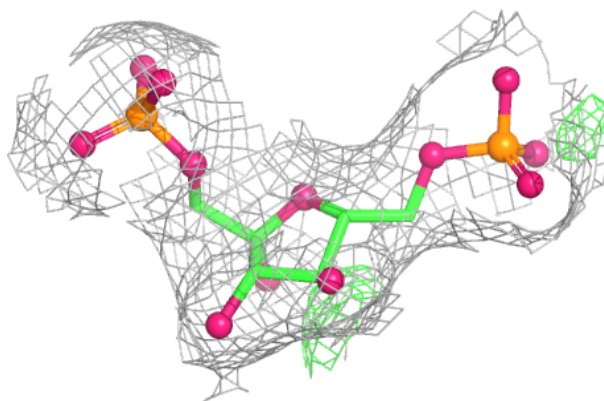
Electron density around FBP B 601:

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

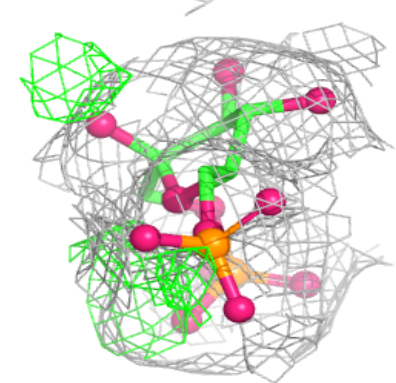
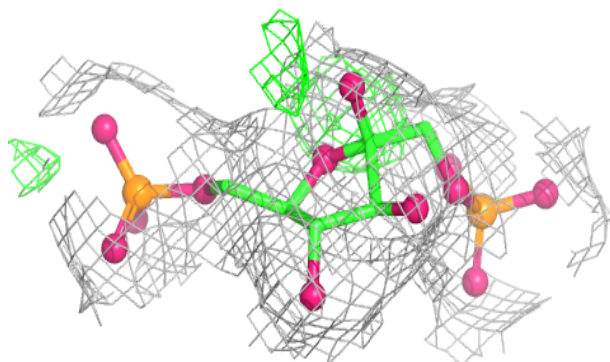
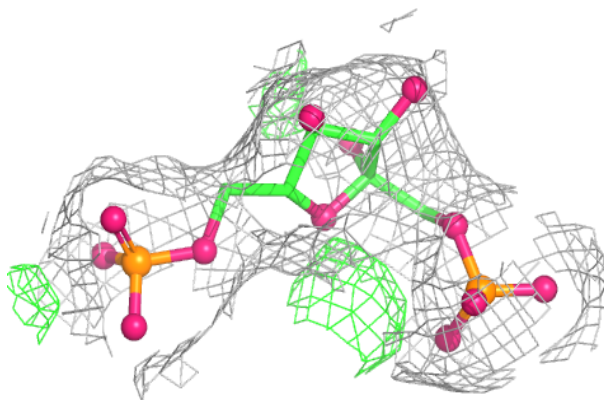


Electron density around FBP G 601:

