



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:03 AM UTC

PDB ID : 7AUD / pdb_00007aud
Title : Structure of an engineered helicase domain construct for human Bloom syndrome protein (BLM)
Authors : Chen, X.; Oliver, A.W.
Deposited on : 2020-11-02
Resolution : 2.96 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

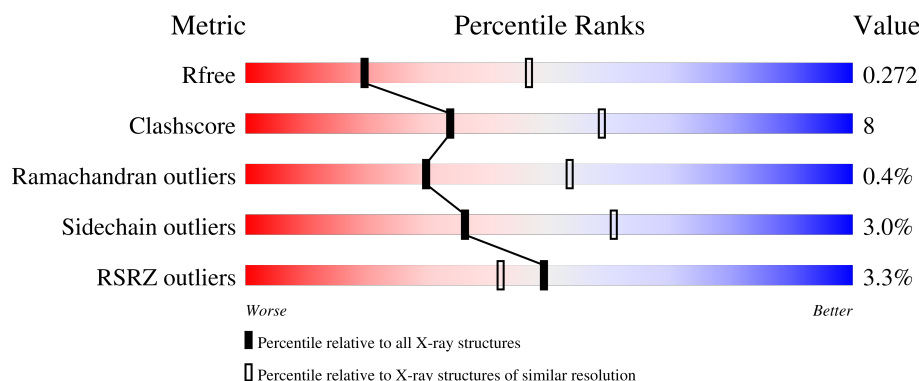
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.96 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	563	73% 19% • 7%
1	B	563	2% 75% 15% • 9%
1	C	563	69% 21% • 9%
1	D	563	4% 73% 17% 9%
1	E	563	4% 74% 16% • 9%

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Mol	Chain	Length	Quality of chain
1	F	563	6% 73% 18% 8%
2	M	15	47% 40% 13%
2	N	15	53% 33% 13%
2	O	15	73% 13% 13%
2	P	15	87% 13%
2	Q	15	60% 27% 13%
2	R	15	67% 20% 13%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 25271 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 Bloom syndrome protein, Bloom syndrome protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	522	Total	C	N	O	S	1	2	0
			4023	2551	703	737	32			
1	B	513	Total	C	N	O	S	1	0	0
			3890	2477	670	712	31			
1	C	513	Total	C	N	O	S	1	1	0
			3908	2483	683	710	32			
1	D	510	Total	C	N	O	S	4	1	0
			3792	2411	657	692	32			
1	E	513	Total	C	N	O	S	2	1	0
			3825	2427	665	703	30			
1	F	516	Total	C	N	O	S	4	1	0
			3750	2384	645	693	28			

There are 186 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	610	MET	-	initiating methionine	UNP P54132
A	611	GLY	-	expression tag	UNP P54132
A	612	SER	-	expression tag	UNP P54132
A	613	ALA	-	expression tag	UNP P54132
A	614	TRP	-	expression tag	UNP P54132
A	615	SER	-	expression tag	UNP P54132
A	616	HIS	-	expression tag	UNP P54132
A	617	PRO	-	expression tag	UNP P54132
A	618	GLN	-	expression tag	UNP P54132
A	619	PHE	-	expression tag	UNP P54132
A	620	GLU	-	expression tag	UNP P54132
A	621	LYS	-	expression tag	UNP P54132
A	622	SER	-	expression tag	UNP P54132
A	623	SER	-	expression tag	UNP P54132
A	624	GLY	-	expression tag	UNP P54132
A	625	LEU	-	expression tag	UNP P54132
A	626	GLU	-	expression tag	UNP P54132

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Chain	Residue	Modelled	Actual	Comment	Reference
A	627	VAL	-	expression tag	UNP P54132
A	628	LEU	-	expression tag	UNP P54132
A	629	PHE	-	expression tag	UNP P54132
A	630	GLN	-	expression tag	UNP P54132
A	631	GLY	-	expression tag	UNP P54132
A	632	GLY	-	expression tag	UNP P54132
A	633	PRO	-	expression tag	UNP P54132
A	634	HIS	-	expression tag	UNP P54132
A	635	MET	-	expression tag	UNP P54132
A	1197	GLY	-	linker	UNP P54132
A	1198	SER	-	linker	UNP P54132
A	1199	GLY	-	linker	UNP P54132
A	1200	GLY	-	linker	UNP P54132
A	1201	SER	-	linker	UNP P54132
B	610	MET	-	initiating methionine	UNP P54132
B	611	GLY	-	expression tag	UNP P54132
B	612	SER	-	expression tag	UNP P54132
B	613	ALA	-	expression tag	UNP P54132
B	614	TRP	-	expression tag	UNP P54132
B	615	SER	-	expression tag	UNP P54132
B	616	HIS	-	expression tag	UNP P54132
B	617	PRO	-	expression tag	UNP P54132
B	618	GLN	-	expression tag	UNP P54132
B	619	PHE	-	expression tag	UNP P54132
B	620	GLU	-	expression tag	UNP P54132
B	621	LYS	-	expression tag	UNP P54132
B	622	SER	-	expression tag	UNP P54132
B	623	SER	-	expression tag	UNP P54132
B	624	GLY	-	expression tag	UNP P54132
B	625	LEU	-	expression tag	UNP P54132
B	626	GLU	-	expression tag	UNP P54132
B	627	VAL	-	expression tag	UNP P54132
B	628	LEU	-	expression tag	UNP P54132
B	629	PHE	-	expression tag	UNP P54132
B	630	GLN	-	expression tag	UNP P54132
B	631	GLY	-	expression tag	UNP P54132
B	632	GLY	-	expression tag	UNP P54132
B	633	PRO	-	expression tag	UNP P54132
B	634	HIS	-	expression tag	UNP P54132
B	635	MET	-	expression tag	UNP P54132
B	1197	GLY	-	linker	UNP P54132
B	1198	SER	-	linker	UNP P54132

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Chain	Residue	Modelled	Actual	Comment	Reference
B	1199	GLY	-	linker	UNP P54132
B	1200	GLY	-	linker	UNP P54132
B	1201	SER	-	linker	UNP P54132
C	610	MET	-	initiating methionine	UNP P54132
C	611	GLY	-	expression tag	UNP P54132
C	612	SER	-	expression tag	UNP P54132
C	613	ALA	-	expression tag	UNP P54132
C	614	TRP	-	expression tag	UNP P54132
C	615	SER	-	expression tag	UNP P54132
C	616	HIS	-	expression tag	UNP P54132
C	617	PRO	-	expression tag	UNP P54132
C	618	GLN	-	expression tag	UNP P54132
C	619	PHE	-	expression tag	UNP P54132
C	620	GLU	-	expression tag	UNP P54132
C	621	LYS	-	expression tag	UNP P54132
C	622	SER	-	expression tag	UNP P54132
C	623	SER	-	expression tag	UNP P54132
C	624	GLY	-	expression tag	UNP P54132
C	625	LEU	-	expression tag	UNP P54132
C	626	GLU	-	expression tag	UNP P54132
C	627	VAL	-	expression tag	UNP P54132
C	628	LEU	-	expression tag	UNP P54132
C	629	PHE	-	expression tag	UNP P54132
C	630	GLN	-	expression tag	UNP P54132
C	631	GLY	-	expression tag	UNP P54132
C	632	GLY	-	expression tag	UNP P54132
C	633	PRO	-	expression tag	UNP P54132
C	634	HIS	-	expression tag	UNP P54132
C	635	MET	-	expression tag	UNP P54132
C	1197	GLY	-	linker	UNP P54132
C	1198	SER	-	linker	UNP P54132
C	1199	GLY	-	linker	UNP P54132
C	1200	GLY	-	linker	UNP P54132
C	1201	SER	-	linker	UNP P54132
D	610	MET	-	initiating methionine	UNP P54132
D	611	GLY	-	expression tag	UNP P54132
D	612	SER	-	expression tag	UNP P54132
D	613	ALA	-	expression tag	UNP P54132
D	614	TRP	-	expression tag	UNP P54132
D	615	SER	-	expression tag	UNP P54132
D	616	HIS	-	expression tag	UNP P54132
D	617	PRO	-	expression tag	UNP P54132

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Chain	Residue	Modelled	Actual	Comment	Reference
D	618	GLN	-	expression tag	UNP P54132
D	619	PHE	-	expression tag	UNP P54132
D	620	GLU	-	expression tag	UNP P54132
D	621	LYS	-	expression tag	UNP P54132
D	622	SER	-	expression tag	UNP P54132
D	623	SER	-	expression tag	UNP P54132
D	624	GLY	-	expression tag	UNP P54132
D	625	LEU	-	expression tag	UNP P54132
D	626	GLU	-	expression tag	UNP P54132
D	627	VAL	-	expression tag	UNP P54132
D	628	LEU	-	expression tag	UNP P54132
D	629	PHE	-	expression tag	UNP P54132
D	630	GLN	-	expression tag	UNP P54132
D	631	GLY	-	expression tag	UNP P54132
D	632	GLY	-	expression tag	UNP P54132
D	633	PRO	-	expression tag	UNP P54132
D	634	HIS	-	expression tag	UNP P54132
D	635	MET	-	expression tag	UNP P54132
D	1197	GLY	-	linker	UNP P54132
D	1198	SER	-	linker	UNP P54132
D	1199	GLY	-	linker	UNP P54132
D	1200	GLY	-	linker	UNP P54132
D	1201	SER	-	linker	UNP P54132
E	610	MET	-	initiating methionine	UNP P54132
E	611	GLY	-	expression tag	UNP P54132
E	612	SER	-	expression tag	UNP P54132
E	613	ALA	-	expression tag	UNP P54132
E	614	TRP	-	expression tag	UNP P54132
E	615	SER	-	expression tag	UNP P54132
E	616	HIS	-	expression tag	UNP P54132
E	617	PRO	-	expression tag	UNP P54132
E	618	GLN	-	expression tag	UNP P54132
E	619	PHE	-	expression tag	UNP P54132
E	620	GLU	-	expression tag	UNP P54132
E	621	LYS	-	expression tag	UNP P54132
E	622	SER	-	expression tag	UNP P54132
E	623	SER	-	expression tag	UNP P54132
E	624	GLY	-	expression tag	UNP P54132
E	625	LEU	-	expression tag	UNP P54132
E	626	GLU	-	expression tag	UNP P54132
E	627	VAL	-	expression tag	UNP P54132
E	628	LEU	-	expression tag	UNP P54132

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Chain	Residue	Modelled	Actual	Comment	Reference
E	629	PHE	-	expression tag	UNP P54132
E	630	GLN	-	expression tag	UNP P54132
E	631	GLY	-	expression tag	UNP P54132
E	632	GLY	-	expression tag	UNP P54132
E	633	PRO	-	expression tag	UNP P54132
E	634	HIS	-	expression tag	UNP P54132
E	635	MET	-	expression tag	UNP P54132
E	1197	GLY	-	linker	UNP P54132
E	1198	SER	-	linker	UNP P54132
E	1199	GLY	-	linker	UNP P54132
E	1200	GLY	-	linker	UNP P54132
E	1201	SER	-	linker	UNP P54132
F	610	MET	-	initiating methionine	UNP P54132
F	611	GLY	-	expression tag	UNP P54132
F	612	SER	-	expression tag	UNP P54132
F	613	ALA	-	expression tag	UNP P54132
F	614	TRP	-	expression tag	UNP P54132
F	615	SER	-	expression tag	UNP P54132
F	616	HIS	-	expression tag	UNP P54132
F	617	PRO	-	expression tag	UNP P54132
F	618	GLN	-	expression tag	UNP P54132
F	619	PHE	-	expression tag	UNP P54132
F	620	GLU	-	expression tag	UNP P54132
F	621	LYS	-	expression tag	UNP P54132
F	622	SER	-	expression tag	UNP P54132
F	623	SER	-	expression tag	UNP P54132
F	624	GLY	-	expression tag	UNP P54132
F	625	LEU	-	expression tag	UNP P54132
F	626	GLU	-	expression tag	UNP P54132
F	627	VAL	-	expression tag	UNP P54132
F	628	LEU	-	expression tag	UNP P54132
F	629	PHE	-	expression tag	UNP P54132
F	630	GLN	-	expression tag	UNP P54132
F	631	GLY	-	expression tag	UNP P54132
F	632	GLY	-	expression tag	UNP P54132
F	633	PRO	-	expression tag	UNP P54132
F	634	HIS	-	expression tag	UNP P54132
F	635	MET	-	expression tag	UNP P54132
F	1197	GLY	-	linker	UNP P54132
F	1198	SER	-	linker	UNP P54132
F	1199	GLY	-	linker	UNP P54132
F	1200	GLY	-	linker	UNP P54132

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Chain	Residue	Modelled	Actual	Comment	Reference
F	1201	SER	-	linker	UNP P54132

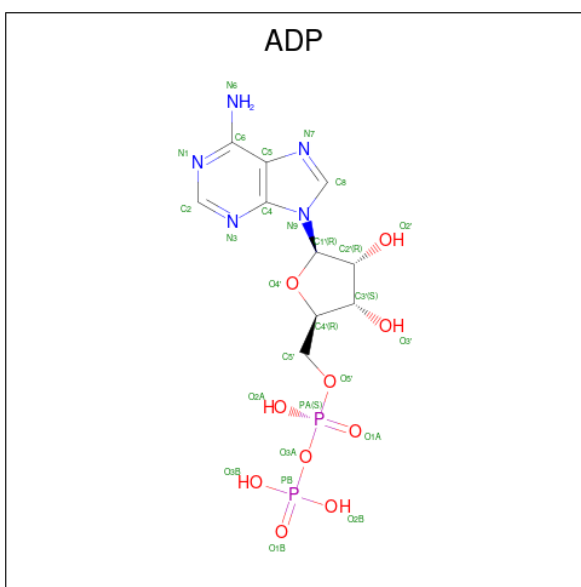
- Molecule 2 is a DNA chain called DNA (5'-D(*GP*TP*AP*CP*CP*CP*GP*AP*TP*GP*TP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	13	Total	C	N	O	P	0	0	0
			264	127	47	78	12			
2	N	13	Total	C	N	O	P	0	0	0
			264	127	47	78	12			
2	O	13	Total	C	N	O	P	0	0	0
			264	127	47	78	12			
2	P	13	Total	C	N	O	P	0	0	0
			267	127	47	80	13			
2	Q	13	Total	C	N	O	P	0	0	0
			267	127	47	80	13			
2	R	13	Total	C	N	O	P	0	0	0
			264	127	47	78	12			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

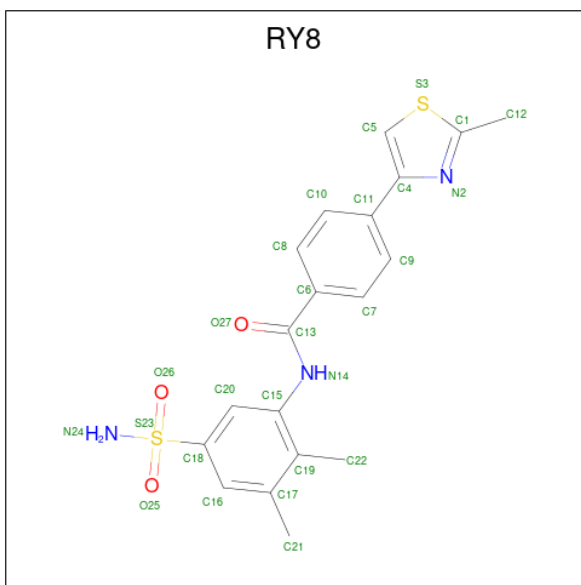
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		
3	E	1	Total	Zn	0	0
			1	1		
3	F	1	Total	Zn	0	0
			1	1		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



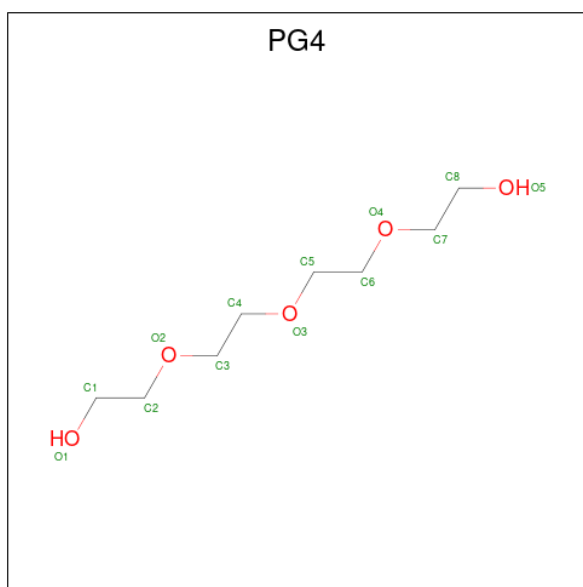
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 5 is N-(2,3-dimethyl-5-sulfamoylphenyl)-4-(2-methylthiazol-4-yl)benzamide (CCD ID: RY8) (formula: C₁₉H₁₉N₃O₃S₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			27	19	3	3	2		
5	B	1	Total	C	N	O	S	0	0
			27	19	3	3	2		
5	C	1	Total	C	N	O	S	0	0
			27	19	3	3	2		
5	D	1	Total	C	N	O	S	0	0
			27	19	3	3	2		
5	E	1	Total	C	N	O	S	0	0
			27	19	3	3	2		
5	F	1	Total	C	N	O	S	0	0
			27	19	3	3	2		

- Molecule 6 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			13	8	5		
6	A	1	Total	C	O	0	0
			13	8	5		
6	B	1	Total	C	O	0	0
			13	8	5		
6	C	1	Total	C	O	0	0
			13	8	5		
6	C	1	Total	C	O	0	0
			13	8	5		
6	D	1	Total	C	O	0	0
			13	8	5		
6	E	1	Total	C	O	0	0
			13	8	5		

- Molecule 7 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	F	1	Total	Mg	0	0
			1	1		
7	M	1	Total	Mg	0	0
			1	1		
7	N	1	Total	Mg	0	0
			1	1		
7	P	1	Total	Mg	0	0
			1	1		
7	Q	2	Total	Mg	0	0
			2	2		

