



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:03 AM UTC

PDB ID : 4K3N / pdb_00004k3n
Title : Phosphonic Arginine Mimetics as Inhibitors of the M17 Aminopeptidases from Plasmodium falciparum
Authors : McGowan, S.
Deposited on : 2013-04-11
Resolution : 2.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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

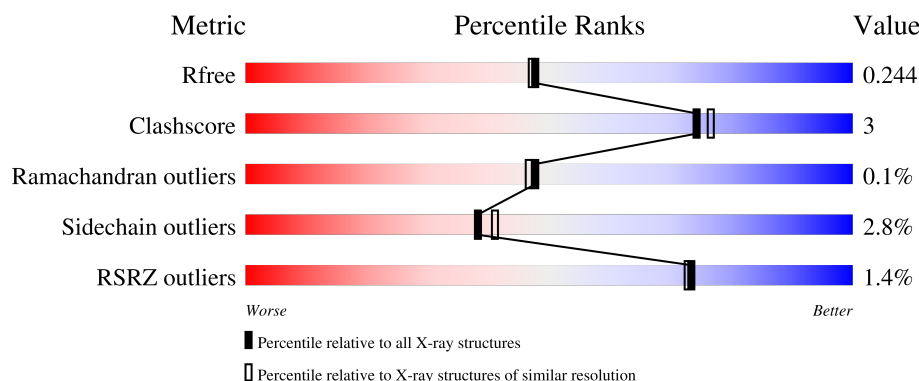
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	528	% 91% 8% .
1	B	528	2% 89% 9% ..
1	C	528	% 89% 9% .
1	D	528	2% 89% 8% ..
1	E	528	% 90% 7% .

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Mol	Chain	Length	Quality of chain
1	F	528	2%88%9%.
1	G	528	%91%7%..
1	H	528	2%89%8%..
1	I	528	%91%7%.
1	J	528	2%90%7%..
1	K	528	%90%6%.
1	L	528	2%89%7%..

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 49847 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 M17 leucyl aminopeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	519	Total	C	N	O	S	0	1	0
			3995	2564	642	769	20			
1	B	518	Total	C	N	O	S	0	0	0
			3939	2529	638	752	20			
1	C	518	Total	C	N	O	S	0	1	0
			3960	2547	639	754	20			
1	D	514	Total	C	N	O	S	0	0	0
			3947	2539	635	753	20			
1	E	510	Total	C	N	O	S	0	0	0
			3905	2513	626	747	19			
1	F	511	Total	C	N	O	S	0	0	0
			3849	2472	624	734	19			
1	G	519	Total	C	N	O	S	0	0	0
			3993	2562	642	770	19			
1	H	516	Total	C	N	O	S	0	0	0
			3914	2513	632	750	19			
1	I	518	Total	C	N	O	S	0	0	0
			3964	2549	638	757	20			
1	J	513	Total	C	N	O	S	0	0	0
			3939	2536	636	747	20			
1	K	509	Total	C	N	O	S	0	0	0
			3906	2515	625	747	19			
1	L	513	Total	C	N	O	S	0	0	0
			3885	2494	628	744	19			

There are 108 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	152	GLN	ASN	engineered mutation	UNP Q8IL11
A	515	GLN	ASN	engineered mutation	UNP Q8IL11
A	546	GLN	ASN	engineered mutation	UNP Q8IL11
A	606	HIS	-	expression tag	UNP Q8IL11
A	607	HIS	-	expression tag	UNP Q8IL11

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Chain	Residue	Modelled	Actual	Comment	Reference
A	608	HIS	-	expression tag	UNP Q8IL11
A	609	HIS	-	expression tag	UNP Q8IL11
A	610	HIS	-	expression tag	UNP Q8IL11
A	611	HIS	-	expression tag	UNP Q8IL11
B	152	GLN	ASN	engineered mutation	UNP Q8IL11
B	515	GLN	ASN	engineered mutation	UNP Q8IL11
B	546	GLN	ASN	engineered mutation	UNP Q8IL11
B	606	HIS	-	expression tag	UNP Q8IL11
B	607	HIS	-	expression tag	UNP Q8IL11
B	608	HIS	-	expression tag	UNP Q8IL11
B	609	HIS	-	expression tag	UNP Q8IL11
B	610	HIS	-	expression tag	UNP Q8IL11
B	611	HIS	-	expression tag	UNP Q8IL11
C	152	GLN	ASN	engineered mutation	UNP Q8IL11
C	515	GLN	ASN	engineered mutation	UNP Q8IL11
C	546	GLN	ASN	engineered mutation	UNP Q8IL11
C	606	HIS	-	expression tag	UNP Q8IL11
C	607	HIS	-	expression tag	UNP Q8IL11
C	608	HIS	-	expression tag	UNP Q8IL11
C	609	HIS	-	expression tag	UNP Q8IL11
C	610	HIS	-	expression tag	UNP Q8IL11
C	611	HIS	-	expression tag	UNP Q8IL11
D	152	GLN	ASN	engineered mutation	UNP Q8IL11
D	515	GLN	ASN	engineered mutation	UNP Q8IL11
D	546	GLN	ASN	engineered mutation	UNP Q8IL11
D	606	HIS	-	expression tag	UNP Q8IL11
D	607	HIS	-	expression tag	UNP Q8IL11
D	608	HIS	-	expression tag	UNP Q8IL11
D	609	HIS	-	expression tag	UNP Q8IL11
D	610	HIS	-	expression tag	UNP Q8IL11
D	611	HIS	-	expression tag	UNP Q8IL11
E	152	GLN	ASN	engineered mutation	UNP Q8IL11
E	515	GLN	ASN	engineered mutation	UNP Q8IL11
E	546	GLN	ASN	engineered mutation	UNP Q8IL11
E	606	HIS	-	expression tag	UNP Q8IL11
E	607	HIS	-	expression tag	UNP Q8IL11
E	608	HIS	-	expression tag	UNP Q8IL11
E	609	HIS	-	expression tag	UNP Q8IL11
E	610	HIS	-	expression tag	UNP Q8IL11
E	611	HIS	-	expression tag	UNP Q8IL11
F	152	GLN	ASN	engineered mutation	UNP Q8IL11
F	515	GLN	ASN	engineered mutation	UNP Q8IL11