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

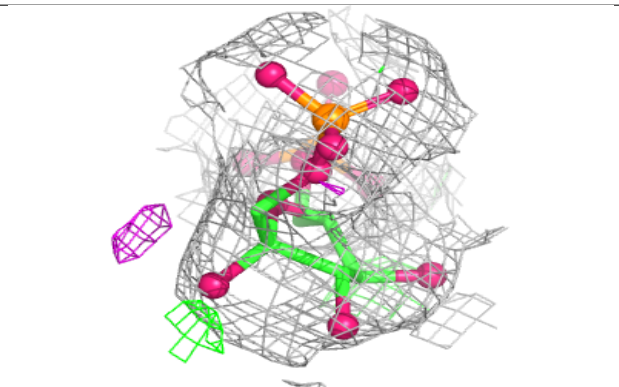
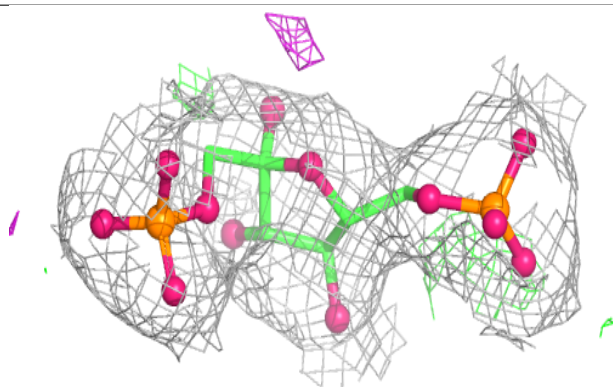
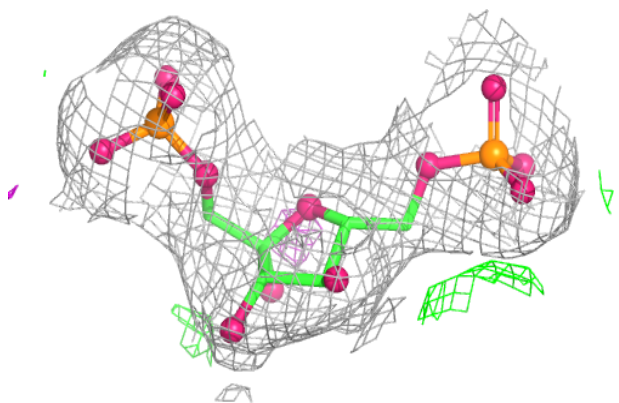
**Electron density around FBP C 601:**

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

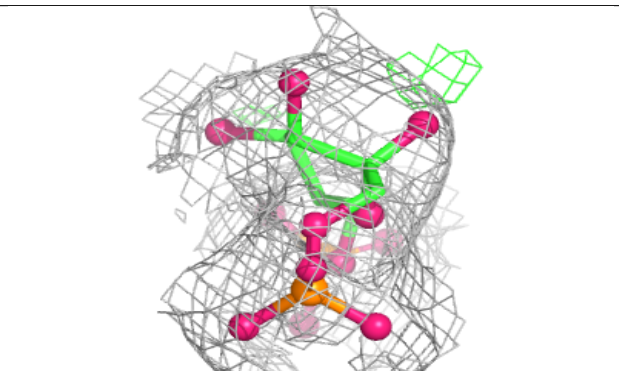
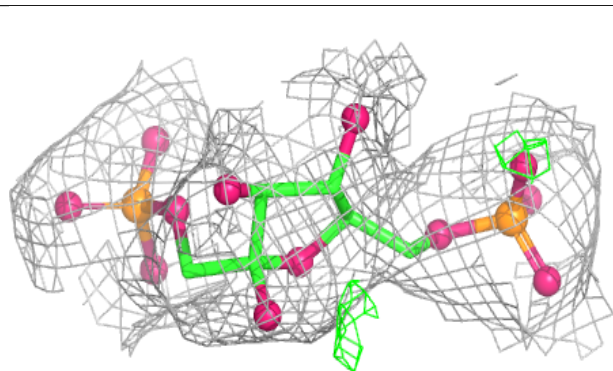
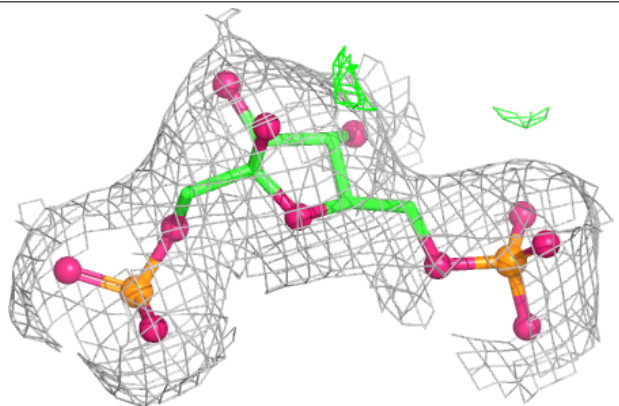


Electron density around FBP H 601:

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

**Electron density around FBP D 601:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.