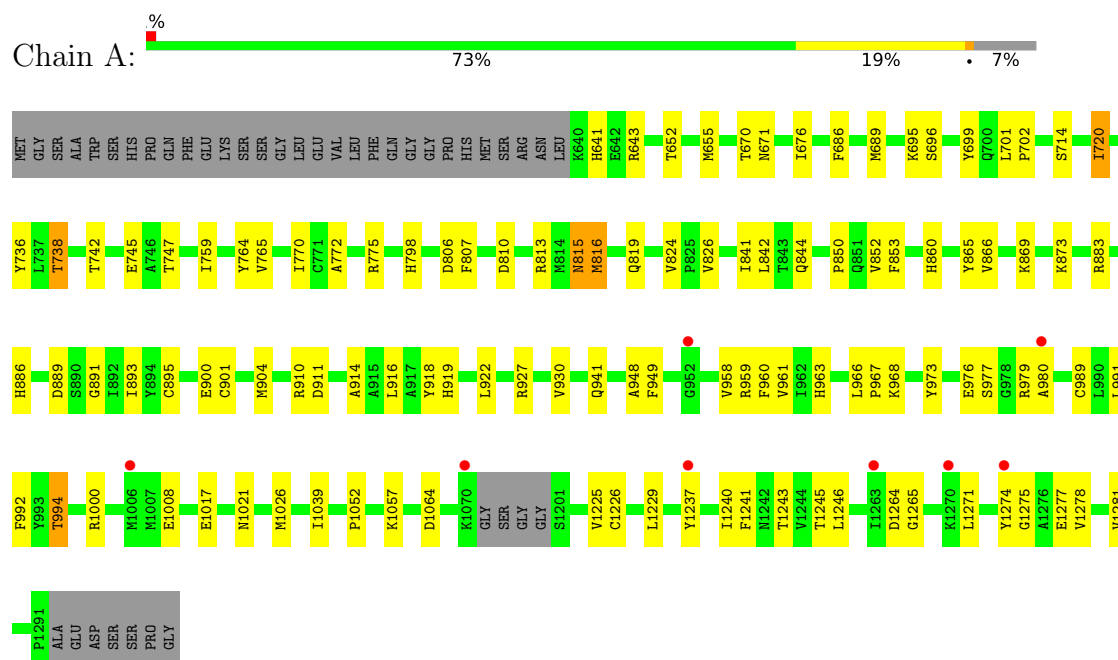
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	31	Total O 31 31	0	0
8	B	16	Total O 16 16	0	0
8	C	5	Total O 5 5	0	0
8	D	4	Total O 4 4	0	0
8	E	6	Total O 6 6	0	0
8	F	1	Total O 1 1	0	0
8	M	1	Total O 1 1	0	0
8	Q	2	Total O 2 2	0	0

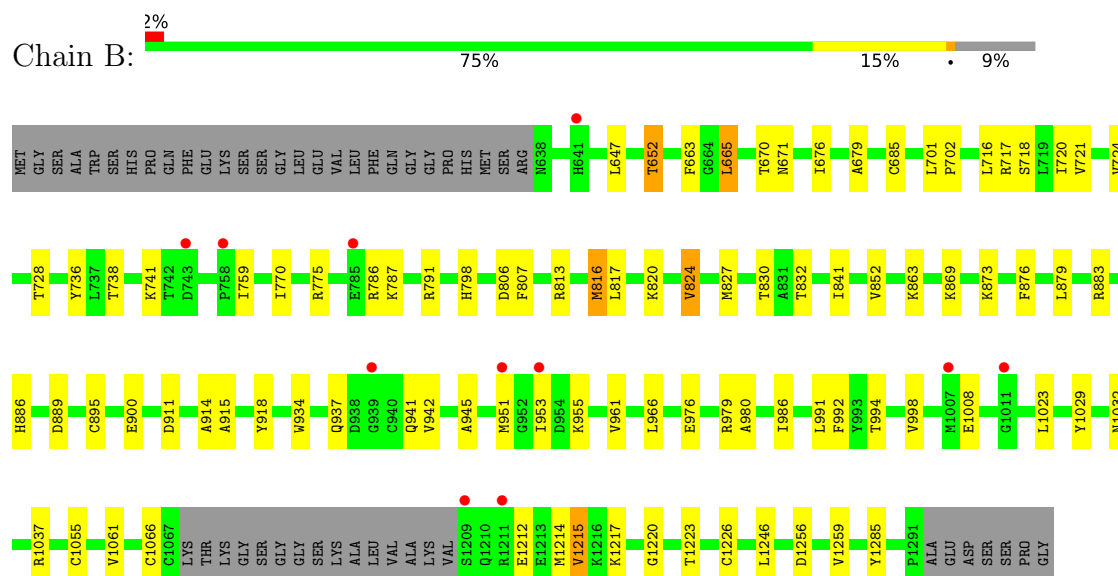
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

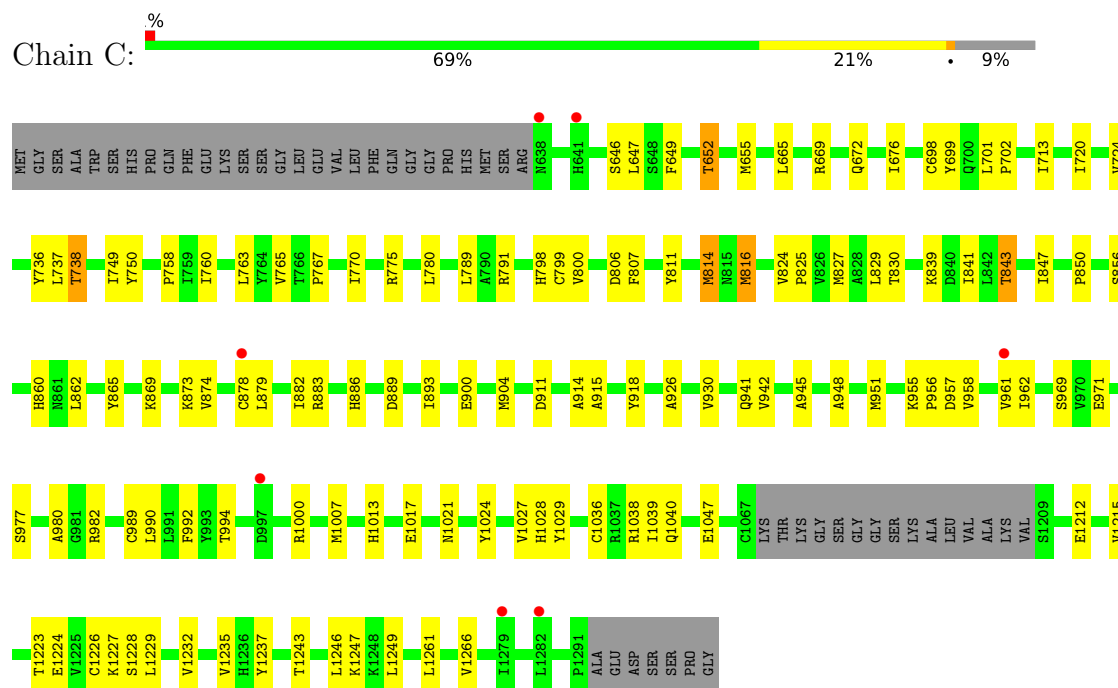
- Molecule 1: Bloom syndrome protein,Bloom syndrome protein



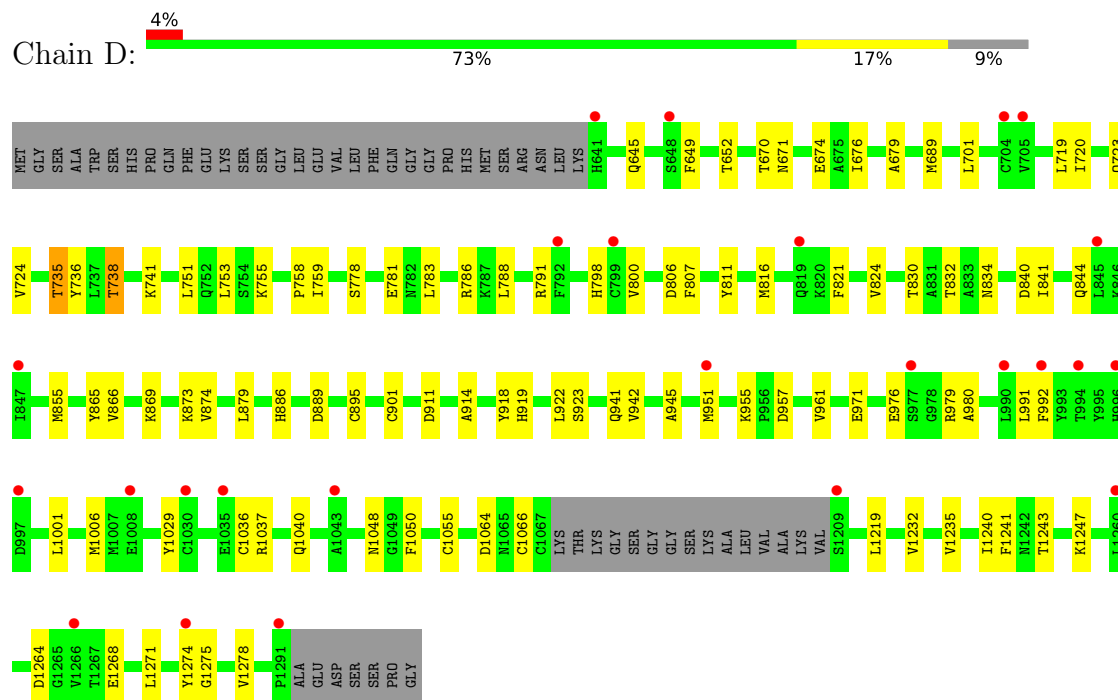
- Molecule 1: Bloom syndrome protein,Bloom syndrome protein



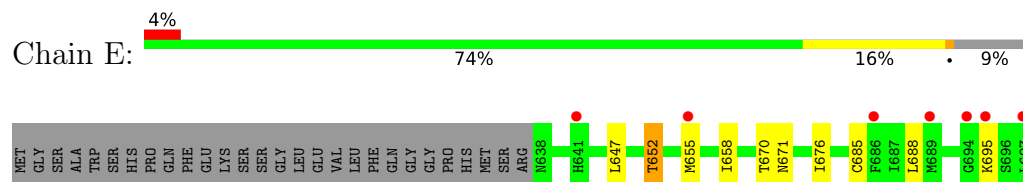
• Molecule 1: Bloom syndrome protein,Bloom syndrome protein

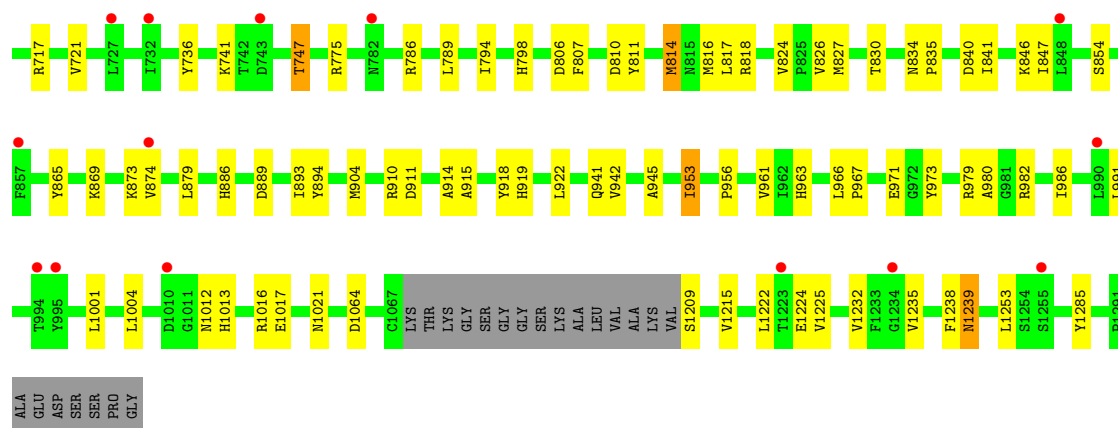


• Molecule 1: Bloom syndrome protein,Bloom syndrome protein

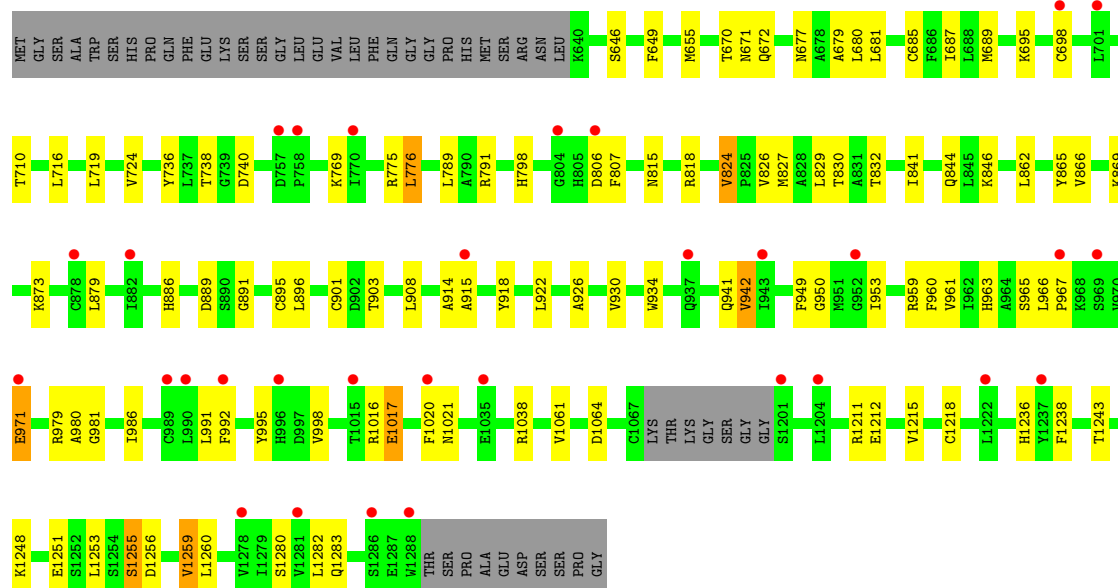
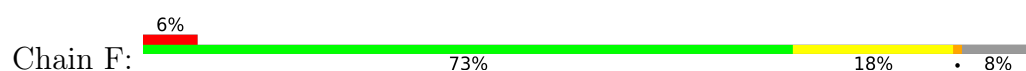


• Molecule 1: Bloom syndrome protein,Bloom syndrome protein





- Molecule 1: Bloom syndrome protein, Bloom syndrome protein



- Molecule 2: DNA (5'-D(*GP*TP*AP*CP*CP*CP*GP*AP*TP*GP*TP*GP*T)-3')



- Molecule 2: DNA (5'-D(*GP*TP*AP*CP*CP*CP*GP*AP*TP*GP*TP*GP*T)-3')



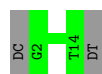
- Molecule 2: DNA (5'-D(*GP*TP*AP*CP*CP*CP*GP*AP*TP*GP*TP*GP*T)-3')

Chain O:  73% 13% 13%



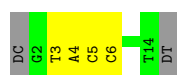
● Molecule 2: DNA (5'-D(*GP*TP*AP*CP*CP*CP*GP*AP*TP*GP*TP*GP*T)-3')

Chain P:  87% 13%



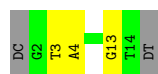
● Molecule 2: DNA (5'-D(*GP*TP*AP*CP*CP*CP*GP*AP*TP*GP*TP*GP*T)-3')

Chain Q:  60% 27% 13%



● Molecule 2: DNA (5'-D(*GP*TP*AP*CP*CP*CP*GP*AP*TP*GP*TP*GP*T)-3')

Chain R:  67% 20% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	84.92Å 111.71Å 132.34Å 72.64° 80.18° 79.22°	Depositor
Resolution (Å)	47.28 – 2.96 47.28 – 2.96	Depositor EDS
% Data completeness (in resolution range)	97.6 (47.28-2.96) 92.3 (47.28-2.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.229 , 0.269 0.233 , 0.272	Depositor DCC
R_{free} test set	4642 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	69.9	Xtriage
Anisotropy	0.428	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 73.2	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	0.92	EDS
Total number of atoms	25271	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.76% 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: RY8, MG, ADP, PG4, ZN

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	0.13	0/4117	0.35	0/5590
1	B	0.12	0/3978	0.34	0/5415
1	C	0.14	0/3999	0.36	0/5439
1	D	0.12	0/3882	0.34	0/5290
1	E	0.13	0/3915	0.35	0/5338
1	F	0.14	0/3838	0.35	0/5247
2	M	0.20	0/295	0.42	0/454
2	N	0.21	0/295	0.47	0/454
2	O	0.22	0/295	0.43	0/454
2	P	0.21	0/298	0.42	0/458
2	Q	0.21	0/298	0.42	0/458
2	R	0.23	0/295	0.42	0/454
All	All	0.14	0/25505	0.35	0/35051