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Chain	Residue	Modelled	Actual	Comment	Reference
F	546	GLN	ASN	engineered mutation	UNP Q8IL11
F	606	HIS	-	expression tag	UNP Q8IL11
F	607	HIS	-	expression tag	UNP Q8IL11
F	608	HIS	-	expression tag	UNP Q8IL11
F	609	HIS	-	expression tag	UNP Q8IL11
F	610	HIS	-	expression tag	UNP Q8IL11
F	611	HIS	-	expression tag	UNP Q8IL11
G	152	GLN	ASN	engineered mutation	UNP Q8IL11
G	515	GLN	ASN	engineered mutation	UNP Q8IL11
G	546	GLN	ASN	engineered mutation	UNP Q8IL11
G	606	HIS	-	expression tag	UNP Q8IL11
G	607	HIS	-	expression tag	UNP Q8IL11
G	608	HIS	-	expression tag	UNP Q8IL11
G	609	HIS	-	expression tag	UNP Q8IL11
G	610	HIS	-	expression tag	UNP Q8IL11
G	611	HIS	-	expression tag	UNP Q8IL11
H	152	GLN	ASN	engineered mutation	UNP Q8IL11
H	515	GLN	ASN	engineered mutation	UNP Q8IL11
H	546	GLN	ASN	engineered mutation	UNP Q8IL11
H	606	HIS	-	expression tag	UNP Q8IL11
H	607	HIS	-	expression tag	UNP Q8IL11
H	608	HIS	-	expression tag	UNP Q8IL11
H	609	HIS	-	expression tag	UNP Q8IL11
H	610	HIS	-	expression tag	UNP Q8IL11
H	611	HIS	-	expression tag	UNP Q8IL11
I	152	GLN	ASN	engineered mutation	UNP Q8IL11
I	515	GLN	ASN	engineered mutation	UNP Q8IL11
I	546	GLN	ASN	engineered mutation	UNP Q8IL11
I	606	HIS	-	expression tag	UNP Q8IL11
I	607	HIS	-	expression tag	UNP Q8IL11
I	608	HIS	-	expression tag	UNP Q8IL11
I	609	HIS	-	expression tag	UNP Q8IL11
I	610	HIS	-	expression tag	UNP Q8IL11
I	611	HIS	-	expression tag	UNP Q8IL11
J	152	GLN	ASN	engineered mutation	UNP Q8IL11
J	515	GLN	ASN	engineered mutation	UNP Q8IL11
J	546	GLN	ASN	engineered mutation	UNP Q8IL11
J	606	HIS	-	expression tag	UNP Q8IL11
J	607	HIS	-	expression tag	UNP Q8IL11
J	608	HIS	-	expression tag	UNP Q8IL11
J	609	HIS	-	expression tag	UNP Q8IL11
J	610	HIS	-	expression tag	UNP Q8IL11

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Chain	Residue	Modelled	Actual	Comment	Reference
J	611	HIS	-	expression tag	UNP Q8IL11
K	152	GLN	ASN	engineered mutation	UNP Q8IL11
K	515	GLN	ASN	engineered mutation	UNP Q8IL11
K	546	GLN	ASN	engineered mutation	UNP Q8IL11
K	606	HIS	-	expression tag	UNP Q8IL11
K	607	HIS	-	expression tag	UNP Q8IL11
K	608	HIS	-	expression tag	UNP Q8IL11
K	609	HIS	-	expression tag	UNP Q8IL11
K	610	HIS	-	expression tag	UNP Q8IL11
K	611	HIS	-	expression tag	UNP Q8IL11
L	152	GLN	ASN	engineered mutation	UNP Q8IL11
L	515	GLN	ASN	engineered mutation	UNP Q8IL11
L	546	GLN	ASN	engineered mutation	UNP Q8IL11
L	606	HIS	-	expression tag	UNP Q8IL11
L	607	HIS	-	expression tag	UNP Q8IL11
L	608	HIS	-	expression tag	UNP Q8IL11
L	609	HIS	-	expression tag	UNP Q8IL11
L	610	HIS	-	expression tag	UNP Q8IL11
L	611	HIS	-	expression tag	UNP Q8IL11

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

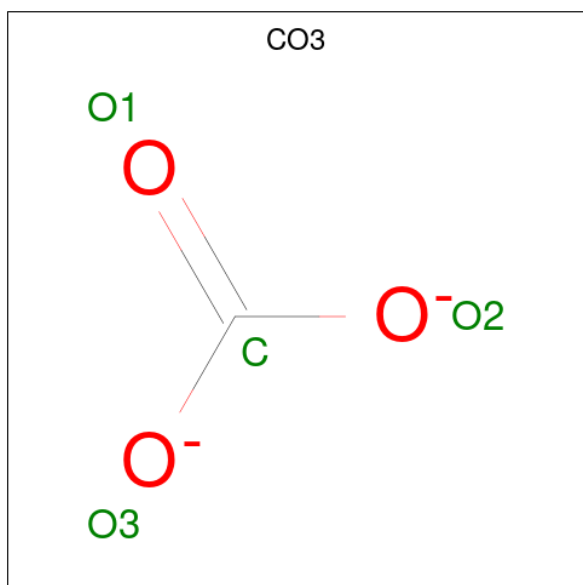
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Zn 2 2	0	0
2	B	2	Total Zn 2 2	0	0
2	C	2	Total Zn 2 2	0	0
2	D	2	Total Zn 2 2	0	0
2	E	2	Total Zn 2 2	0	0
2	F	2	Total Zn 2 2	0	0
2	G	2	Total Zn 2 2	0	0
2	H	2	Total Zn 2 2	0	0
2	I	2	Total Zn 2 2	0	0
2	J	2	Total Zn 2 2	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	K	2	Total	Zn	0	0
			2	2		
2	L	2	Total	Zn	0	0
			2	2		

- Molecule 3 is CARBONATE ION (CCD ID: CO3) (formula: CO₃).



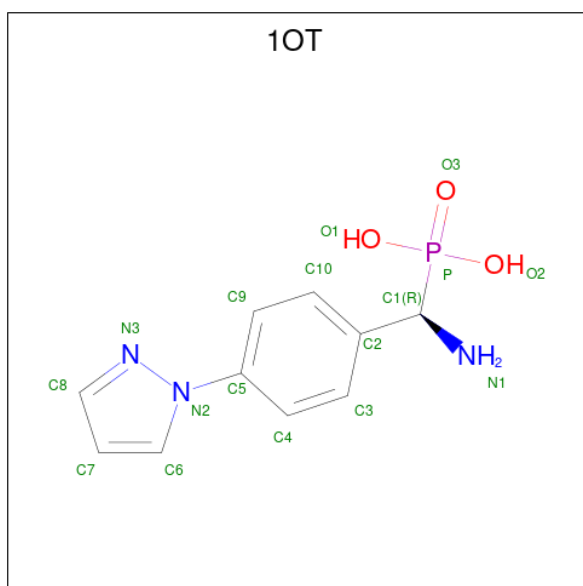
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	1	3		
3	B	1	Total	C	O	0	0
			4	1	3		
3	C	1	Total	C	O	0	0
			4	1	3		
3	D	1	Total	C	O	0	0
			4	1	3		
3	E	1	Total	C	O	0	0
			4	1	3		
3	F	1	Total	C	O	0	0
			4	1	3		
3	G	1	Total	C	O	0	0
			4	1	3		
3	H	1	Total	C	O	0	0
			4	1	3		
3	I	1	Total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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	J	1	Total	C	O	0	0
			4	1	3		
3	K	1	Total	C	O	0	0
			4	1	3		
3	L	1	Total	C	O	0	0
			4	1	3		

- Molecule 4 is {(R)-amino[4-(1H-pyrazol-1-yl)phenyl]methyl}phosphonic acid (CCD ID: 1OT) (formula: C₁₀H₁₂N₃O₃P).



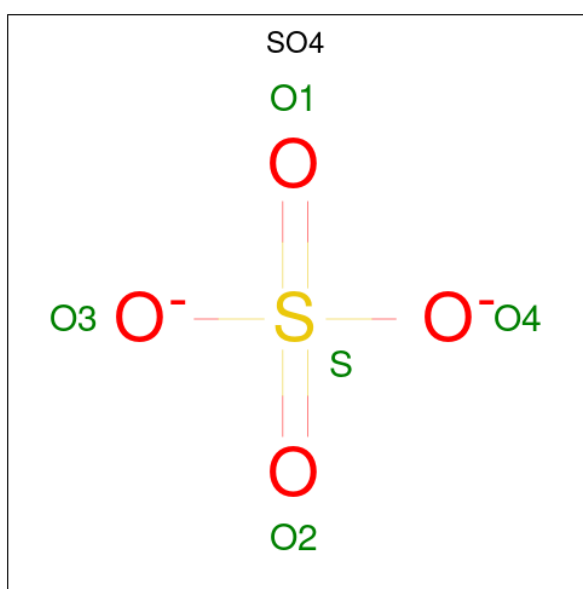
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	B	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	C	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	D	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	E	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	F	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	G	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	H	1	Total	C	N	O	P	0	0
			17	10	3	3	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	I	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	J	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	K	1	Total	C	N	O	P	0	0
			17	10	3	3	1		
4	L	1	Total	C	N	O	P	0	0
			17	10	3	3	1		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



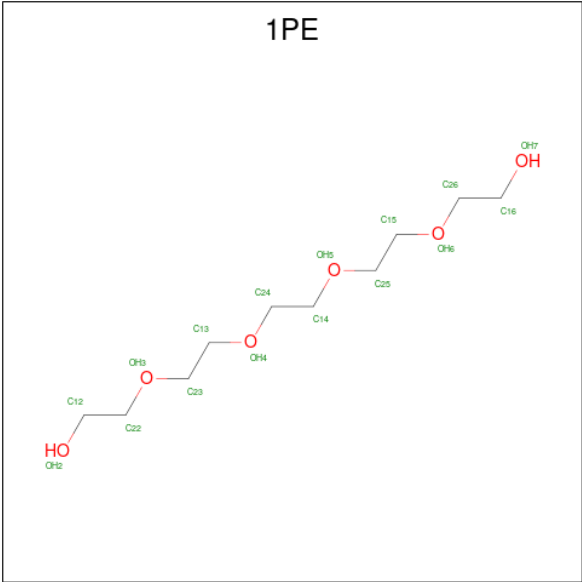
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	O	S	0	0
			5	4	1		
5	E	1	Total	O	S	0	0
			5	4	1		
5	E	1	Total	O	S	0	0
			5	4	1		
5	F	1	Total	O	S	0	0
			5	4	1		
5	G	1	Total	O	S	0	0
			5	4	1		
5	G	1	Total	O	S	0	0
			5	4	1		
5	G	1	Total	O	S	0	0
			5	4	1		
5	H	1	Total	O	S	0	0
			5	4	1		
5	I	1	Total	O	S	0	0
			5	4	1		
5	I	1	Total	O	S	0	0
			5	4	1		
5	J	1	Total	O	S	0	0
			5	4	1		
5	K	1	Total	O	S	0	0
			5	4	1		
5	K	1	Total	O	S	0	0
			5	4	1		
5	L	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).



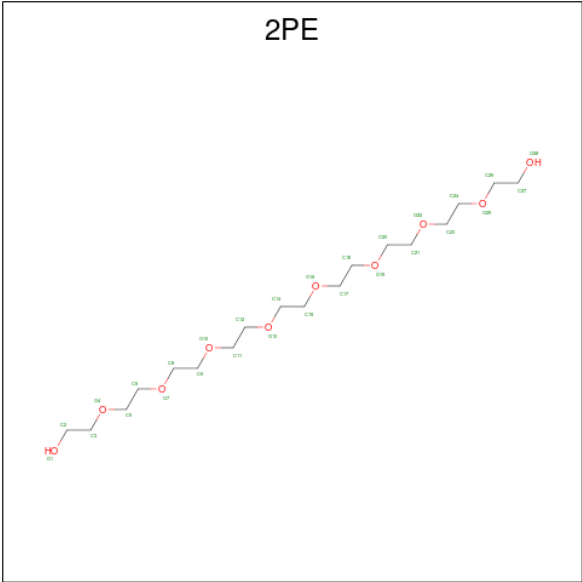
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			9	6	3		
6	A	1	Total	C	O	0	0
			12	8	4		
6	C	1	Total	C	O	0	0
			13	9	4		
6	C	1	Total	C	O	0	0
			9	6	3		
6	C	1	Total	C	O	0	0
			7	5	2		
6	D	1	Total	C	O	0	0
			10	7	3		
6	D	1	Total	C	O	0	0
			9	6	3		
6	D	1	Total	C	O	0	0
			11	8	3		
6	E	1	Total	C	O	0	0
			12	8	4		
6	E	1	Total	C	O	0	0
			12	8	4		
6	E	1	Total	C	O	0	0
			8	5	3		
6	F	1	Total	C	O	0	0
			10	6	4		
6	F	1	Total	C	O	0	0
			10	6	4		
6	F	1	Total	C	O	0	0
			10	6	4		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	F	1	Total C O 11 7 4	0	0
6	G	1	Total C O 9 6 3	0	0
6	G	1	Total C O 6 4 2	0	0
6	G	1	Total C O 6 4 2	0	0
6	G	1	Total C O 15 10 5	0	0
6	I	1	Total C O 15 10 5	0	0
6	I	1	Total C O 11 8 3	0	0
6	I	1	Total C O 5 3 2	0	0
6	J	1	Total C O 11 7 4	0	0
6	J	1	Total C O 10 6 4	0	0
6	J	1	Total C O 11 8 3	0	0
6	J	1	Total C O 9 6 3	0	0
6	K	1	Total C O 12 8 4	0	0
6	K	1	Total C O 12 8 4	0	0
6	K	1	Total C O 11 7 4	0	0
6	K	1	Total C O 6 4 2	0	0
6	L	1	Total C O 10 6 4	0	0
6	L	1	Total C O 12 8 4	0	0
6	L	1	Total C O 8 5 3	0	0
6	L	1	Total C O 11 7 4	0	0