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	4023	0	3844	66	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3890	0	3651	54	0
1	C	3908	0	3689	71	0
1	D	3792	0	3481	57	0
1	E	3825	0	3517	51	0
1	F	3750	0	3370	63	0
2	M	264	0	149	7	0
2	N	264	0	149	4	0
2	O	264	0	149	2	0
2	P	267	0	148	0	0
2	Q	267	0	148	3	0
2	R	264	0	149	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	27	0	12	0	0
4	B	27	0	12	0	0
4	C	27	0	12	1	0
4	D	27	0	12	0	0
4	E	27	0	12	0	0
4	F	27	0	12	1	0
5	A	27	0	0	0	0
5	B	27	0	0	0	0
5	C	27	0	0	0	0
5	D	27	0	0	1	0
5	E	27	0	0	2	0
5	F	27	0	0	0	0
6	A	26	0	36	2	0
6	B	13	0	18	1	0
6	C	26	0	36	2	0
6	D	13	0	18	1	0
6	E	13	0	18	1	0
7	F	1	0	0	0	0
7	M	1	0	0	0	0
7	N	1	0	0	0	0
7	P	1	0	0	0	0
7	Q	2	0	0	0	0
8	A	31	0	0	0	0
8	B	16	0	0	0	0
8	C	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	D	4	0	0	0	0
8	E	6	0	0	0	0
8	F	1	0	0	0	0
8	M	1	0	0	0	0
8	Q	2	0	0	0	0
All	All	25271	0	22642	367	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 (367) 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:669:ARG:H	1:C:672:GLN:HE21	1.25	0.83
1:A:1225:VAL:HG21	1:A:1281:VAL:HG11	1.59	0.83
1:C:737:LEU:HD13	1:C:763:LEU:HD11	1.68	0.75
1:F:679:ALA:O	1:F:791:ARG:NH2	2.21	0.74
1:B:1008:GLU:OE2	2:N:7:DC:N4	2.22	0.73
1:C:791[B]:ARG:HG2	1:C:825:PRO:HG2	1.70	0.72
1:C:672:GLN:HE22	4:C:1302:ADP:HN61	1.39	0.71
1:C:737:LEU:HD11	1:C:749:ILE:HG21	1.73	0.69
1:F:886:HIS:HB3	1:F:889:ASP:HB2	1.75	0.69
1:F:710:THR:HG23	1:F:791:ARG:HB2	1.75	0.68
1:E:816:MET:HE2	6:E:1304:PG4:H32	1.75	0.68
1:B:679:ALA:O	1:B:791:ARG:NH2	2.26	0.68
1:F:738:THR:HG23	1:F:740:ASP:H	1.59	0.68
1:F:866:VAL:HG22	1:F:991:LEU:HB3	1.77	0.67
1:D:676:ILE:HD13	1:D:701:LEU:HD23	1.76	0.67
1:A:815:ASN:ND2	1:A:844:GLN:O	2.26	0.66
1:E:956:PRO:HA	1:E:982:ARG:HB2	1.78	0.66
1:F:966:LEU:HA	1:F:991:LEU:HD11	1.77	0.66
1:D:755:LYS:O	1:D:786:ARG:NH2	2.22	0.65
1:C:839:LYS:O	1:C:843:THR:OG1	2.15	0.65
1:D:679:ALA:O	1:D:791:ARG:NH2	2.31	0.64
1:A:747:THR:HG22	1:A:775:ARG:HH22	1.63	0.63
1:D:649:PHE:HB2	1:D:652:THR:HG23	1.81	0.63
1:A:961:VAL:HG23	1:A:980:ALA:HB2	1.78	0.62
1:B:869:LYS:HD3	1:B:994:THR:HG22	1.81	0.62
1:F:915:ALA:HA	1:F:942:VAL:O	1.98	0.62
1:C:758:PRO:O	1:C:760:ILE:N	2.31	0.62
1:C:869:LYS:HE3	1:C:874:VAL:HG11	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:869:LYS:HB2	1:C:992:PHE:HB3	1.82	0.62
1:D:886:HIS:HB3	1:D:889:ASP:HB2	1.80	0.62
1:F:1236:HIS:HD2	1:F:1238:PHE:H	1.45	0.61
1:A:1237:TYR:HA	1:A:1240:ILE:HD12	1.82	0.61
1:E:655:MET:HG3	1:E:701:LEU:HD21	1.82	0.61
1:A:916:LEU:HD23	1:A:930:VAL:HG22	1.83	0.61
1:A:900:GLU:OE1	1:A:1000:ARG:NH2	2.27	0.61
1:D:786:ARG:HE	1:D:788:LEU:HD12	1.66	0.60
1:A:886:HIS:HB3	1:A:889:ASP:HB2	1.83	0.60
1:B:1029:TYR:O	1:B:1037:ARG:NH2	2.35	0.60
1:C:1243:THR:OG1	1:C:1247:LYS:NZ	2.35	0.60
1:A:772:ALA:HA	6:A:1304:PG4:H72	1.83	0.60
1:E:717:ARG:O	1:E:721:VAL:HG23	2.01	0.60
1:C:676:ILE:HD13	1:C:701:LEU:HD23	1.83	0.59
1:E:1239:ASN:O	1:E:1239:ASN:ND2	2.30	0.59
1:A:676:ILE:HD13	1:A:701:LEU:HD23	1.84	0.59
1:B:883:ARG:NH1	1:B:911:ASP:O	2.35	0.59
1:A:860:HIS:HB2	6:A:1305:PG4:H52	1.85	0.59
1:F:959:ARG:NH1	1:F:981:GLY:HA2	2.17	0.58
1:A:742:THR:HG23	1:A:745:GLU:H	1.69	0.58
1:B:976:GLU:HA	1:B:979:ARG:NH1	2.19	0.58
1:C:713:ILE:HG21	1:C:814:MET:HE2	1.86	0.58
1:F:934:TRP:HZ3	1:F:953:ILE:HG21	1.68	0.58
1:B:1217:LYS:HE2	1:B:1285:TYR:CZ	2.38	0.58
1:C:955:LYS:NZ	1:C:957:ASP:OD1	2.35	0.58
1:E:961:VAL:HG23	1:E:980:ALA:HB2	1.84	0.58
1:F:687:ILE:HD12	1:F:827:MET:HE3	1.84	0.58
1:D:1055:CYS:CB	1:D:1066:CYS:SG	2.92	0.58
1:A:948:ALA:HB2	2:M:8:DG:H5"	1.85	0.57
1:C:816:MET:HE2	6:C:1305:PG4:H51	1.86	0.57
1:E:685:CYS:HB2	1:E:827:MET:HG3	1.86	0.57
1:C:736:TYR:HD1	1:C:738:THR:HG23	1.68	0.57
1:F:869:LYS:HB2	1:F:992:PHE:HB3	1.86	0.57
1:F:815:ASN:ND2	1:F:844:GLN:O	2.37	0.57
1:F:979:ARG:HH21	1:F:979:ARG:HG3	1.69	0.57
1:A:1008:GLU:OE2	2:M:7:DC:N4	2.38	0.57
1:D:1264:ASP:OD1	1:D:1264:ASP:N	2.37	0.57
1:A:869:LYS:HB2	1:A:992:PHE:HB3	1.87	0.56
1:C:961:VAL:HG23	1:C:980:ALA:HB2	1.87	0.56
1:B:813:ARG:HD3	6:B:1304:PG4:H22	1.86	0.56
1:D:778:SER:O	1:D:781:GLU:HG2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:919:HIS:H	1:D:922:LEU:HD13	1.69	0.56
1:E:846:LYS:NZ	1:F:889:ASP:OD1	2.33	0.56
1:A:652:THR:HA	1:A:655:MET:HE2	1.88	0.56
1:A:967:PRO:HG2	1:A:973:TYR:HB2	1.88	0.56
1:B:1212:GLU:O	1:B:1215:VAL:HG12	2.06	0.56
1:C:1013:HIS:O	1:C:1017:GLU:HG2	2.06	0.55
1:E:647:LEU:HA	1:E:652:THR:HG21	1.88	0.55
1:E:798:HIS:HA	1:E:841:ILE:HD11	1.87	0.55
1:E:966:LEU:HA	1:E:991:LEU:HD11	1.87	0.55
1:C:926:ALA:O	1:C:930:VAL:HG12	2.06	0.55
1:C:652:THR:HA	1:C:655:MET:HE2	1.87	0.55
1:F:775:ARG:HD2	2:R:13:DG:C5	2.42	0.55
1:F:926:ALA:O	1:F:930:VAL:HG12	2.07	0.55
1:C:1028:HIS:O	1:C:1029:TYR:HB3	2.07	0.55
1:A:736:TYR:HD2	1:A:738:THR:HG23	1.72	0.54
1:F:685:CYS:HB2	1:F:827:MET:HG3	1.90	0.54
1:F:1211:ARG:O	1:F:1215:VAL:HG13	2.08	0.54
1:C:869:LYS:HD3	1:C:994:THR:HG22	1.89	0.54
1:F:914:ALA:O	1:F:941:GLN:N	2.38	0.54
1:A:869:LYS:HD3	1:A:994:THR:HG22	1.88	0.54
1:D:830:THR:HG22	1:D:832:THR:H	1.72	0.54
1:A:676:ILE:HG23	1:A:702:PRO:HG3	1.90	0.54
1:C:669:ARG:H	1:C:672:GLN:NE2	2.01	0.54
1:D:1029:TYR:O	1:D:1037:ARG:NH1	2.40	0.54
1:A:686:PHE:HB3	1:A:852:VAL:HG12	1.90	0.54
1:B:770:ILE:HD13	1:B:817:LEU:HD11	1.89	0.54
1:C:798:HIS:HA	1:C:841:ILE:HD11	1.89	0.53
1:D:869:LYS:HB2	1:D:992:PHE:HB3	1.90	0.53
1:A:966:LEU:HA	1:A:991:LEU:HD11	1.90	0.53
1:F:769:LYS:HG3	1:F:776:LEU:HD12	1.89	0.53
1:B:961:VAL:HG23	1:B:980:ALA:HB2	1.90	0.53
1:C:1243:THR:O	1:C:1247:LYS:HG2	2.09	0.53
1:E:1012:ASN:O	1:E:1016:ARG:HD3	2.09	0.53
1:B:869:LYS:HB2	1:B:992:PHE:HB3	1.91	0.52
1:A:866:VAL:HG22	1:A:991:LEU:HB3	1.91	0.52
1:B:716:LEU:O	1:B:720:ILE:HG23	2.09	0.52
1:E:658:ILE:HD13	1:E:705:VAL:HG22	1.91	0.52
1:A:689:MET:HE3	1:A:695:LYS:HA	1.92	0.52
1:A:798:HIS:HA	1:A:841:ILE:HD11	1.90	0.52
1:B:886:HIS:HB3	1:B:889:ASP:HB2	1.89	0.52
2:R:3:DT:H2"	2:R:4:DA:C8	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:816:MET:SD	1:B:820:LYS:HE2	2.50	0.52
1:F:950:GLY:HA3	1:F:979:ARG:NH2	2.25	0.52
1:A:893:ILE:HG21	1:A:904:MET:HE3	1.92	0.52
1:A:914:ALA:O	1:A:941:GLN:N	2.41	0.52
1:C:948:ALA:HB2	2:O:8:DG:H5"	1.92	0.52
1:D:1055:CYS:HB2	1:D:1066:CYS:SG	2.50	0.52
2:M:3:DT:H2"	2:M:4:DA:C8	2.45	0.52
1:D:869:LYS:HE3	1:D:874:VAL:HG11	1.91	0.51
2:Q:3:DT:H2"	2:Q:4:DA:N7	2.25	0.51
1:B:717:ARG:O	1:B:720:ILE:HG12	2.11	0.51
1:A:641:HIS:HD2	1:A:853:PHE:CE1	2.27	0.51
1:D:914:ALA:O	1:D:941:GLN:N	2.34	0.51
1:C:886:HIS:HB3	1:C:889:ASP:HB2	1.92	0.51
1:D:976:GLU:HA	1:D:979:ARG:NH2	2.26	0.51
1:D:736:TYR:HD1	1:D:738:THR:HG23	1.75	0.51
1:E:886:HIS:HB3	1:E:889:ASP:HB2	1.91	0.51
1:E:794:ILE:HD12	1:E:814:MET:HE1	1.93	0.51
1:C:883:ARG:NH1	1:C:911:ASP:O	2.43	0.51
1:D:736:TYR:CD1	1:D:738:THR:HG23	2.47	0.50
1:C:893:ILE:HG21	1:C:904:MET:HE3	1.94	0.50
1:B:1256:ASP:HB3	1:B:1259:VAL:HG12	1.93	0.50
1:D:976:GLU:HA	1:D:979:ARG:HH21	1.77	0.50
1:E:967:PRO:HG2	1:E:973:TYR:HB2	1.93	0.50
1:F:716:LEU:HD12	1:F:719:LEU:HD23	1.93	0.50
1:E:1222:LEU:O	1:E:1225:VAL:HG22	2.12	0.50
1:B:676:ILE:HD13	1:B:701:LEU:HD23	1.93	0.50
1:E:894:TYR:OH	1:E:979:ARG:NH1	2.45	0.50
1:B:879:LEU:HD11	1:B:911:ASP:HB3	1.94	0.49
1:E:834:ASN:HB2	1:E:835:PRO:HD2	1.93	0.49
1:B:685:CYS:HB2	1:B:827:MET:HG3	1.92	0.49
1:E:919:HIS:H	1:E:922:LEU:HD23	1.77	0.49
1:A:883[A]:ARG:NH2	1:A:911:ASP:OD2	2.46	0.49
1:D:751:LEU:O	1:D:755:LYS:HG3	2.12	0.49
1:F:862:LEU:O	1:F:1038:ARG:NH1	2.46	0.49
1:C:1036:CYS:O	1:C:1039:ILE:HG22	2.13	0.49
1:A:736:TYR:CD2	1:A:738:THR:HG23	2.48	0.49
1:F:966:LEU:HD21	1:F:998:VAL:HG22	1.95	0.49
1:A:810:ASP:HA	1:A:813:ARG:HB2	1.95	0.49
1:D:689:MET:HE2	1:D:855:MET:HE1	1.95	0.49
1:B:998:VAL:HG13	1:B:1023:LEU:HD21	1.94	0.49
1:E:1013:HIS:O	1:E:1017:GLU:HG2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:655:MET:HE2	1:F:680:LEU:HD12	1.95	0.48
1:A:968:LYS:NZ	2:M:7:DC:N3	2.58	0.48
1:C:765:VAL:HG21	1:C:770:ILE:HD11	1.94	0.48
1:B:895:CYS:HB3	1:B:900:GLU:HB3	1.95	0.48
1:B:798:HIS:HA	1:B:841:ILE:HD11	1.94	0.48
1:D:1268:GLU:HA	1:D:1271:LEU:HG	1.96	0.47
1:F:818:ARG:HD3	1:F:826:VAL:HG21	1.95	0.47
1:C:915:ALA:HA	1:C:942:VAL:O	2.14	0.47
1:E:806:ASP:O	1:E:807:PHE:HB3	2.15	0.47
1:A:720:ILE:HD13	1:A:764:TYR:HB3	1.95	0.47
2:Q:5:DC:H5"	2:Q:6:DC:OP2	2.14	0.47
1:C:1249:LEU:HD21	1:C:1266:VAL:HG21	1.96	0.47
1:E:676:ILE:HG23	1:E:702:PRO:HG3	1.95	0.47
1:F:896:LEU:HB2	1:F:965:SER:OG	2.15	0.47
1:B:966:LEU:HA	1:B:991:LEU:HD11	1.97	0.47
1:C:1024:TYR:HA	1:C:1027:VAL:HG12	1.95	0.47
1:F:689:MET:O	1:F:695:LYS:HE3	2.14	0.47
1:D:1275:GLY:HA2	1:D:1278:VAL:HG12	1.97	0.47
1:E:953:ILE:O	1:E:953:ILE:HG13	2.15	0.47
1:F:959:ARG:HH11	1:F:981:GLY:HA2	1.80	0.47
1:B:1220:GLY:O	1:B:1223:THR:HG22	2.15	0.47
1:B:863:LYS:HG2	1:B:1061:VAL:HA	1.97	0.46
1:F:995:TYR:HA	1:F:998:VAL:HG23	1.97	0.46
1:C:1228:SER:O	1:C:1232:VAL:HG13	2.15	0.46
1:F:1016:ARG:HB3	1:F:1020:PHE:CE2	2.50	0.46
1:C:698:CYS:HB3	1:C:827:MET:HE1	1.97	0.46
1:E:915:ALA:HA	1:E:942:VAL:O	2.16	0.46
1:C:806:ASP:O	1:C:807:PHE:HB3	2.16	0.46
1:A:891:GLY:HA3	1:A:960:PHE:CZ	2.50	0.46
1:B:934:TRP:CH2	1:B:955:LYS:HD2	2.50	0.46
1:D:806:ASP:O	1:D:807:PHE:HB3	2.16	0.46
1:A:976:GLU:HA	1:A:979:ARG:HH21	1.81	0.46
1:C:736:TYR:CD1	1:C:738:THR:HG23	2.48	0.46
1:D:865:TYR:CD1	1:D:1064:ASP:HB3	2.51	0.46
1:D:1240:ILE:HG22	1:D:1241:PHE:CD1	2.51	0.46
1:F:966:LEU:HD21	1:F:998:VAL:CG2	2.46	0.46
1:C:958:VAL:HB	1:C:980:ALA:HA	1.98	0.46
1:D:961:VAL:HG23	1:D:980:ALA:HB2	1.98	0.46
1:F:961:VAL:HG23	1:F:980:ALA:HB2	1.97	0.46
1:D:719:LEU:HD22	1:D:723:GLN:HE21	1.81	0.46
1:F:775:ARG:HD2	2:R:13:DG:C4	2.52	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:891:GLY:HA3	1:F:960:PHE:CZ	2.51	0.46
1:F:1212:GLU:O	1:F:1215:VAL:HG22	2.16	0.46
1:A:670:THR:HG22	1:A:671:ASN:ND2	2.31	0.45
1:E:816:MET:HE3	1:E:816:MET:HB3	1.89	0.45
1:B:791:ARG:HA	1:B:824:VAL:HG13	1.98	0.45
1:D:1048:ASN:C	1:D:1050:PHE:H	2.23	0.45
1:F:670:THR:HG22	1:F:671:ASN:ND2	2.31	0.45
1:F:865:TYR:CD2	1:F:1064:ASP:HB3	2.51	0.45
1:A:895:CYS:HB2	1:A:901:CYS:SG	2.57	0.45
1:E:810:ASP:O	1:E:814:MET:HG2	2.17	0.45
1:F:963:HIS:CD2	1:F:967:PRO:HG3	2.52	0.45
1:A:720:ILE:HG23	1:A:764:TYR:HB3	1.98	0.45
1:F:798:HIS:HA	1:F:841:ILE:HD11	1.97	0.45
2:Q:3:DT:H2"	2:Q:4:DA:C8	2.51	0.45
1:F:966:LEU:HD12	1:F:967:PRO:O	2.16	0.45
1:B:736:TYR:CZ	1:B:741:LYS:HE2	2.52	0.45
1:B:775:ARG:HD3	2:N:13:DG:C8	2.52	0.45
1:D:798:HIS:HA	1:D:841:ILE:HD11	1.97	0.45
1:B:1032:ASN:O	1:B:1037:ARG:NH1	2.50	0.44
1:D:840:ASP:O	1:D:844:GLN:HG2	2.17	0.44
1:E:893:ILE:HG21	1:E:904:MET:HE2	1.97	0.44
1:F:1017:GLU:O	1:F:1021:ASN:ND2	2.50	0.44
1:A:865:TYR:CD1	1:A:1064:ASP:HB3	2.52	0.44
1:B:876:PHE:O	1:B:879:LEU:HB3	2.17	0.44
1:B:915:ALA:HA	1:B:942:VAL:O	2.16	0.44
1:C:1224:GLU:O	1:C:1227:LYS:HG2	2.16	0.44
1:D:674:GLU:OE1	1:D:674:GLU:N	2.40	0.44
1:E:879:LEU:HD11	1:E:911:ASP:HB2	1.98	0.44
1:D:798:HIS:CD2	1:D:830:THR:HG23	2.52	0.44
1:E:1001:LEU:HD23	1:E:1004:LEU:HD12	2.00	0.44
1:A:1240:ILE:HG23	1:A:1274:TYR:CD2	2.52	0.44
1:B:806:ASP:O	1:B:807:PHE:HB3	2.18	0.44
1:C:1226:CYS:HB3	1:C:1237:TYR:HB2	1.99	0.44
1:E:670:THR:HG22	1:E:671:ASN:ND2	2.33	0.44
1:F:724:VAL:HG11	1:F:736:TYR:CD1	2.52	0.44
1:F:895:CYS:HB2	1:F:901:CYS:SG	2.58	0.44
1:C:1007:MET:HE1	2:O:6:DC:C2	2.53	0.44
1:C:1223:THR:HG22	1:C:1246:LEU:HD13	1.99	0.44
1:D:1006:MET:HE2	1:D:1006:MET:HB3	1.89	0.44
1:E:789:LEU:HD23	1:E:824:VAL:HG21	2.00	0.44
1:B:670:THR:HG22	1:B:671:ASN:ND2	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:1248:LYS:O	1:F:1251:GLU:HG2	2.17	0.44
1:A:806:ASP:O	1:A:807:PHE:HB3	2.17	0.43
1:A:1039:ILE:HD11	1:A:1052:PRO:HA	2.00	0.43
1:B:676:ILE:HG23	1:B:702:PRO:HG3	1.99	0.43
1:C:647:LEU:HA	1:C:652:THR:HG21	2.00	0.43
1:C:962:ILE:HG12	1:C:990:LEU:HB2	2.00	0.43
1:E:1239:ASN:OD1	2:M:13:DG:H5'	2.18	0.43
1:C:814:MET:O	1:C:814:MET:HG2	2.16	0.43
1:D:821:PHE:O	1:D:824:VAL:HG12	2.18	0.43
1:B:718:SER:O	1:B:721:VAL:HG22	2.19	0.43
1:D:645:GLN:HG2	1:D:674:GLU:HG3	2.00	0.43
1:E:695:LYS:HE2	1:E:695:LYS:HB2	1.86	0.43
1:D:670:THR:HG22	1:D:671:ASN:ND2	2.33	0.43
1:A:1226:CYS:SG	1:A:1246:LEU:HD21	2.57	0.43
1:A:1240:ILE:HG22	1:A:1241:PHE:CD1	2.54	0.43
1:C:750:TYR:CE2	1:C:775:ARG:HD3	2.53	0.43
1:D:1240:ILE:HG23	1:D:1274:TYR:CE2	2.54	0.43
1:C:914:ALA:O	1:C:941:GLN:N	2.44	0.43
1:D:879:LEU:HD11	1:D:911:ASP:HB3	2.00	0.43
1:E:818:ARG:HH21	1:E:847:ILE:HG22	1.84	0.43
1:C:900:GLU:OE2	1:C:1000:ARG:NH1	2.49	0.43
1:C:977:SER:HB2	1:C:989:CYS:SG	2.59	0.43
1:D:800:VAL:HG22	1:D:811:TYR:CD2	2.54	0.43
1:F:879:LEU:HD12	1:F:908:LEU:HA	2.01	0.43
1:C:646:SER:HB3	1:C:649:PHE:CE2	2.54	0.43
1:C:1212:GLU:HA	1:C:1215:VAL:HG12	2.01	0.43
1:E:914:ALA:O	1:E:941:GLN:N	2.44	0.43
1:B:717:ARG:O	1:B:721:VAL:HG13	2.19	0.43
1:B:976:GLU:HA	1:B:979:ARG:HH11	1.84	0.43
1:C:847:ILE:HG21	1:C:850:PRO:HB3	2.00	0.43
1:C:865:TYR:HB2	1:C:990:LEU:HD23	2.01	0.43
1:F:646:SER:HB3	1:F:649:PHE:CE2	2.53	0.43
1:D:918:TYR:O	1:D:945:ALA:HA	2.19	0.42
1:A:910:ARG:HD3	1:E:910:ARG:HG3	1.99	0.42
1:A:922:LEU:HB2	1:A:927:ARG:HG3	2.00	0.42
1:B:914:ALA:O	1:B:941:GLN:N	2.44	0.42
1:C:767:PRO:HG2	1:C:811:TYR:CE2	2.54	0.42
1:D:798:HIS:HD2	1:D:830:THR:HG23	1.83	0.42
1:A:977:SER:HB2	1:A:989:CYS:SG	2.59	0.42
1:B:816:MET:HE3	1:B:816:MET:HB3	1.93	0.42
1:F:918:TYR:CD1	1:F:949:PHE:HB2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1225:VAL:O	1:A:1229:LEU:HD23	2.19	0.42
1:B:787:LYS:HD3	1:B:787:LYS:HA	1.76	0.42
1:D:955:LYS:NZ	1:D:957:ASP:OD1	2.49	0.42
1:A:1017:GLU:O	1:A:1021:ASN:ND2	2.52	0.42
1:B:986:ILE:H	1:B:986:ILE:HD12	1.84	0.42
1:C:1017:GLU:O	1:C:1021:ASN:ND2	2.52	0.42
1:D:735:THR:HG22	1:D:736:TYR:H	1.84	0.42
1:E:869:LYS:HG3	1:E:874:VAL:HB	2.00	0.42
1:C:918:TYR:O	1:C:945:ALA:HA	2.20	0.42
1:D:720:ILE:O	1:D:724:VAL:HG23	2.19	0.42
1:D:1240:ILE:HG22	1:D:1241:PHE:HD1	1.84	0.42
1:E:840:ASP:OD1	5:E:1303:RY8:N24	2.53	0.42
1:E:963:HIS:CD2	1:E:967:PRO:HG3	2.54	0.42
1:F:818:ARG:HG3	1:F:846:LYS:HB2	2.00	0.42
1:F:1236:HIS:CD2	1:F:1238:PHE:H	2.33	0.42
1:F:1256:ASP:O	1:F:1259:VAL:HG22	2.19	0.42
1:D:895:CYS:HB2	1:D:901:CYS:SG	2.59	0.42
6:D:1304:PG4:H31	6:D:1304:PG4:H51	1.88	0.42
1:E:865:TYR:CD1	1:E:1064:ASP:HB3	2.54	0.42
1:E:893:ILE:HD13	1:E:904:MET:HE2	2.02	0.42
1:A:963:HIS:CD2	1:A:967:PRO:HG3	2.55	0.42
1:D:834:ASN:ND2	1:D:971:GLU:OE2	2.52	0.42
1:D:866:VAL:HG12	1:D:991:LEU:HB3	2.01	0.42
1:D:1219:LEU:HD12	1:D:1247:LYS:HA	2.01	0.42
1:A:967:PRO:O	1:A:1026:MET:HG2	2.20	0.41
1:A:1271:LEU:HD23	1:A:1271:LEU:HA	1.77	0.41
1:B:720:ILE:HD12	1:B:736:TYR:HB2	2.02	0.41
1:C:676:ILE:HG23	1:C:702:PRO:HG3	2.02	0.41
1:C:879:LEU:HD11	1:C:911:ASP:CB	2.50	0.41
1:E:1238:PHE:CD2	2:M:11:DG:H2'	2.55	0.41
1:F:789:LEU:O	1:F:824:VAL:HG22	2.20	0.41
1:F:986:ILE:HG23	1:F:1061:VAL:HG12	2.02	0.41
1:B:647:LEU:HA	1:B:652:THR:HG21	2.02	0.41
1:C:799:CYS:SG	1:C:951:MET:HE1	2.61	0.41
1:A:889:ASP:HB3	1:A:959:ARG:HB2	2.02	0.41
1:C:860:HIS:HB2	6:C:1304:PG4:H32	2.03	0.41
1:E:986:ILE:HD12	1:E:986:ILE:H	1.85	0.41
1:A:1245:THR:HG22	1:A:1264:ASP:O	2.21	0.41
1:B:738:THR:HG22	2:N:11:DG:OP1	2.21	0.41
1:E:918:TYR:O	1:E:945:ALA:HA	2.21	0.41
1:F:672:GLN:OE1	4:F:1302:ADP:N6	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:953:ILE:HG22	1:B:955:LYS:H	1.86	0.41
1:C:1229:LEU:O	1:C:1232:VAL:HG22	2.20	0.41
1:F:966:LEU:HA	1:F:967:PRO:HD3	1.91	0.41
1:A:1245:THR:HG21	1:A:1265:GLY:HA3	2.02	0.41
1:B:1226:CYS:SG	1:B:1246:LEU:HD21	2.60	0.41
1:C:789:LEU:HD23	1:C:824:VAL:HG21	2.02	0.41
1:C:956:PRO:HA	1:C:982:ARG:HB3	2.02	0.41
1:A:696:SER:HA	1:A:699:TYR:CE1	2.56	0.41
1:A:1275:GLY:HA2	1:A:1278:VAL:HG12	2.03	0.41
1:C:856:SER:HA	1:C:1047:GLU:OE2	2.20	0.41
1:D:1036:CYS:O	1:D:1040:GLN:HG3	2.21	0.41
1:E:736:TYR:CZ	1:E:741:LYS:HE2	2.56	0.41
1:E:971:GLU:HG3	5:E:1303:RY8:C4	2.50	0.41
1:F:827:MET:HE2	1:F:829:LEU:HD11	2.01	0.41
1:A:816:MET:O	1:A:819:GLN:HG3	2.20	0.41
1:A:842:LEU:HD22	1:A:850:PRO:HG3	2.02	0.41
1:A:918:TYR:CD1	1:A:949:PHE:HB2	2.56	0.41
1:A:1277:GLU:O	1:A:1281:VAL:HG13	2.21	0.41
1:B:918:TYR:O	1:B:945:ALA:HA	2.20	0.41
1:C:800:VAL:HG22	1:C:811:TYR:CD1	2.56	0.41
1:D:840:ASP:OD1	5:D:1303:RY8:N24	2.54	0.41
1:F:655:MET:HE1	1:F:677:ASN:OD1	2.20	0.41
1:F:1218:CYS:SG	1:F:1282:LEU:HD13	2.61	0.41
1:B:720:ILE:O	1:B:724:VAL:HG23	2.21	0.41
1:C:1036:CYS:O	1:C:1040:GLN:HG3	2.21	0.41
1:D:736:TYR:CZ	1:D:741:LYS:HE2	2.56	0.41
1:E:1253:LEU:HD23	1:E:1285:TYR:HB3	2.03	0.41
2:M:3:DT:H2''	2:M:4:DA:H8	1.86	0.41
2:N:3:DT:H2''	2:N:4:DA:C8	2.56	0.41
1:A:765:VAL:HG21	1:A:770:ILE:HD11	2.03	0.40
1:C:699:TYR:CD1	1:C:829:LEU:HD11	2.55	0.40
1:A:714:SER:HB3	1:A:720:ILE:HD11	2.02	0.40
1:A:958:VAL:HB	1:A:980:ALA:HA	2.04	0.40
1:B:663:PHE:HB3	1:B:665:LEU:HD13	2.03	0.40
1:C:862:LEU:O	1:C:1038:ARG:NH2	2.54	0.40
1:E:688:LEU:O	1:E:854:SER:HA	2.21	0.40
1:E:747:THR:OG1	1:E:775:ARG:NH2	2.54	0.40
1:F:1255:SER:HB3	1:F:1283:GLN:HA	2.04	0.40
1:A:1057:LYS:O	1:B:786:ARG:HD3	2.21	0.40
1:C:878:CYS:O	1:C:882:ILE:HG13	2.22	0.40
1:C:969:SER:HB2	1:C:971:GLU:OE1	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:1017:GLU:O	1:E:1021:ASN:ND2	2.54	0.40
1:A:919:HIS:H	1:A:922:LEU:HD23	1.86	0.40
1:B:1055:CYS:HB2	1:B:1066:CYS:SG	2.61	0.40
1:B:1214:MET:HA	1:B:1217:LYS:NZ	2.37	0.40
1:F:698:CYS:C	1:F:827:MET:HE1	2.47	0.40
1:C:720:ILE:O	1:C:724:VAL:HG23	2.22	0.40
1:D:758:PRO:HG2	1:D:788:LEU:HD13	2.04	0.40
1:D:816:MET:HE3	1:D:816:MET:HB3	1.94	0.40
1:F:963:HIS:HB2	1:F:991:LEU:HD13	2.03	0.40
1:F:971:GLU:H	1:F:971:GLU:HG3	1.41	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	520/563 (92%)	498 (96%)	20 (4%)	2 (0%)	30	53
1	B	509/563 (90%)	488 (96%)	19 (4%)	2 (0%)	30	53
1	C	510/563 (91%)	490 (96%)	19 (4%)	1 (0%)	43	66
1	D	507/563 (90%)	488 (96%)	17 (3%)	2 (0%)	30	53
1	E	510/563 (91%)	492 (96%)	17 (3%)	1 (0%)	43	66
1	F	513/563 (91%)	488 (95%)	20 (4%)	5 (1%)	12	32
All	All	3069/3378 (91%)	2944 (96%)	112 (4%)	13 (0%)	30	53