- Molecule 7 is NONAETHYLENE GLYCOL (CCD ID: 2PE) (formula: C₁₈H₃₈O₁₀).

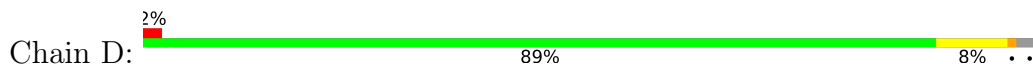


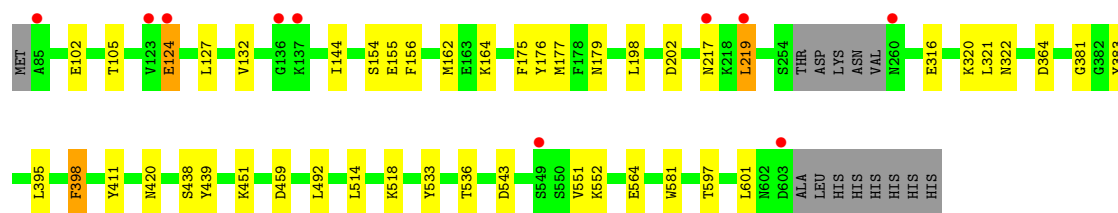
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			26	17	9		
7	B	1	Total	C	O	0	0
			10	6	4		
7	C	1	Total	C	O	0	0
			9	6	3		
7	D	1	Total	C	O	0	0
			13	9	4		
7	D	1	Total	C	O	0	0
			5	3	2		
7	D	1	Total	C	O	0	0
			9	6	3		
7	E	1	Total	C	O	0	0
			9	6	3		
7	F	1	Total	C	O	0	0
			9	6	3		
7	H	1	Total	C	O	0	0
			25	16	9		
7	H	1	Total	C	O	0	0
			10	7	3		
7	I	1	Total	C	O	0	0
			8	5	3		
7	J	1	Total	C	O	0	0
			10	7	3		

- Molecule 8 is water.

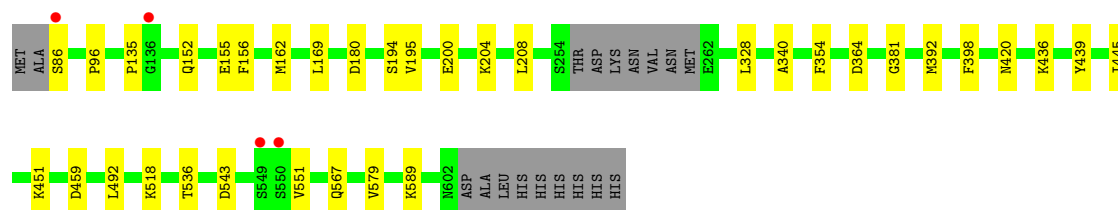
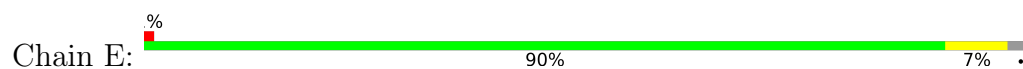
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	173	Total 173	O 173	0	0
8	B	125	Total 125	O 125	0	0
8	C	168	Total 168	O 168	0	0
8	D	150	Total 150	O 150	0	0
8	E	193	Total 193	O 193	0	0
8	F	117	Total 117	O 117	0	0
8	G	145	Total 145	O 145	0	0
8	H	129	Total 129	O 129	0	0
8	I	146	Total 146	O 146	0	0
8	J	137	Total 137	O 137	0	0
8	K	175	Total 175	O 175	0	0
8	L	126	Total 126	O 126	0	0

- Molecule 1: M17 leucyl aminopeptidase

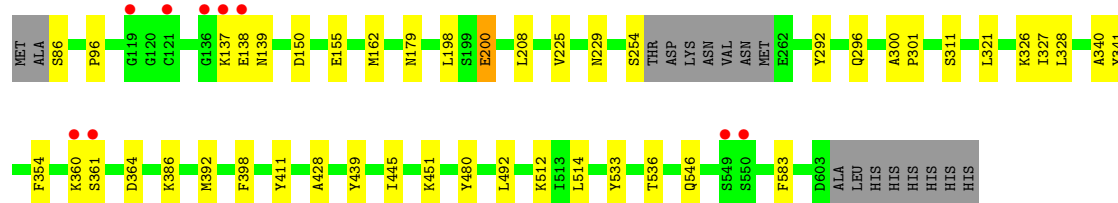
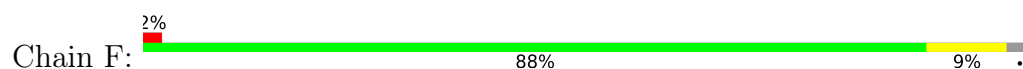