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	759	ILE
1	B	759	ILE

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Mol	Chain	Res	Type
1	D	759	ILE
1	F	807	PHE
1	F	1280	SER
1	A	873	LYS
1	C	873	LYS
1	D	873	LYS
1	F	873	LYS
1	B	873	LYS
1	E	873	LYS
1	F	806	ASP
1	F	1255	SER

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	421/499 (84%)	412 (98%)	9 (2%)	47	69
1	B	398/499 (80%)	387 (97%)	11 (3%)	38	62
1	C	402/499 (81%)	392 (98%)	10 (2%)	42	66
1	D	374/499 (75%)	364 (97%)	10 (3%)	39	64
1	E	381/499 (76%)	366 (96%)	15 (4%)	28	54
1	F	360/499 (72%)	346 (96%)	14 (4%)	28	54
All	All	2336/2994 (78%)	2267 (97%)	69 (3%)	36	60

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	643	ARG
1	A	720	ILE
1	A	738	THR
1	A	815	ASN
1	A	816	MET
1	A	824	VAL
1	A	826	VAL

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Mol	Chain	Res	Type
1	A	994	THR
1	A	1243	THR
1	B	652	THR
1	B	665	LEU
1	B	728	THR
1	B	816	MET
1	B	824	VAL
1	B	830	THR
1	B	832	THR
1	B	852	VAL
1	B	937	GLN
1	B	951	MET
1	B	1215	VAL
1	C	652	THR
1	C	665	LEU
1	C	738	THR
1	C	780	LEU
1	C	814	MET
1	C	816	MET
1	C	830	THR
1	C	843	THR
1	C	1235	VAL
1	C	1261	LEU
1	D	735	THR
1	D	738	THR
1	D	753	LEU
1	D	783	LEU
1	D	923	SER
1	D	942	VAL
1	D	951	MET
1	D	1001	LEU
1	D	1235	VAL
1	D	1243	THR
1	E	652	THR
1	E	747	THR
1	E	786	ARG
1	E	811	TYR
1	E	814	MET
1	E	817	LEU
1	E	826	VAL
1	E	830	THR
1	E	953	ILE

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Mol	Chain	Res	Type
1	E	1209	SER
1	E	1215	VAL
1	E	1224	GLU
1	E	1232	VAL
1	E	1235	VAL
1	E	1239	ASN
1	F	681	LEU
1	F	776	LEU
1	F	824	VAL
1	F	830	THR
1	F	832	THR
1	F	903	THR
1	F	922	LEU
1	F	942	VAL
1	F	971	GLU
1	F	1017	GLU
1	F	1243	THR
1	F	1253	LEU
1	F	1259	VAL
1	F	1260	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	641	HIS
1	A	671	ASN
1	A	752	GLN
1	A	798	HIS
1	A	802	GLN
1	A	885	HIS
1	A	1013	HIS
1	A	1021	ASN
1	A	1239	ASN
1	B	651	HIS
1	B	671	ASN
1	B	798	HIS
1	B	1239	ASN
1	C	672	GLN
1	C	752	GLN
1	C	802	GLN
1	C	996	HIS
1	C	1013	HIS