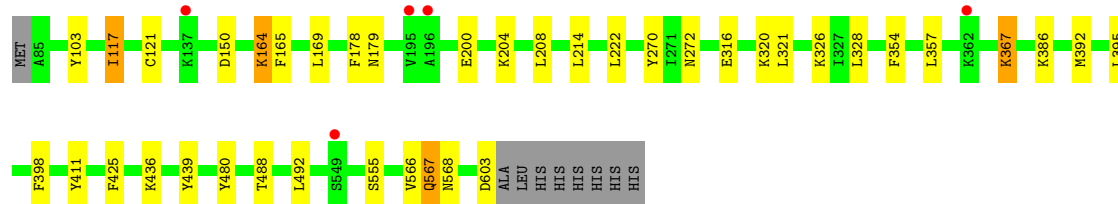
• Molecule 1: M17 leucyl aminopeptidase



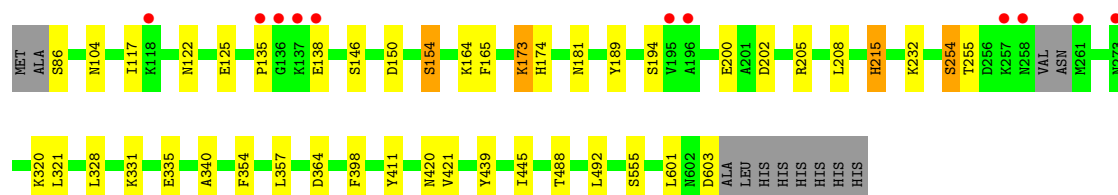
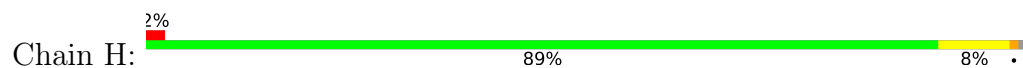
• Molecule 1: M17 leucyl aminopeptidase



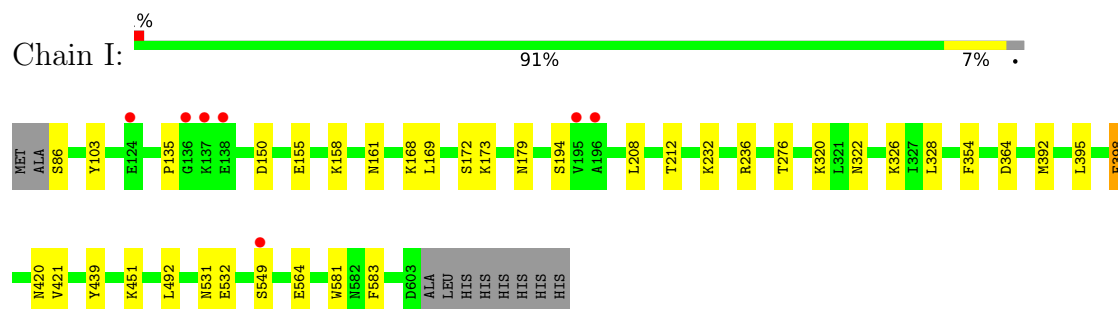
• Molecule 1: M17 leucyl aminopeptidase



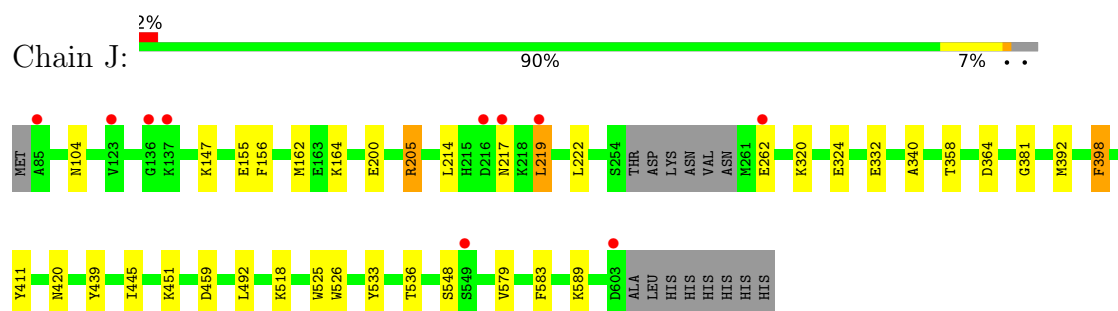
• Molecule 1: M17 leucyl aminopeptidase



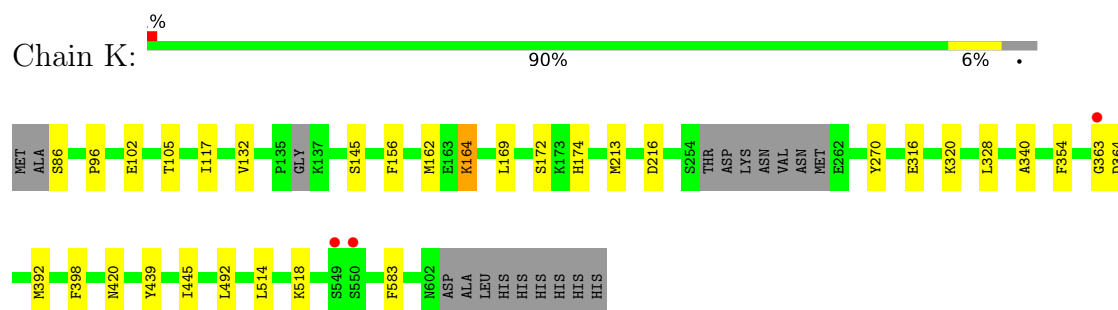
- Molecule 1: M17 leucyl aminopeptidase



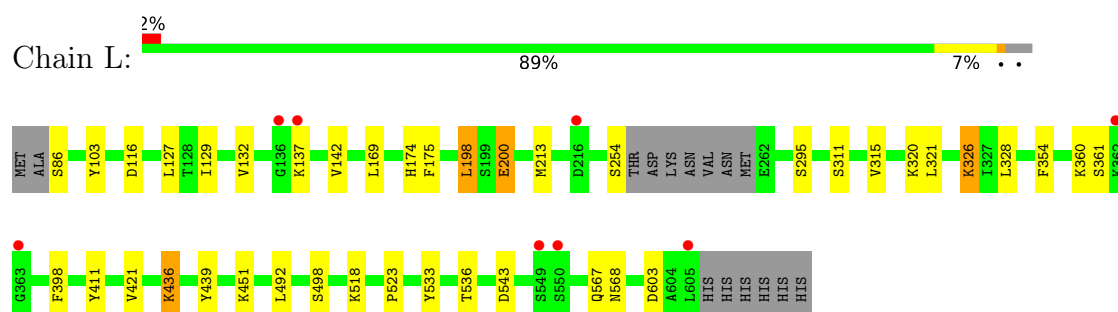
- Molecule 1: M17 leucyl aminopeptidase



- Molecule 1: M17 leucyl aminopeptidase



- Molecule 1: M17 leucyl aminopeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	173.72Å 177.08Å 229.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.11 – 2.00 45.11 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.11-2.00) 99.9 (45.11-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.00Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.196 , 0.240 0.205 , 0.244	Depositor DCC
R_{free} test set	23744 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.725	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	49847	wwPDB-VP
Average B, all atoms (Å ²)	16.0	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 43.71 % 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 1.6963e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: 1PE, ZN, 2PE, CO3, SO4, 1OT