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Mol	Chain	Res	Type
1	C	1021	ASN
1	C	1048	ASN
1	C	1065	ASN
1	D	666	HIS
1	D	671	ASN
1	D	723	GLN
1	D	752	GLN
1	D	774	ASN
1	D	798	HIS
1	D	909	GLN
1	D	931	GLN
1	D	1021	ASN
1	E	748	ASN
1	E	752	GLN
1	E	774	ASN
1	E	802	GLN
1	E	858	ASN
1	E	996	HIS
1	E	1021	ASN
1	E	1058	HIS
1	F	667	ASN
1	F	671	ASN
1	F	798	HIS
1	F	802	GLN
1	F	885	HIS
1	F	1019	HIS
1	F	1021	ASN
1	F	1236	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 31 ligands modelled in this entry, 12 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
4	ADP	D	1302	-	28,29,29	1.42	4 (14%)	43,45,45	1.82	9 (20%)
5	RY8	A	1303	-	29,29,29	0.15	0	42,43,43	0.32	0
4	ADP	E	1302	-	28,29,29	1.41	4 (14%)	43,45,45	1.83	9 (20%)
6	PG4	B	1304	-	12,12,12	0.11	0	11,11,11	0.12	0
6	PG4	E	1304	-	12,12,12	0.11	0	11,11,11	0.11	0
6	PG4	A	1304	-	12,12,12	0.12	0	11,11,11	0.11	0
6	PG4	C	1304	-	12,12,12	0.12	0	11,11,11	0.10	0
5	RY8	E	1303	-	29,29,29	0.10	0	42,43,43	0.25	0
6	PG4	D	1304	-	12,12,12	0.13	0	11,11,11	0.09	0
4	ADP	F	1302	-	28,29,29	1.40	4 (14%)	43,45,45	1.82	9 (20%)
4	ADP	B	1302	-	28,29,29	1.41	4 (14%)	43,45,45	1.80	9 (20%)
5	RY8	C	1303	-	29,29,29	0.10	0	42,43,43	0.24	0
6	PG4	C	1305	-	12,12,12	0.11	0	11,11,11	0.11	0
6	PG4	A	1305	-	12,12,12	0.12	0	11,11,11	0.11	0
4	ADP	A	1302	-	28,29,29	1.40	4 (14%)	43,45,45	1.81	9 (20%)
5	RY8	D	1303	-	29,29,29	0.12	0	42,43,43	0.27	0
5	RY8	B	1303	-	29,29,29	0.11	0	42,43,43	0.26	0
4	ADP	C	1302	-	28,29,29	1.41	4 (14%)	43,45,45	1.80	9 (20%)
5	RY8	F	1303	-	29,29,29	0.11	0	42,43,43	0.25	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ADP	D	1302	-	-	4/16/32/32	0/3/3/3
5	RY8	A	1303	-	-	2/18/18/18	0/3/3/3
4	ADP	E	1302	-	-	1/16/32/32	0/3/3/3
6	PG4	B	1304	-	-	4/10/10/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PG4	E	1304	-	-	5/10/10/10	-
6	PG4	A	1304	-	-	2/10/10/10	-
6	PG4	C	1304	-	-	4/10/10/10	-
5	RY8	E	1303	-	-	6/18/18/18	0/3/3/3
6	PG4	D	1304	-	-	4/10/10/10	-
4	ADP	F	1302	-	-	1/16/32/32	0/3/3/3
4	ADP	B	1302	-	-	5/16/32/32	0/3/3/3
5	RY8	C	1303	-	-	2/18/18/18	0/3/3/3
6	PG4	C	1305	-	-	3/10/10/10	-
6	PG4	A	1305	-	-	7/10/10/10	-
4	ADP	A	1302	-	-	5/16/32/32	0/3/3/3
5	RY8	D	1303	-	-	6/18/18/18	0/3/3/3
5	RY8	B	1303	-	-	6/18/18/18	0/3/3/3
4	ADP	C	1302	-	-	3/16/32/32	0/3/3/3
5	RY8	F	1303	-	-	5/18/18/18	0/3/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1302	ADP	C5-C4	4.73	1.47	1.39
4	D	1302	ADP	C5-C4	4.72	1.47	1.39
4	A	1302	ADP	C5-C4	4.69	1.47	1.39
4	C	1302	ADP	C5-C4	4.69	1.47	1.39
4	F	1302	ADP	C5-C4	4.69	1.47	1.39
4	B	1302	ADP	C5-C4	4.68	1.47	1.39
4	D	1302	ADP	C5-C6	2.77	1.48	1.41
4	F	1302	ADP	C5-C6	2.76	1.48	1.41
4	E	1302	ADP	C5-C6	2.75	1.48	1.41
4	C	1302	ADP	C5-C6	2.75	1.48	1.41
4	B	1302	ADP	C5-C6	2.74	1.48	1.41
4	A	1302	ADP	C5-C6	2.74	1.48	1.41
4	D	1302	ADP	C8-N7	2.44	1.36	1.31
4	C	1302	ADP	C8-N7	2.42	1.36	1.31
4	F	1302	ADP	C8-N7	2.41	1.36	1.31
4	E	1302	ADP	C8-N7	2.39	1.36	1.31
4	B	1302	ADP	C8-N7	2.39	1.36	1.31
4	A	1302	ADP	C8-N7	2.38	1.36	1.31
4	B	1302	ADP	C5-N7	-2.25	1.35	1.39
4	D	1302	ADP	C5-N7	-2.23	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1302	ADP	C5-N7	-2.22	1.35	1.39
4	C	1302	ADP	C5-N7	-2.19	1.35	1.39
4	A	1302	ADP	C5-N7	-2.16	1.35	1.39
4	F	1302	ADP	C5-N7	-2.13	1.35	1.39

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	1302	ADP	C5-C4-N3	-5.79	118.75	126.72
4	F	1302	ADP	C5-C4-N3	-5.68	118.90	126.72
4	B	1302	ADP	C5-C4-N3	-5.63	118.96	126.72
4	A	1302	ADP	C5-C4-N3	-5.63	118.96	126.72
4	D	1302	ADP	C5-C4-N3	-5.62	118.97	126.72
4	C	1302	ADP	C5-C4-N3	-5.57	119.05	126.72
4	E	1302	ADP	N3-C4-N9	4.51	134.84	127.17
4	B	1302	ADP	N3-C4-N9	4.49	134.81	127.17
4	F	1302	ADP	N3-C4-N9	4.47	134.77	127.17
4	D	1302	ADP	N3-C4-N9	4.40	134.66	127.17
4	A	1302	ADP	N3-C4-N9	4.40	134.65	127.17
4	C	1302	ADP	N3-C4-N9	4.34	134.55	127.17
4	E	1302	ADP	C2-N3-C4	3.73	120.93	111.83
4	F	1302	ADP	C2-N3-C4	3.68	120.82	111.83
4	C	1302	ADP	C2-N3-C4	3.66	120.77	111.83
4	B	1302	ADP	C2-N3-C4	3.65	120.75	111.83
4	D	1302	ADP	C2-N3-C4	3.65	120.75	111.83
4	A	1302	ADP	C2-N3-C4	3.65	120.75	111.83
4	E	1302	ADP	C4-C5-N7	-3.57	106.50	110.58
4	C	1302	ADP	C4-C5-N7	-3.56	106.51	110.58
4	A	1302	ADP	C4-C5-N7	-3.55	106.53	110.58
4	D	1302	ADP	C4-C5-N7	-3.50	106.58	110.58
4	F	1302	ADP	C4-C5-N7	-3.46	106.62	110.58
4	B	1302	ADP	C4-C5-N7	-3.39	106.70	110.58
4	F	1302	ADP	N3-C2-N1	-3.31	123.57	128.58
4	A	1302	ADP	N3-C2-N1	-3.30	123.59	128.58
4	C	1302	ADP	N3-C2-N1	-3.29	123.60	128.58
4	E	1302	ADP	N3-C2-N1	-3.28	123.62	128.58
4	B	1302	ADP	N3-C2-N1	-3.27	123.64	128.58
4	D	1302	ADP	N3-C2-N1	-3.26	123.64	128.58
4	B	1302	ADP	C4-N9-C8	2.77	108.64	105.74
4	C	1302	ADP	C4-N9-C8	2.75	108.63	105.74
4	A	1302	ADP	C4-N9-C8	2.75	108.62	105.74
4	D	1302	ADP	C4-N9-C8	2.72	108.60	105.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	1302	ADP	C4-N9-C8	2.67	108.55	105.74
4	E	1302	ADP	C4-N9-C8	2.64	108.51	105.74
4	E	1302	ADP	C5-N7-C8	2.60	107.54	103.45
4	C	1302	ADP	C5-N7-C8	2.60	107.53	103.45
4	A	1302	ADP	C5-N7-C8	2.57	107.49	103.45
4	F	1302	ADP	C5-N7-C8	2.53	107.43	103.45
4	D	1302	ADP	C5-N7-C8	2.53	107.43	103.45
4	B	1302	ADP	C5-N7-C8	2.52	107.41	103.45
4	C	1302	ADP	C6-C5-N7	2.30	136.52	132.09
4	E	1302	ADP	C6-C5-N7	2.23	136.38	132.09
4	A	1302	ADP	C6-C5-N7	2.21	136.35	132.09
4	D	1302	ADP	C6-C5-N7	2.20	136.33	132.09
4	F	1302	ADP	C6-C5-N7	2.15	136.24	132.09
4	B	1302	ADP	C6-C5-N7	2.13	136.20	132.09
4	C	1302	ADP	N9-C8-N7	-2.12	110.92	113.94
4	A	1302	ADP	N9-C8-N7	-2.09	110.98	113.94
4	B	1302	ADP	N9-C8-N7	-2.06	111.01	113.94
4	D	1302	ADP	N9-C8-N7	-2.05	111.02	113.94
4	E	1302	ADP	N9-C8-N7	-2.04	111.04	113.94
4	F	1302	ADP	N9-C8-N7	-2.02	111.07	113.94

There are no chirality outliers.

All (75) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1302	ADP	C5'-O5'-PA-O1A
4	A	1302	ADP	C5'-O5'-PA-O3A
4	B	1302	ADP	PA-O3A-PB-O2B
4	B	1302	ADP	PA-O3A-PB-O3B
4	C	1302	ADP	PA-O3A-PB-O2B
4	D	1302	ADP	PA-O3A-PB-O2B
4	D	1302	ADP	C5'-O5'-PA-O1A
5	F	1303	RY8	C10-C11-C4-N2
4	D	1302	ADP	O4'-C4'-C5'-O5'
6	A	1305	PG4	O3-C5-C6-O4
5	B	1303	RY8	C10-C11-C4-N2
5	F	1303	RY8	C9-C11-C4-N2
5	B	1303	RY8	C9-C11-C4-N2
4	A	1302	ADP	O4'-C4'-C5'-O5'
4	A	1302	ADP	C3'-C4'-C5'-O5'
4	B	1302	ADP	O4'-C4'-C5'-O5'
4	C	1302	ADP	O4'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
5	F	1303	RY8	C10-C11-C4-C5
5	F	1303	RY8	C9-C11-C4-C5
4	D	1302	ADP	C3'-C4'-C5'-O5'
5	B	1303	RY8	C10-C11-C4-C5
5	B	1303	RY8	C9-C11-C4-C5
4	B	1302	ADP	C3'-C4'-C5'-O5'
4	C	1302	ADP	C3'-C4'-C5'-O5'
5	C	1303	RY8	C20-C15-N14-C13
5	A	1303	RY8	C20-C15-N14-C13
5	E	1303	RY8	C19-C15-N14-C13
5	B	1303	RY8	C20-C15-N14-C13
5	B	1303	RY8	C19-C15-N14-C13
5	E	1303	RY8	C20-C15-N14-C13
6	B	1304	PG4	C5-C6-O4-C7
6	E	1304	PG4	C8-C7-O4-C6
6	C	1305	PG4	C4-C3-O2-C2
6	B	1304	PG4	C4-C3-O2-C2
6	C	1304	PG4	C4-C3-O2-C2
6	C	1305	PG4	C1-C2-O2-C3
6	E	1304	PG4	C5-C6-O4-C7
5	C	1303	RY8	C19-C15-N14-C13
6	C	1304	PG4	C8-C7-O4-C6
6	A	1305	PG4	C5-C6-O4-C7
6	E	1304	PG4	C6-C5-O3-C4
6	D	1304	PG4	C8-C7-O4-C6
5	D	1303	RY8	C20-C15-N14-C13
6	B	1304	PG4	C6-C5-O3-C4
5	A	1303	RY8	C19-C15-N14-C13
4	A	1302	ADP	C5'-O5'-PA-O2A
4	B	1302	ADP	C5'-O5'-PA-O1A
6	D	1304	PG4	C5-C6-O4-C7
5	D	1303	RY8	C10-C11-C4-N2
6	A	1305	PG4	C1-C2-O2-C3
6	D	1304	PG4	C3-C4-O3-C5
6	C	1305	PG4	C5-C6-O4-C7
5	D	1303	RY8	C19-C15-N14-C13
6	A	1305	PG4	C4-C3-O2-C2
5	D	1303	RY8	C10-C11-C4-C5
5	E	1303	RY8	C10-C11-C4-C5
6	B	1304	PG4	O2-C3-C4-O3
5	D	1303	RY8	C9-C11-C4-N2
6	A	1305	PG4	C3-C4-O3-C5

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Mol	Chain	Res	Type	Atoms
4	F	1302	ADP	O4'-C4'-C5'-O5'
5	E	1303	RY8	C9-C11-C4-N2
6	D	1304	PG4	C4-C3-O2-C2
6	A	1305	PG4	C6-C5-O3-C4
5	E	1303	RY8	C10-C11-C4-N2
6	E	1304	PG4	C3-C4-O3-C5
5	D	1303	RY8	C9-C11-C4-C5
5	E	1303	RY8	C9-C11-C4-C5
6	A	1304	PG4	O2-C3-C4-O3
6	E	1304	PG4	C1-C2-O2-C3
6	C	1304	PG4	C6-C5-O3-C4
4	E	1302	ADP	C3'-C4'-C5'-O5'
6	A	1304	PG4	C1-C2-O2-C3
6	C	1304	PG4	O2-C3-C4-O3
6	A	1305	PG4	O2-C3-C4-O3
5	F	1303	RY8	C20-C15-N14-C13

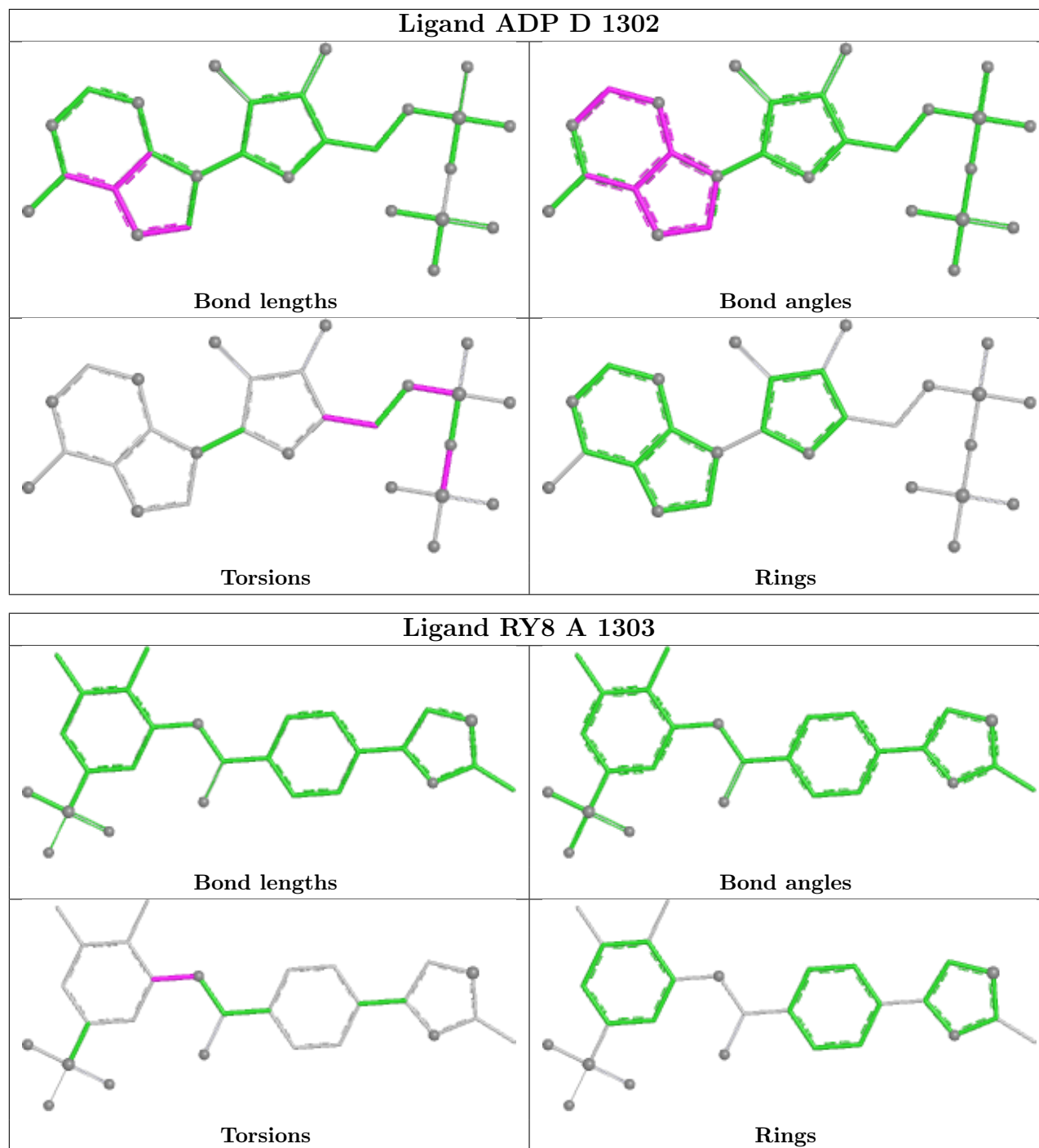
There are no ring outliers.

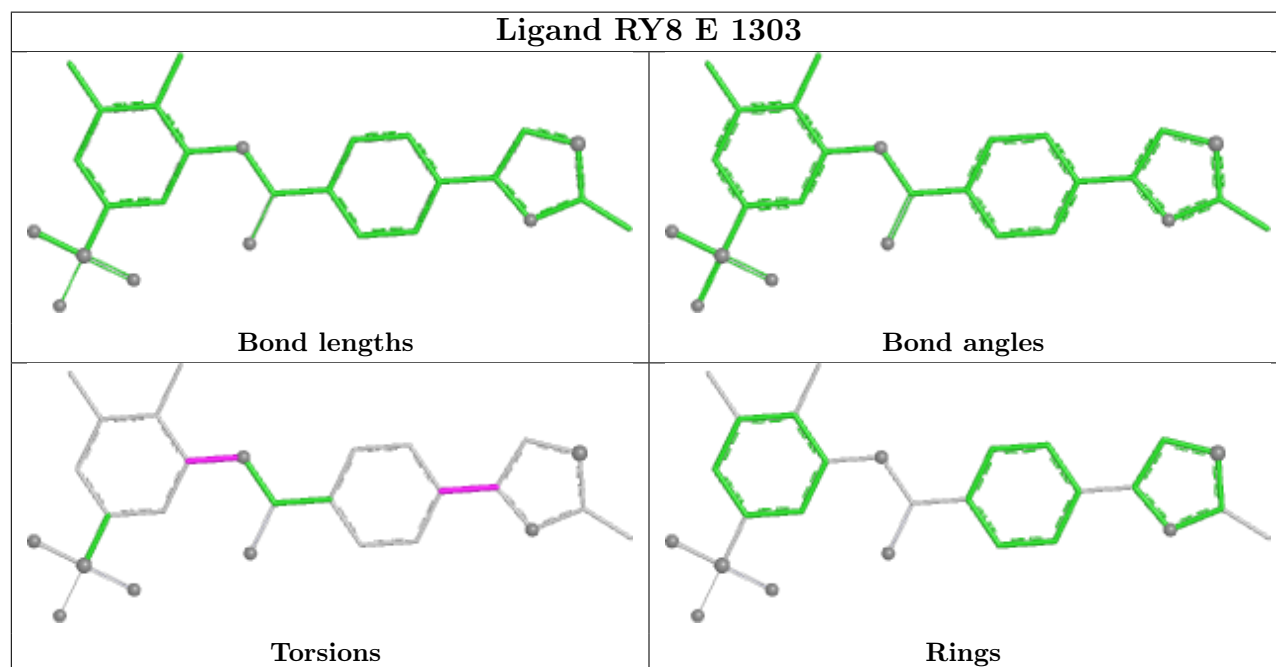
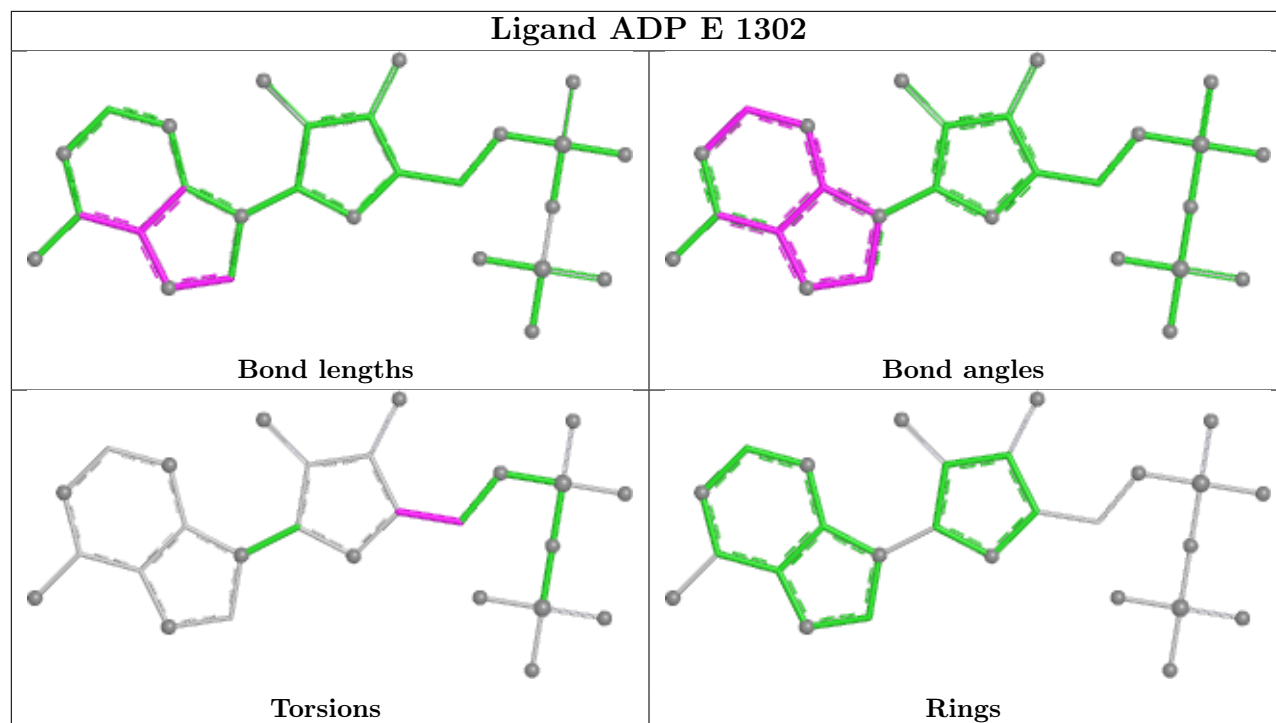
11 monomers are involved in 12 short contacts:

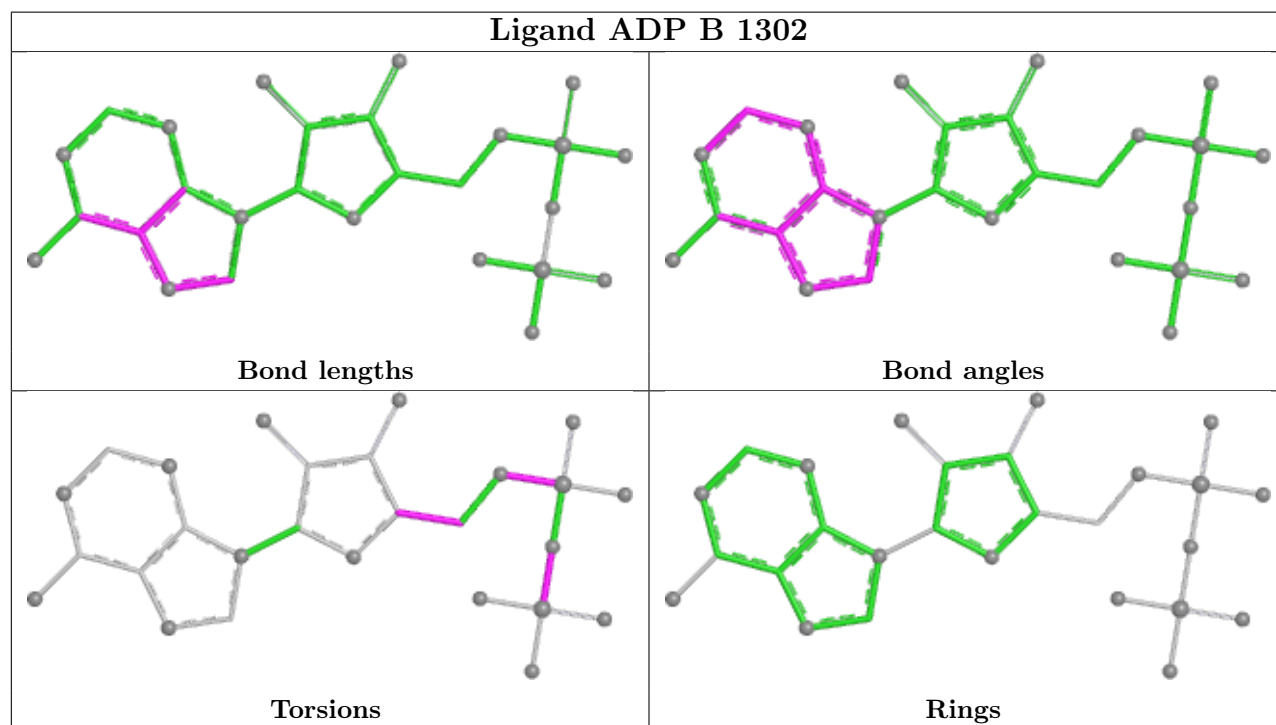
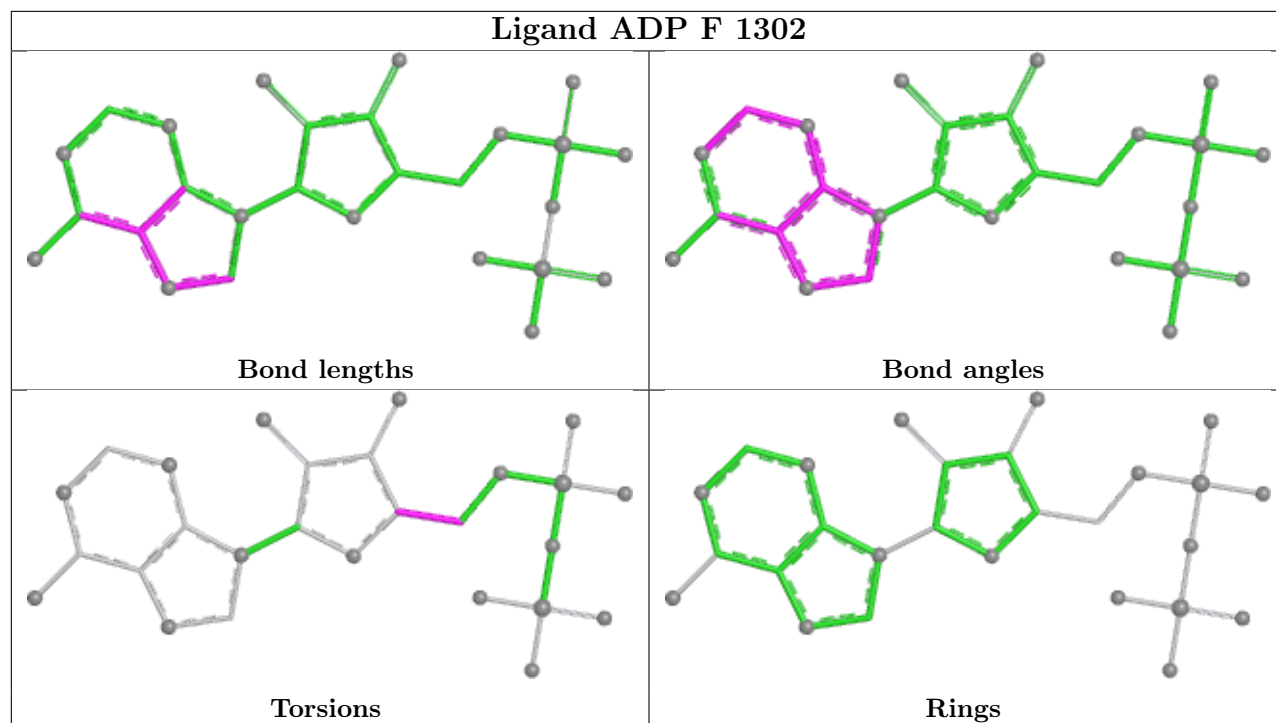
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1304	PG4	1	0
6	E	1304	PG4	1	0
6	A	1304	PG4	1	0
6	C	1304	PG4	1	0
5	E	1303	RY8	2	0
6	D	1304	PG4	1	0
4	F	1302	ADP	1	0
6	C	1305	PG4	1	0
6	A	1305	PG4	1	0
5	D	1303	RY8	1	0
4	C	1302	ADP	1	0

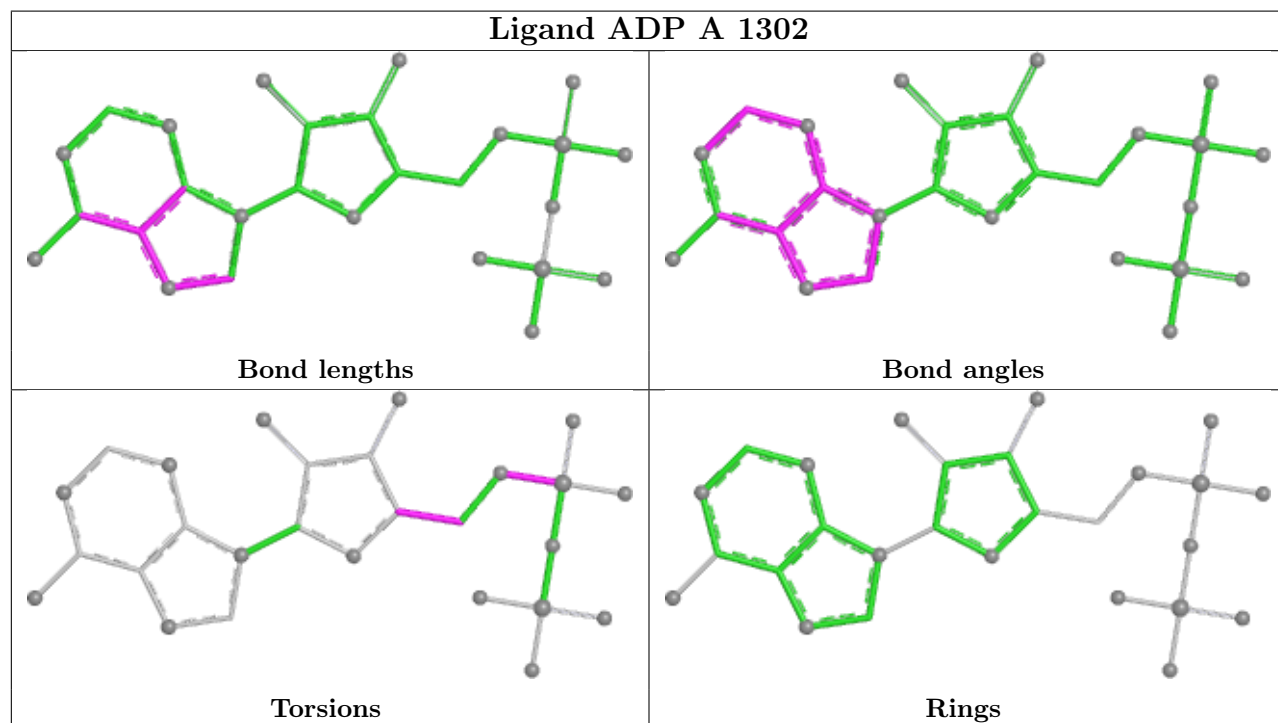
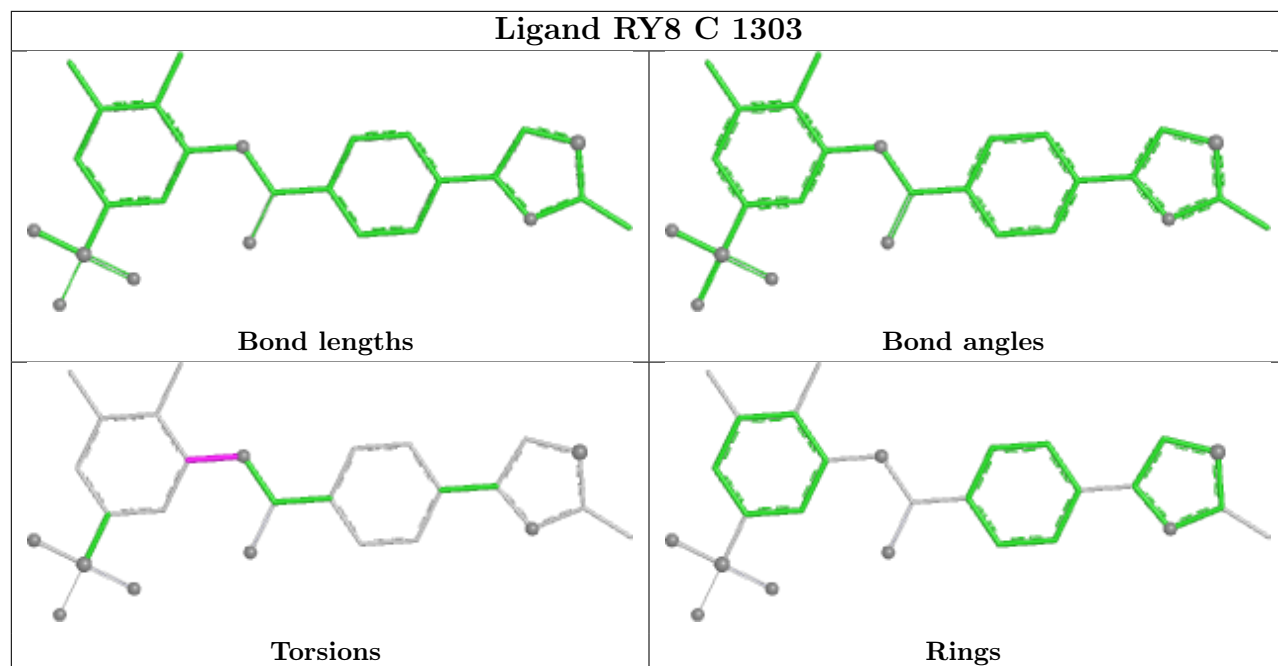
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