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	0.43	0/4075	0.76	0/5525
1	B	0.42	0/4017	0.77	1/5455 (0.0%)
1	C	0.44	0/4041	0.75	0/5483
1	D	0.43	0/4024	0.77	1/5455 (0.0%)
1	E	0.44	0/3982	0.76	0/5403
1	F	0.41	0/3925	0.80	5/5336 (0.1%)
1	G	0.47	2/4071 (0.0%)	0.76	1/5521 (0.0%)
1	H	0.42	0/3991	0.77	1/5422 (0.0%)
1	I	0.43	0/4042	0.75	0/5484
1	J	0.43	0/4016	0.76	0/5443
1	K	0.44	0/3982	0.77	0/5400
1	L	0.43	0/3962	0.75	0/5384
All	All	0.43	2/48128 (0.0%)	0.76	9/65311 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	567	GLN	C-O	-5.33	1.17	1.24
1	G	566	VAL	C-O	-5.16	1.18	1.24

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	137	LYS	CB-CA-C	-10.55	104.41	116.54
1	F	137	LYS	N-CA-C	7.66	119.96	108.31
1	B	215	HIS	N-CA-C	6.85	119.76	111.82
1	F	138	GLU	CA-C-N	6.74	133.83	121.70
1	F	138	GLU	C-N-CA	6.74	133.83	121.70
1	F	139	ASN	N-CA-C	5.60	126.68	111.00
1	H	215	HIS	N-CA-C	5.55	117.41	111.36
1	G	568	ASN	N-CA-C	5.26	117.19	110.61
1	D	551	VAL	N-CA-C	5.07	115.63	109.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	257	LYS	Peptide

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	3995	0	3929	23	0
1	B	3939	0	3835	22	0
1	C	3960	0	3890	20	0
1	D	3947	0	3887	42	0
1	E	3905	0	3830	21	0
1	F	3849	0	3722	25	0
1	G	3993	0	3922	27	0
1	H	3914	0	3793	18	0
1	I	3964	0	3897	24	0
1	J	3939	0	3892	29	0
1	K	3906	0	3837	17	0
1	L	3885	0	3760	26	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	2	0	0	0	0
2	H	2	0	0	0	0
2	I	2	0	0	0	0
2	J	2	0	0	0	0
2	K	2	0	0	0	0
2	L	2	0	0	0	0
3	A	4	0	0	0	0
3	B	4	0	0	0	0
3	C	4	0	0	0	0
3	D	4	0	0	1	0
3	E	4	0	0	0	0
3	F	4	0	0	0	0
3	G	4	0	0	0	0
3	H	4	0	0	0	0
3	I	4	0	0	0	0
3	J	4	0	0	0	0
3	K	4	0	0	0	0
3	L	4	0	0	1	0
4	A	17	0	10	4	0
4	B	17	0	11	4	0
4	C	17	0	11	3	0
4	D	17	0	12	2	0
4	E	17	0	11	4	0
4	F	17	0	12	4	0
4	G	17	0	10	4	0
4	H	17	0	12	2	0
4	I	17	0	12	3	0
4	J	17	0	11	3	0
4	K	17	0	12	3	0
4	L	17	0	11	4	0
5	A	20	0	0	0	0
5	B	5	0	0	0	0
5	C	5	0	0	0	0
5	D	10	0	0	0	0
5	E	10	0	0	0	0
5	F	5	0	0	0	0
5	G	15	0	0	0	0
5	H	5	0	0	0	0
5	I	10	0	0	0	0
5	J	5	0	0	1	0
5	K	10	0	0	0	0
5	L	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	21	0	22	1	0
6	C	29	0	30	2	0
6	D	30	0	30	7	0
6	E	32	0	36	4	0
6	F	41	0	51	2	0
6	G	36	0	39	8	0
6	I	31	0	35	3	0
6	J	41	0	48	8	0
6	K	41	0	44	3	0
6	L	41	0	48	10	0
7	B	36	0	45	0	0
7	C	9	0	10	2	0
7	D	27	0	28	5	0
7	E	9	0	10	1	0
7	F	9	0	10	1	0
7	H	35	0	43	2	0
7	I	8	0	8	0	0
7	J	10	0	10	0	0
8	A	173	0	0	0	0
8	B	125	0	0	0	0
8	C	168	0	0	1	0
8	D	150	0	0	6	0
8	E	193	0	0	0	0
8	F	117	0	0	2	0
8	G	145	0	0	0	0
8	H	129	0	0	1	0
8	I	146	0	0	3	0
8	J	137	0	0	3	0
8	K	175	0	0	2	0
8	L	126	0	0	2	0
All	All	49847	0	46876	298	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 3.