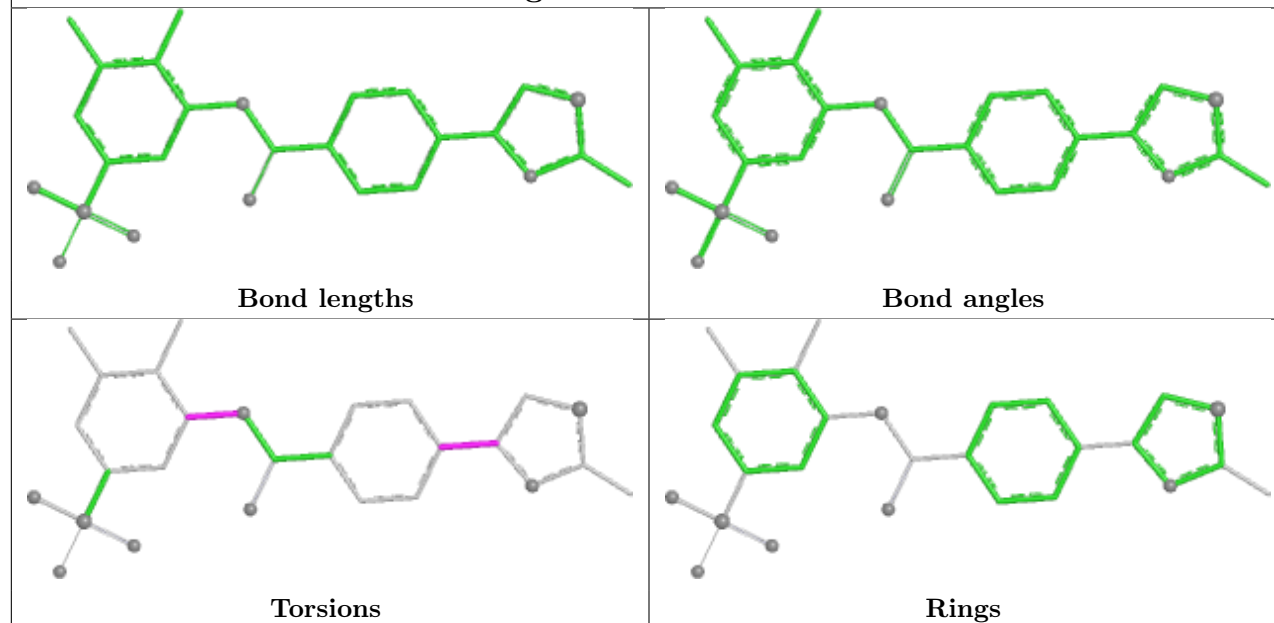




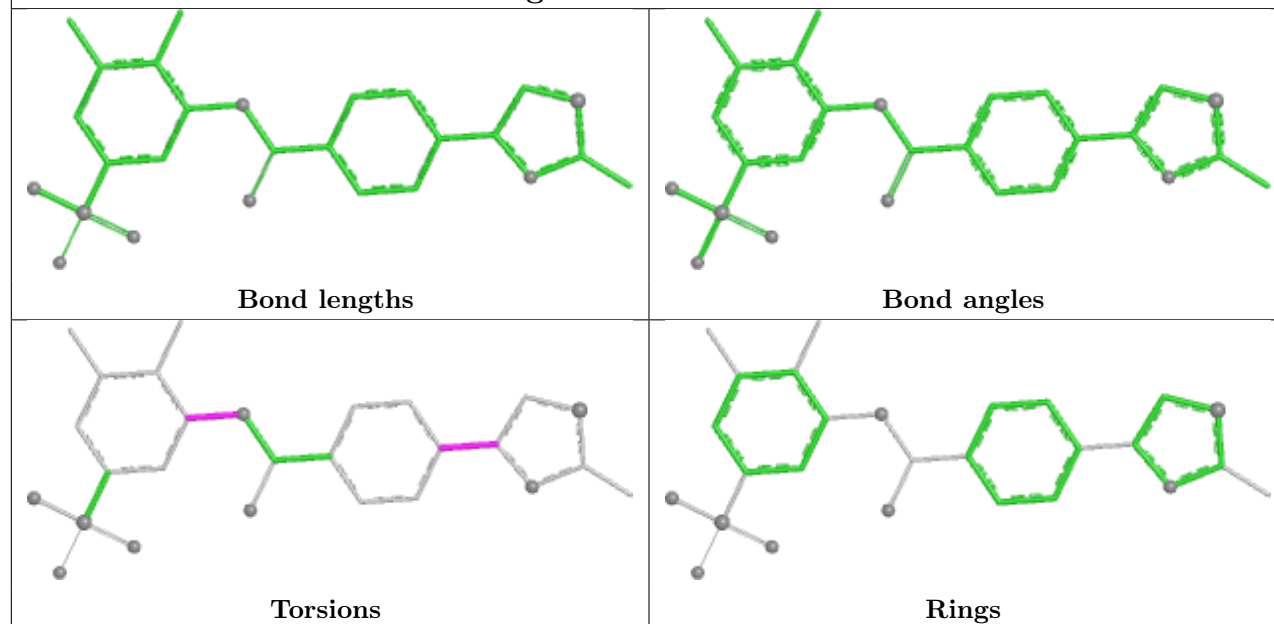


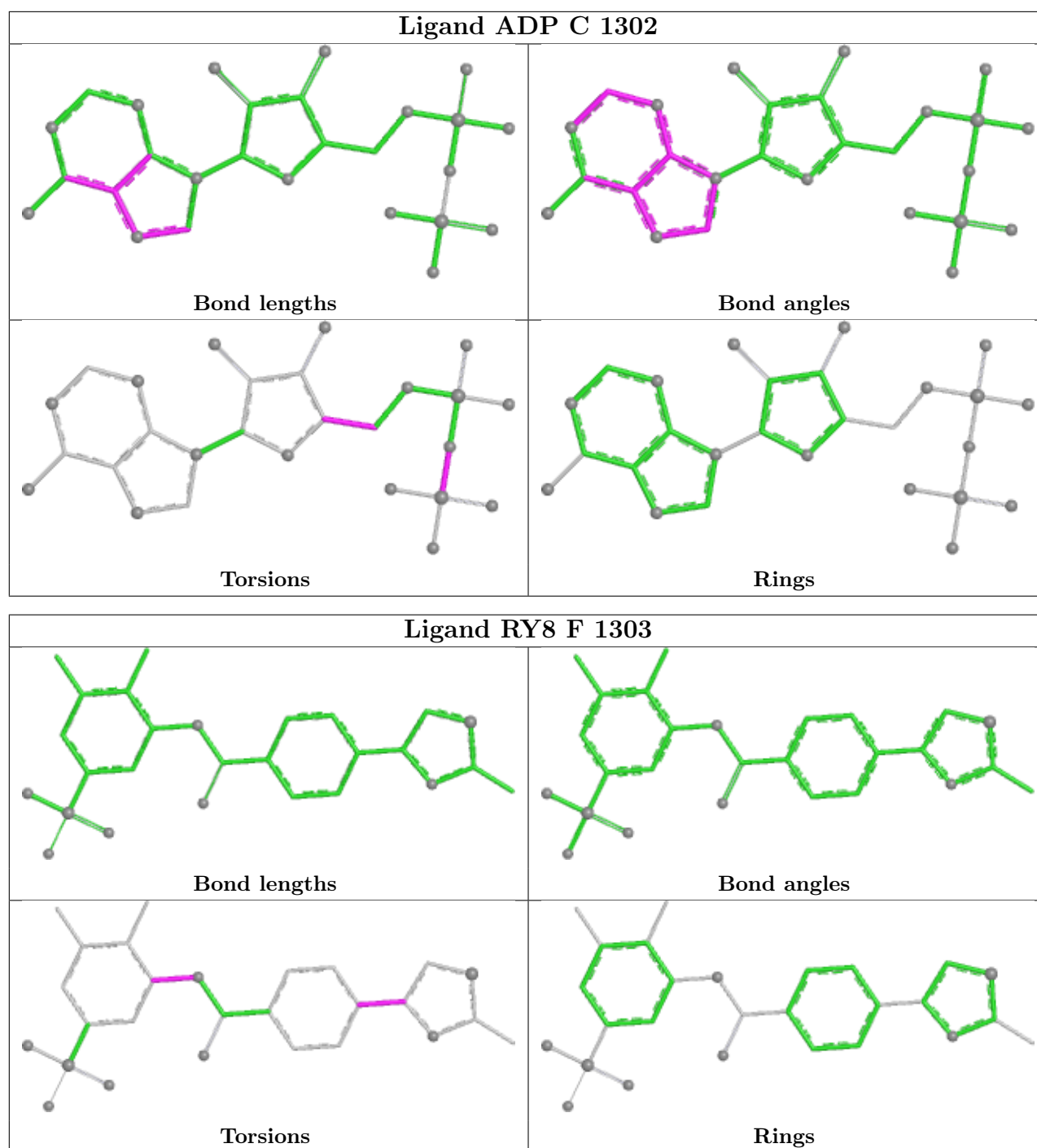


Ligand RY8 D 1303



Ligand RY8 B 1303





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	522/563 (92%)	0.27	8 (1%) 72 65	43, 71, 125, 194	3 (0%)
1	B	513/563 (91%)	0.26	11 (2%) 63 55	51, 79, 143, 197	1 (0%)
1	C	513/563 (91%)	0.37	7 (1%) 73 67	38, 84, 143, 194	2 (0%)
1	D	510/563 (90%)	0.54	25 (4%) 35 29	48, 94, 156, 213	3 (0%)
1	E	513/563 (91%)	0.49	21 (4%) 41 34	52, 94, 147, 193	3 (0%)
1	F	516/563 (91%)	0.66	31 (6%) 27 23	66, 101, 153, 229	3 (0%)
2	M	13/15 (86%)	-0.05	0 100 100	63, 84, 121, 129	0
2	N	13/15 (86%)	-0.07	0 100 100	60, 80, 143, 150	0
2	O	13/15 (86%)	0.07	0 100 100	70, 101, 145, 158	0
2	P	13/15 (86%)	-0.06	0 100 100	68, 93, 124, 153	0
2	Q	13/15 (86%)	0.08	0 100 100	78, 100, 137, 164	0
2	R	13/15 (86%)	0.18	0 100 100	105, 121, 157, 163	0
All	All	3165/3468 (91%)	0.42	103 (3%) 49 42	38, 87, 146, 229	15 (0%)

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	1278	VAL	4.1
1	C	878	CYS	4.1
1	F	878	CYS	3.7
1	D	704	CYS	3.6
1	E	995	TYR	3.4
1	E	694	GLY	3.4
1	B	758	PRO	3.3
1	F	1015	THR	3.3
1	D	990	LEU	3.3
1	B	1209	SER	3.3
1	C	638	ASN	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	1237	TYR	3.1
1	D	1209	SER	3.1
1	F	952	GLY	3.0
1	E	727	LEU	3.0
1	F	943	ILE	3.0
1	D	819	GLN	3.0
1	A	1070	LYS	2.9
1	D	951	MET	2.9
1	F	701	LEU	2.9
1	F	758	PRO	2.7
1	E	686	PHE	2.7
1	E	782	ASN	2.7
1	D	1260	LEU	2.7
1	D	641	HIS	2.6
1	F	882	ILE	2.6
1	F	1288	TRP	2.6
1	D	847	ILE	2.6
1	E	743	ASP	2.6
1	E	697	LEU	2.6
1	D	705	VAL	2.6
1	F	989	CYS	2.6
1	B	785	GLU	2.6
1	C	997	ASP	2.5
1	F	1020	PHE	2.5
1	E	641	HIS	2.5
1	B	743	ASP	2.4
1	E	990	LEU	2.4
1	D	1008	GLU	2.4
1	D	845	LEU	2.4
1	B	939	GLY	2.4
1	C	961	VAL	2.4
1	E	1223	THR	2.4
1	E	655	MET	2.4
1	D	1035	GLU	2.4
1	F	992	PHE	2.4
1	D	1043	ALA	2.4
1	A	1263	ILE	2.4
1	A	952	GLY	2.4
1	F	915	ALA	2.3
1	B	951	MET	2.3
1	F	804	GLY	2.3
1	D	1030	CYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	1204	LEU	2.3
1	A	980	ALA	2.3
1	C	1282	LEU	2.3
1	C	1279	ILE	2.3
1	D	996	HIS	2.3
1	E	1255	SER	2.3
1	F	1286	SER	2.3
1	D	992	PHE	2.3
1	F	996	HIS	2.3
1	F	969	SER	2.3
1	B	1011	GLY	2.3
1	C	641	HIS	2.2
1	D	648	SER	2.2
1	F	1281	VAL	2.2
1	E	1234	GLY	2.2
1	A	1270	LYS	2.2
1	F	770	ILE	2.2
1	D	997	ASP	2.2
1	E	857	PHE	2.2
1	B	1211	ARG	2.2
1	D	1266	VAL	2.2
1	F	806	ASP	2.2
1	E	994	THR	2.2
1	D	1274	TYR	2.2
1	D	977	SER	2.2
1	F	971	GLU	2.2
1	E	695	LYS	2.1
1	F	1222	LEU	2.1
1	A	1274	TYR	2.1
1	D	799	CYS	2.1
1	E	732	ILE	2.1
1	F	1201	SER	2.1
1	E	874	VAL	2.1
1	D	1291	PRO	2.1
1	D	994	THR	2.1
1	E	1010	ASP	2.1
1	F	757	ASP	2.1
1	A	1006	MET	2.1
1	B	953	ILE	2.1
1	B	641	HIS	2.0
1	F	967	PRO	2.0
1	D	792	PHE	2.0

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Mol	Chain	Res	Type	RSRZ
1	F	1035	GLU	2.0
1	E	689	MET	2.0
1	F	937	GLN	2.0
1	F	698	CYS	2.0
1	E	848	LEU	2.0
1	F	990	LEU	2.0
1	B	1007	MET	2.0
1	F	1237	TYR	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.

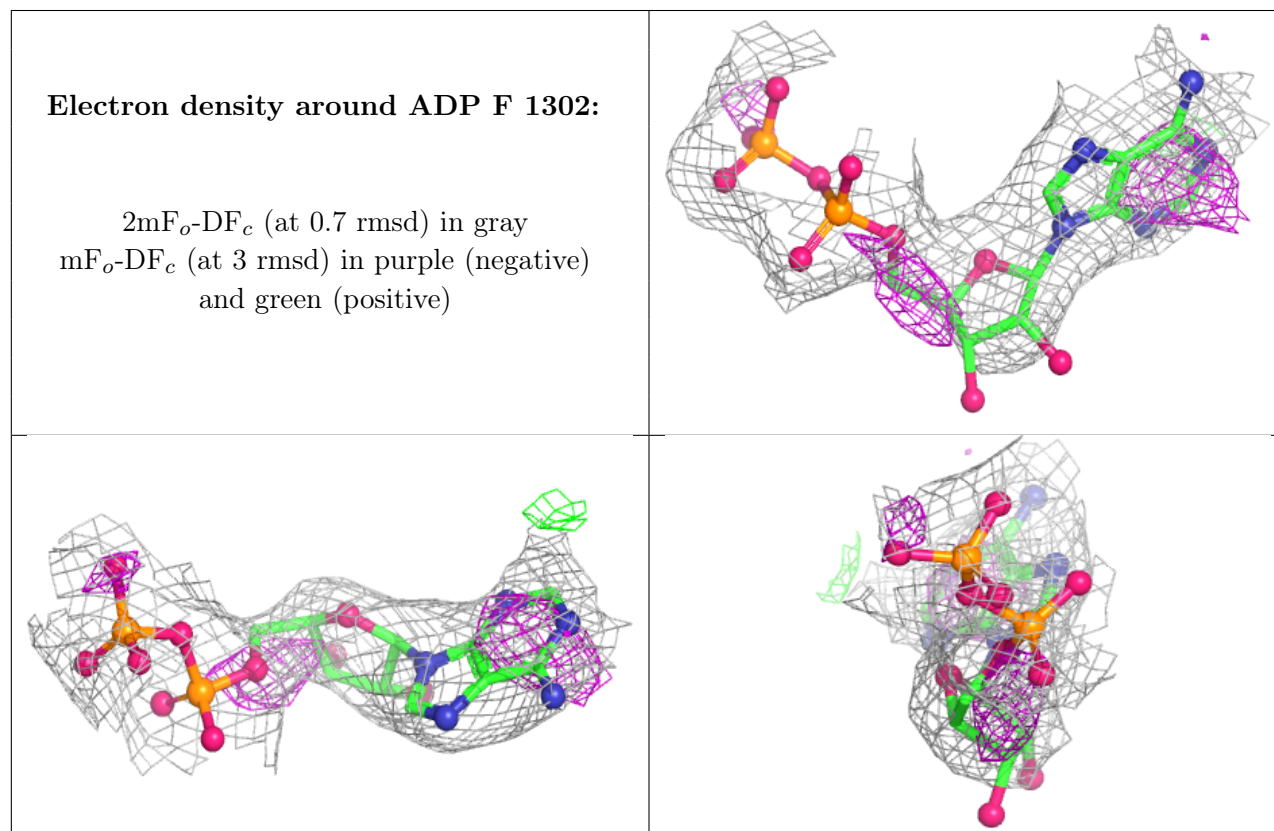
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	MG	F	1304	1/1	0.66	0.31	94,94,94,94	0
7	MG	N	101	1/1	0.71	0.24	69,69,69,69	0
6	PG4	A	1305	13/13	0.77	0.16	84,93,99,102	0
6	PG4	C	1305	13/13	0.77	0.16	83,98,103,104	0
6	PG4	B	1304	13/13	0.79	0.14	79,84,94,95	0
6	PG4	A	1304	13/13	0.79	0.14	74,77,84,89	0
4	ADP	F	1302	27/27	0.81	0.11	70,86,96,110	0
6	PG4	D	1304	13/13	0.84	0.15	86,92,100,101	0
6	PG4	E	1304	13/13	0.84	0.15	83,90,93,97	0
6	PG4	C	1304	13/13	0.85	0.12	85,92,100,101	0
7	MG	Q	101	1/1	0.85	0.17	79,79,79,79	0
4	ADP	E	1302	27/27	0.87	0.11	88,100,111,113	0
7	MG	M	100	1/1	0.87	0.25	69,69,69,69	0
7	MG	P	101	1/1	0.88	0.14	67,67,67,67	0
4	ADP	D	1302	27/27	0.88	0.10	76,99,108,113	0

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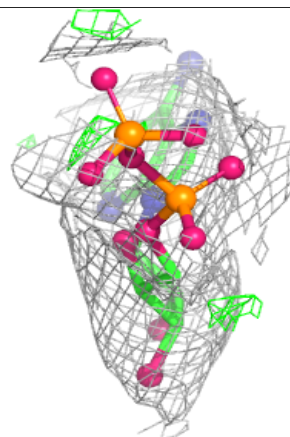
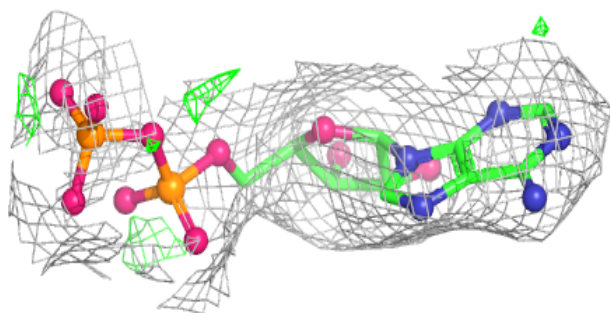
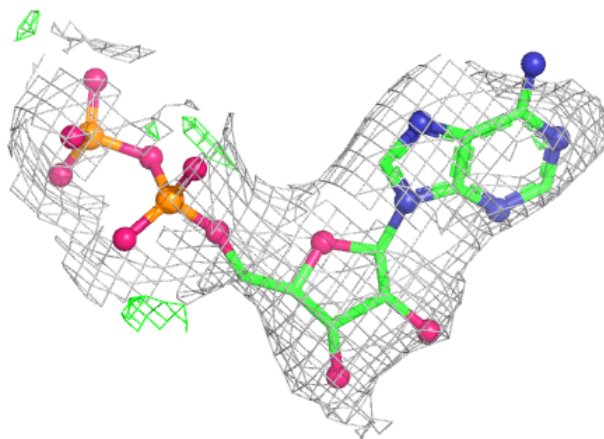
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	MG	Q	102	1/1	0.88	0.26	77,77,77,77	0
5	RY8	F	1303	27/27	0.89	0.13	78,83,89,94	0
5	RY8	D	1303	27/27	0.91	0.12	68,85,99,108	0
4	ADP	C	1302	27/27	0.91	0.09	58,73,82,85	0
5	RY8	C	1303	27/27	0.91	0.13	57,71,80,96	0
5	RY8	E	1303	27/27	0.92	0.11	63,76,79,88	0
4	ADP	A	1302	27/27	0.93	0.08	61,69,81,83	0
4	ADP	B	1302	27/27	0.94	0.07	62,74,80,82	0
5	RY8	B	1303	27/27	0.95	0.09	45,59,66,82	0
5	RY8	A	1303	27/27	0.95	0.09	43,51,61,82	0
3	ZN	E	1301	1/1	0.97	0.06	121,121,121,121	0
3	ZN	F	1301	1/1	0.98	0.03	103,103,103,103	0
3	ZN	C	1301	1/1	0.98	0.05	95,95,95,95	0
3	ZN	B	1301	1/1	0.99	0.05	94,94,94,94	0
3	ZN	A	1301	1/1	0.99	0.05	70,70,70,70	0
3	ZN	D	1301	1/1	0.99	0.03	113,113,113,113	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.