All (298) 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:I:155:GLU:O	8:I:1103:HOH:O	1.97	0.82
1:D:320:LYS:HZ1	7:D:1010:2PE:H121	1.46	0.80
1:I:161:ASN:ND2	8:I:1103:HOH:O	2.17	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:526:TRP:HB3	6:J:1009:1PE:H242	1.67	0.76
1:G:178:PHE:HZ	1:J:155:GLU:HG2	1.53	0.73
1:J:536:THR:OG1	8:J:1221:HOH:O	2.06	0.73
1:A:178:PHE:HZ	1:D:155:GLU:HG2	1.54	0.72
1:G:320:LYS:HZ1	6:G:1011:1PE:H251	1.56	0.71
1:J:411:TYR:HE1	6:J:1006:1PE:H151	1.54	0.71
1:J:525:TRP:HA	6:J:1009:1PE:H132	1.70	0.71
1:L:320:LYS:HZ1	6:L:1009:1PE:H252	1.56	0.70
1:K:316:GLU:HG3	6:K:1007:1PE:H152	1.73	0.70
1:K:328:LEU:HB2	1:K:354:PHE:HB3	1.72	0.69
1:J:320:LYS:HB3	6:J:1006:1PE:H161	1.74	0.69
1:H:320:LYS:HB3	7:H:1006:2PE:H211	1.75	0.69
1:J:533:TYR:O	1:J:536:THR:HG22	1.96	0.66
1:B:236:ARG:NE	1:B:240:GLU:OE2	2.21	0.65
1:L:436:LYS:NZ	8:L:1226:HOH:O	2.17	0.65
1:G:567:GLN:HA	1:G:567:GLN:OE1	1.95	0.65
1:L:328:LEU:HB2	1:L:354:PHE:HB3	1.79	0.64
1:F:546:GLN:O	8:F:1209:HOH:O	2.15	0.64
1:B:392:MET:HE1	1:B:583:PHE:CE2	2.33	0.64
1:C:178:PHE:HZ	1:E:155:GLU:HG2	1.62	0.62
1:G:316:GLU:HG3	6:G:1011:1PE:H231	1.80	0.62
1:J:411:TYR:CE1	6:J:1006:1PE:H151	2.32	0.62
1:A:328:LEU:HB2	1:A:354:PHE:HB3	1.81	0.62
1:A:156:PHE:CD2	1:A:156:PHE:CZ	2.86	0.61
1:E:328:LEU:HB2	1:E:354:PHE:HB3	1.80	0.61
1:E:543:ASP:HB3	6:E:1009:1PE:H232	1.81	0.61
1:D:533:TYR:O	1:D:536:THR:HG22	2.01	0.61
1:B:230:VAL:HG12	1:B:234:LEU:HD23	1.84	0.60
1:A:366:LYS:HG2	1:A:420:ASN:HB3	1.84	0.60
1:J:451:LYS:HE3	6:J:1008:1PE:H261	1.82	0.60
1:L:492:LEU:HD11	4:L:1004:1OT:C7	2.33	0.59
1:F:321:LEU:HD11	1:F:411:TYR:HA	1.84	0.59
1:D:124:GLU:HA	1:D:179:ASN:HD22	1.67	0.59
1:D:177:MET:HE2	8:D:1131:HOH:O	2.01	0.59
1:G:214:LEU:HD21	1:G:222:LEU:HD22	1.84	0.59
1:G:392:MET:HE2	4:G:1004:1OT:C8	2.33	0.59
1:G:411:TYR:HE1	6:G:1008:1PE:H132	1.68	0.59
1:D:411:TYR:HE1	6:D:1007:1PE:H232	1.68	0.59
1:G:328:LEU:HB2	1:G:354:PHE:HB3	1.85	0.58
1:I:392:MET:HE1	1:I:583:PHE:CE2	2.38	0.58
1:D:492:LEU:HD11	4:D:1004:1OT:C7	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:320:LYS:HE2	6:L:1009:1PE:H141	1.86	0.58
1:B:328:LEU:HB2	1:B:354:PHE:HB3	1.86	0.58
1:F:514:LEU:HD22	6:F:1009:1PE:H142	1.85	0.58
1:H:331:LYS:O	1:H:335:GLU:HG3	2.02	0.57
1:L:320:LYS:NZ	6:L:1009:1PE:H252	2.18	0.57
1:F:533:TYR:O	1:F:536:THR:HG22	2.04	0.57
1:A:173:LYS:NZ	1:D:217:ASN:OD1	2.36	0.57
1:A:392:MET:HE2	4:A:1004:1OT:C8	2.35	0.57
1:B:538:ASN:O	1:B:545:ASN:ND2	2.38	0.57
1:E:451:LYS:HG2	6:E:1009:1PE:H231	1.85	0.57
1:D:320:LYS:HZ3	7:D:1010:2PE:H201	1.70	0.57
1:F:451:LYS:HG2	6:F:1008:1PE:H141	1.87	0.57
1:B:492:LEU:HD11	4:B:1004:1OT:H12	1.87	0.57
1:E:392:MET:HE2	4:E:1004:1OT:C8	2.35	0.56
1:I:328:LEU:HB2	1:I:354:PHE:HB3	1.85	0.56
1:E:381:GLY:HA2	1:E:459:ASP:OD1	2.06	0.56
1:J:392:MET:HE2	4:J:1004:1OT:C8	2.34	0.56
1:L:533:TYR:O	1:L:536:THR:HG22	2.05	0.56
1:K:174:HIS:ND1	1:K:213:MET:HE2	2.21	0.56
1:B:492:LEU:HD11	4:B:1004:1OT:C7	2.35	0.56
1:H:321:LEU:HD11	1:H:411:TYR:HA	1.88	0.56
1:D:320:LYS:NZ	7:D:1010:2PE:H121	2.21	0.55
1:E:492:LEU:HD11	4:E:1004:1OT:C7	2.35	0.55
1:K:96:PRO:HA	6:K:1010:1PE:H141	1.87	0.55
1:F:96:PRO:HA	7:F:1010:2PE:H111	1.88	0.55
1:H:492:LEU:HD11	4:H:1004:1OT:H12	1.88	0.55
1:C:172:SER:O	1:C:173:LYS:HD2	2.06	0.55
1:K:492:LEU:HD11	4:K:1004:1OT:H12	1.89	0.55
1:F:340:ALA:HA	1:F:445:ILE:HD12	1.87	0.55
1:B:214:LEU:HD21	1:B:222:LEU:HD22	1.86	0.55
1:F:328:LEU:HB2	1:F:354:PHE:HB3	1.87	0.55
1:D:451:LYS:NZ	6:D:1009:1PE:H242	2.22	0.54
1:K:492:LEU:HD11	4:K:1004:1OT:C7	2.38	0.54
1:A:178:PHE:CZ	1:D:155:GLU:HG2	2.39	0.54
1:C:392:MET:HE1	1:C:583:PHE:CE2	2.43	0.54
1:E:96:PRO:HA	7:E:1010:2PE:H182	1.90	0.54
1:G:492:LEU:HD11	4:G:1004:1OT:H12	1.90	0.54
1:J:340:ALA:HA	1:J:445:ILE:HD12	1.90	0.54
1:K:340:ALA:HA	1:K:445:ILE:HD12	1.90	0.53
1:A:176:TYR:HB3	8:D:1131:HOH:O	2.08	0.53