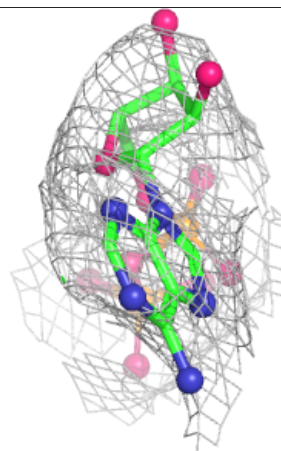
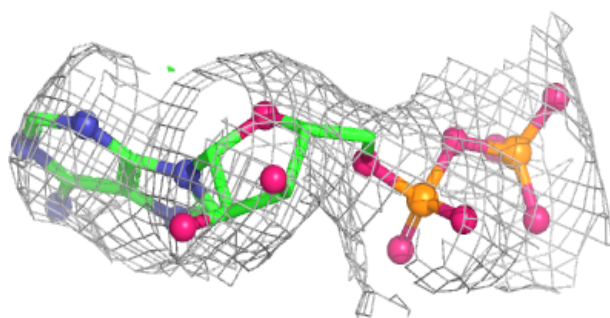
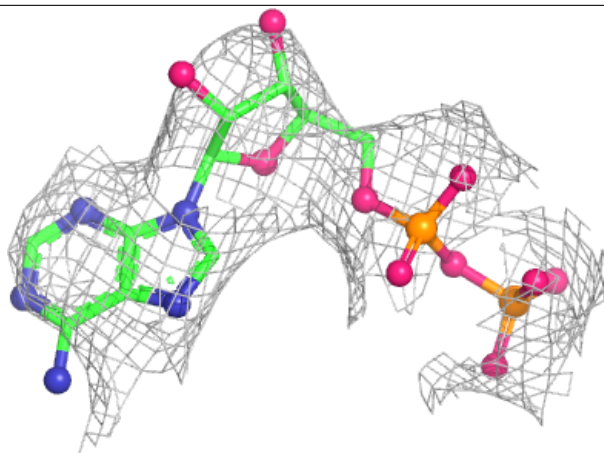
Electron density around ADP E 1302:

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



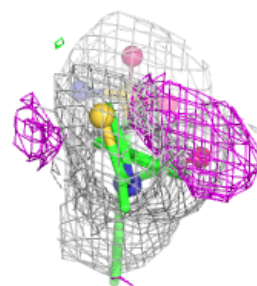
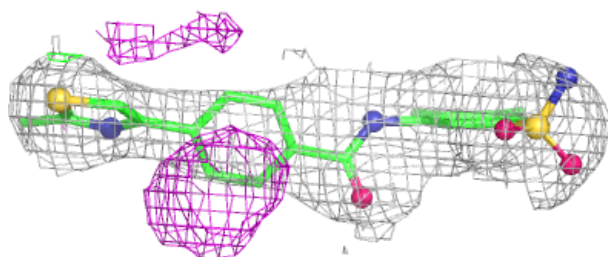
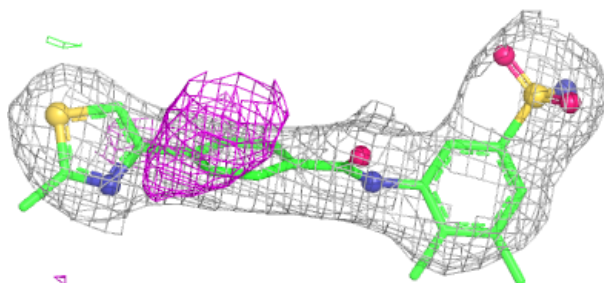
Electron density around ADP D 1302:

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

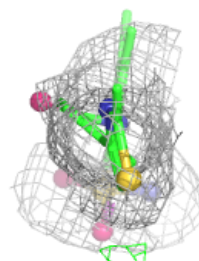
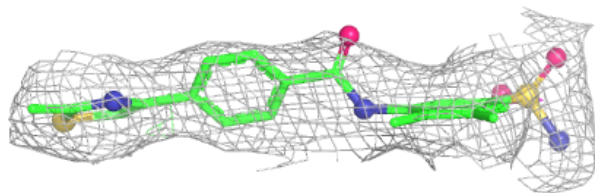
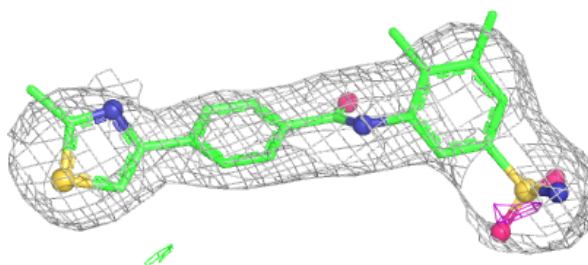


Electron density around RY8 F 1303:

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

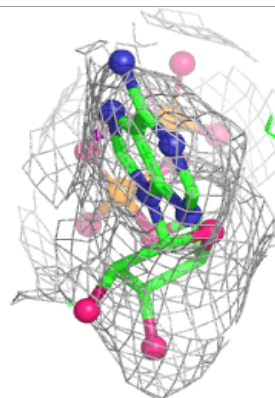
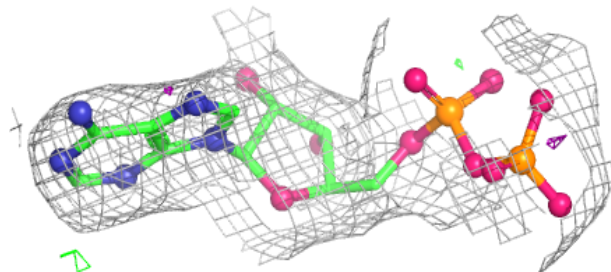
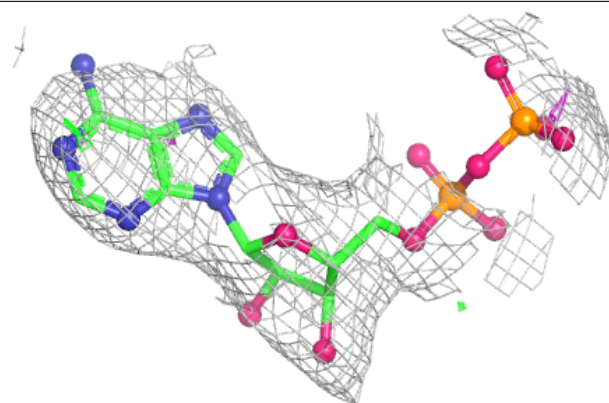
**Electron density around RY8 D 1303:**

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

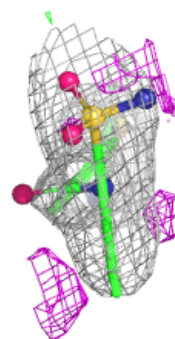
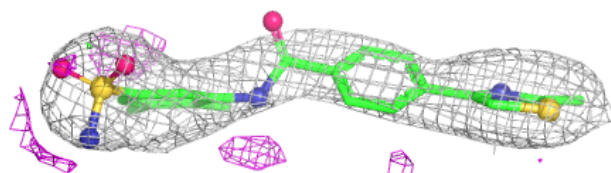
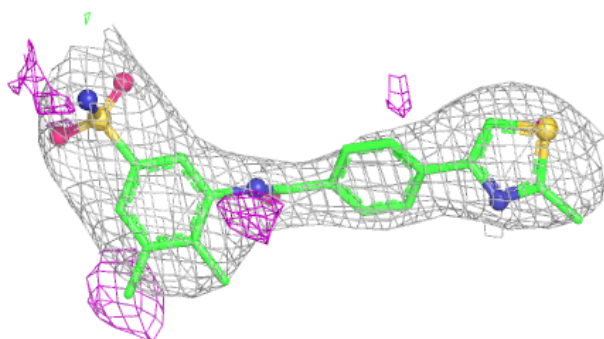


Electron density around ADP C 1302:

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and green (positive)

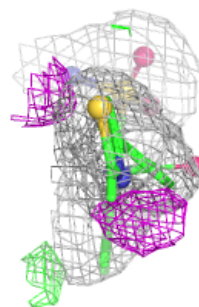
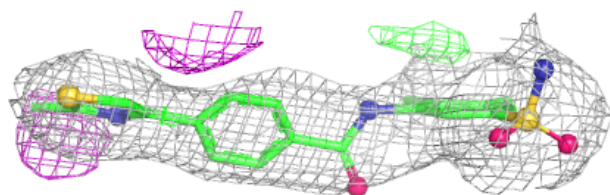
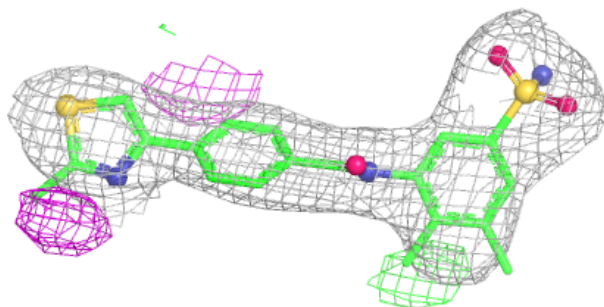
**Electron density around RY8 C 1303:**

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and green (positive)

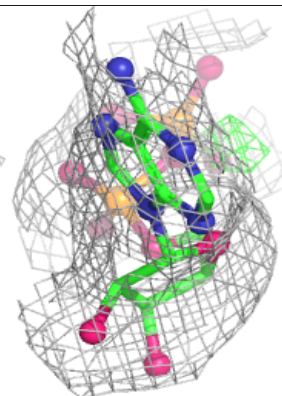
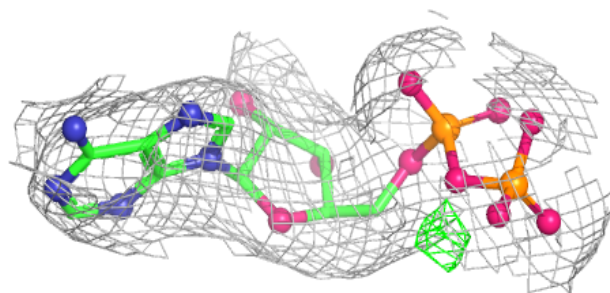
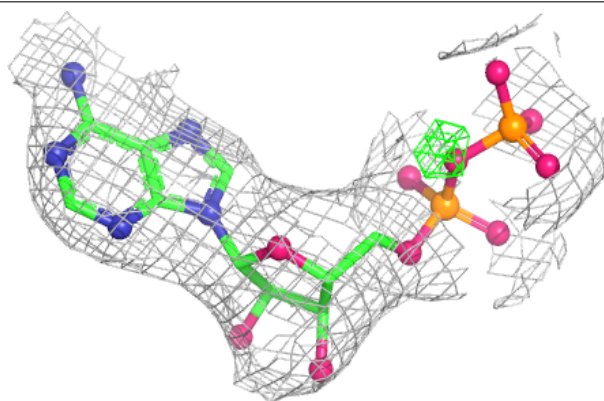


Electron density around RY8 E 1303:

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

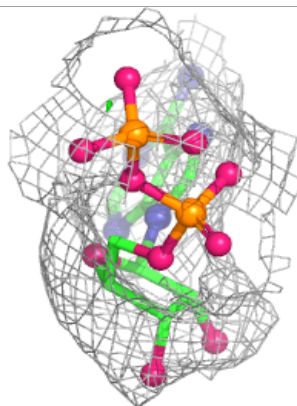
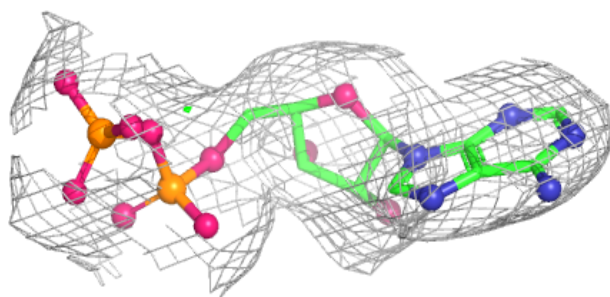
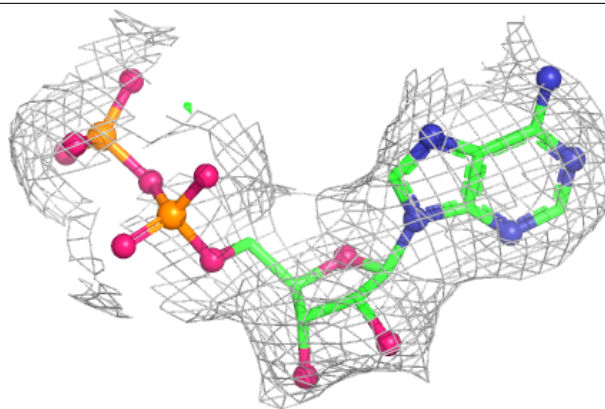
**Electron density around ADP A 1302:**

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and green (positive)

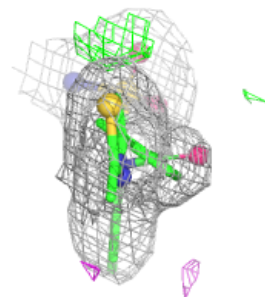
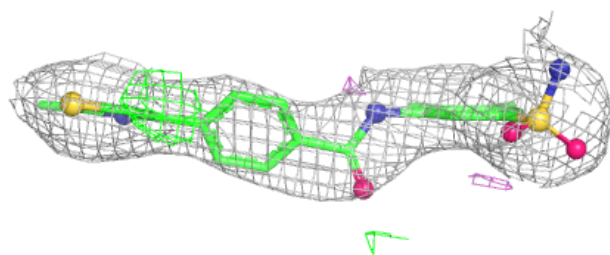
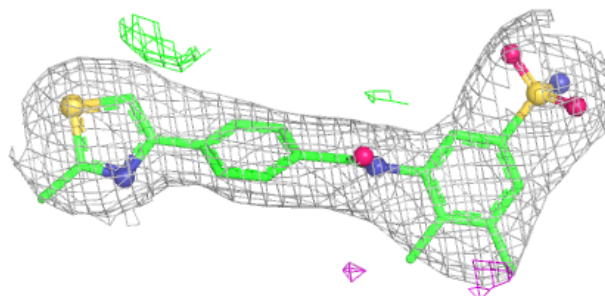


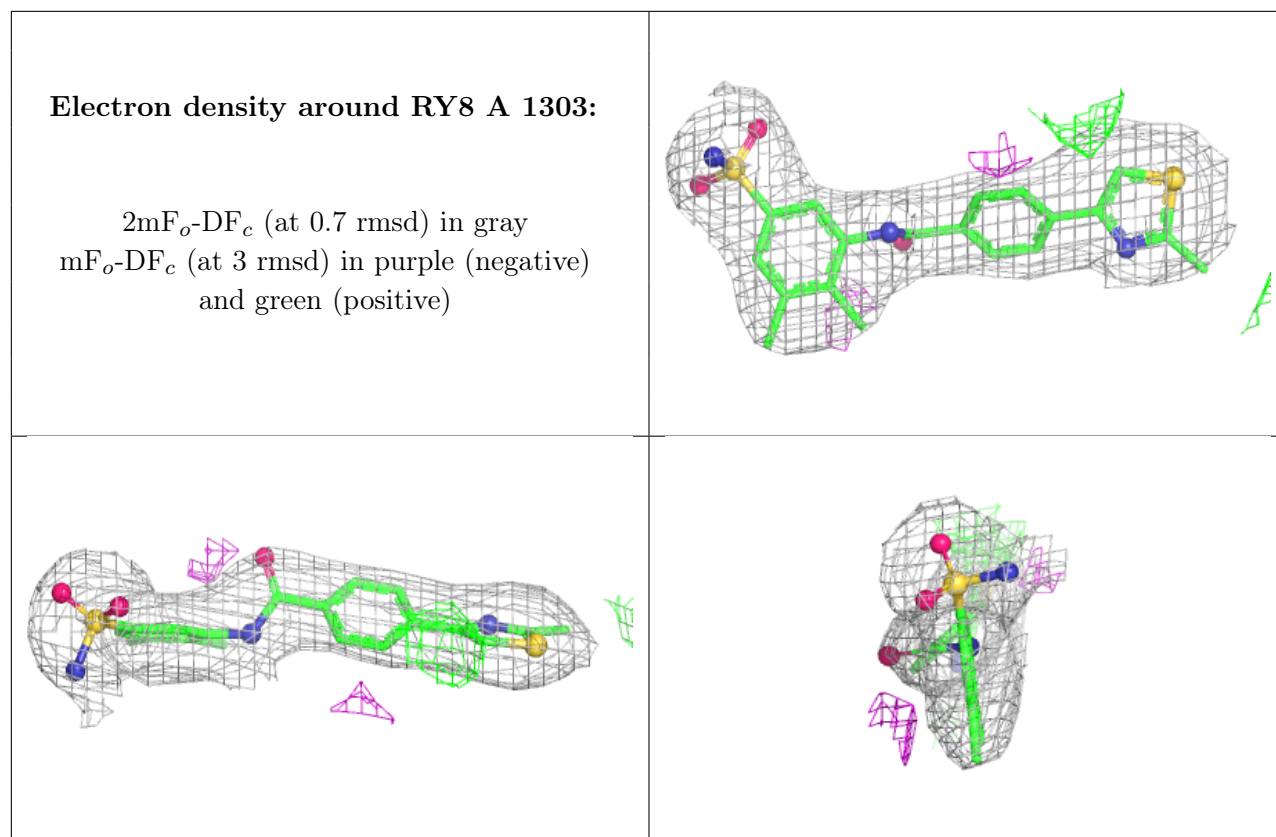
Electron density around ADP B 1302:

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

**Electron density around RY8 B 1303:**

$2mF_o - DF_c$ (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.