1:A:392:MET:HE1	1:A:583:PHE:CE2	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:320:LYS:HD3	6:K:1007:1PE:H222	1.91	0.53
1:G:392:MET:HE3	1:G:395:LEU:HD22	1.89	0.53
1:D:164:LYS:HE2	8:D:1240:HOH:O	2.08	0.53
1:G:386:LYS:NZ	4:G:1004:1OT:O2	2.38	0.53
1:G:326:LYS:HG2	1:G:328:LEU:HD12	1.91	0.53
1:C:328:LEU:HB2	1:C:354:PHE:HB3	1.90	0.52
1:L:174:HIS:ND1	1:L:213:MET:HE2	2.24	0.52
1:J:217:ASN:HB2	1:J:219:LEU:HD21	1.90	0.52
1:L:326:LYS:HE3	1:L:328:LEU:HD11	1.90	0.52
1:B:178:PHE:HZ	1:F:155:GLU:HG2	1.75	0.52
1:B:135:PRO:HA	1:B:194:SER:O	2.08	0.52
1:D:514:LEU:O	1:D:518:LYS:HG2	2.09	0.52
1:D:543:ASP:HB3	6:D:1009:1PE:H132	1.90	0.52
1:I:392:MET:HE2	4:I:1004:1OT:C8	2.40	0.52
4:H:1004:1OT:H10	4:H:1004:1OT:O2	2.10	0.52
1:D:217:ASN:HB3	1:D:219:LEU:HD21	1.92	0.52
1:B:172:SER:O	1:B:173:LYS:HD2	2.10	0.51
1:B:392:MET:HE3	1:B:395:LEU:HD22	1.92	0.51
1:H:340:ALA:HA	1:H:445:ILE:HD12	1.92	0.51
1:J:492:LEU:HD11	4:J:1004:1OT:C7	2.40	0.51
1:D:175:PHE:HB3	8:D:1131:HOH:O	2.09	0.51
1:F:480:TYR:OH	1:F:512:LYS:NZ	2.40	0.51
1:A:150:ASP:OD1	1:A:179:ASN:HB2	2.10	0.51
1:G:117:ILE:HG13	1:G:270:TYR:HB3	1.92	0.51
1:I:150:ASP:OD1	1:I:179:ASN:HB2	2.10	0.51
1:J:324:GLU:CG	1:J:358:THR:HB	2.41	0.51
1:L:103:TYR:HB2	6:L:1009:1PE:H161	1.93	0.50
1:K:392:MET:HE1	1:K:583:PHE:CE2	2.47	0.50
1:A:321:LEU:HD11	1:A:411:TYR:HA	1.94	0.50
1:E:536:THR:HG21	1:E:551:VAL:HG23	1.94	0.50
1:G:320:LYS:HZ2	6:G:1011:1PE:H161	1.77	0.50
1:E:200:GLU:O	1:E:204:LYS:HG3	2.12	0.49
1:A:492:LEU:HD11	4:A:1004:1OT:H12	1.93	0.49
1:C:411:TYR:HE1	6:C:1007:1PE:H231	1.76	0.49
1:I:395:LEU:HD11	1:I:581:TRP:CG	2.48	0.49
1:J:104:ASN:N	5:J:1005:SO4:O2	2.40	0.49
1:A:386:LYS:NZ	4:A:1004:1OT:O2	2.43	0.49
1:D:217:ASN:HB2	1:D:219:LEU:HD11	1.95	0.49
1:C:364:ASP:O	1:C:420:ASN:HA	2.12	0.49
1:F:200:GLU:HG3	8:F:1122:HOH:O	2.13	0.49
1:C:392:MET:HE1	1:C:583:PHE:HE2	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:178:PHE:CZ	1:J:155:GLU:HG2	2.42	0.49
1:J:548:SER:OG	8:J:1221:HOH:O	2.19	0.48
1:H:173:LYS:HB2	1:H:189:TYR:CE1	2.48	0.48
1:H:254:SER:OG	1:H:255:THR:N	2.44	0.48
1:J:156:PHE:O	1:J:162:MET:HE3	2.14	0.48
1:K:156:PHE:O	1:K:162:MET:HE3	2.13	0.48
1:B:364:ASP:O	1:B:420:ASN:HA	2.13	0.48
1:K:164:LYS:NZ	8:K:1188:HOH:O	2.46	0.48
1:H:364:ASP:O	1:H:420:ASN:HA	2.13	0.48
1:L:492:LEU:HD11	4:L:1004:1OT:H12	1.96	0.48
1:G:320:LYS:HD3	6:G:1011:1PE:H162	1.94	0.48
1:I:451:LYS:HE3	1:I:564:GLU:O	2.14	0.48
1:I:135:PRO:HA	1:I:194:SER:O	2.14	0.48
1:I:232:LYS:NZ	1:I:276:THR:O	2.46	0.48
1:I:392:MET:HE1	1:I:583:PHE:HE2	1.78	0.48
1:G:121:CYS:HA	1:G:270:TYR:CE2	2.49	0.47
1:G:320:LYS:NZ	6:G:1011:1PE:H251	2.26	0.47
1:A:316:GLU:HG3	6:A:1010:1PE:H131	1.95	0.47
1:B:340:ALA:HA	1:B:445:ILE:HD12	1.95	0.47
4:B:1004:1OT:H10	4:B:1004:1OT:O2	2.14	0.47
1:J:364:ASP:O	1:J:420:ASN:HA	2.15	0.47
1:H:202:ASP:OD1	1:H:205:ARG:NH1	2.48	0.47
1:K:102:GLU:HG3	8:K:1272:HOH:O	2.15	0.47
1:L:543:ASP:HB3	6:L:1007:1PE:H221	1.96	0.47
1:C:392:MET:HE2	4:C:1004:1OT:C8	2.44	0.47
1:F:492:LEU:HD11	4:F:1004:1OT:H12	1.96	0.47
1:G:103:TYR:CD1	6:G:1011:1PE:H252	2.50	0.47
1:G:367:LYS:HD3	1:G:480:TYR:CE2	2.50	0.47
1:H:104:ASN:HB3	8:H:1207:HOH:O	2.14	0.47
1:I:531:ASN:HB2	6:I:1009:1PE:H241	1.94	0.47
4:J:1004:1OT:H10	4:J:1004:1OT:O2	2.14	0.47
4:E:1004:1OT:H10	4:E:1004:1OT:O2	2.14	0.47
1:K:117:ILE:HG12	1:K:270:TYR:HB3	1.96	0.47
1:D:156:PHE:CD2	1:D:177:MET:HE1	2.50	0.47
1:H:165:PHE:HB3	1:H:189:TYR:OH	2.14	0.47
7:D:1012:2PE:H121	8:D:1222:HOH:O	2.15	0.47
1:B:321:LEU:HD11	1:B:411:TYR:HA	1.97	0.46
1:D:451:LYS:HG2	6:D:1009:1PE:H251	1.97	0.46
1:A:492:LEU:HD11	4:A:1004:1OT:C7	2.45	0.46
1:C:195:VAL:HG12	1:C:197:ASP:H	1.81	0.46
1:D:364:ASP:O	1:D:420:ASN:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:132:VAL:HG21	1:L:142:VAL:HG13	1.97	0.46
1:L:518:LYS:NZ	6:L:1008:1PE:H131	2.31	0.46
1:F:150:ASP:OD1	1:F:179:ASN:HB2	2.16	0.46
1:C:244:TYR:OH	1:C:588:PRO:O	2.32	0.46
1:C:536:THR:HG21	1:C:551:VAL:HG23	1.97	0.46
1:C:340:ALA:HA	1:C:445:ILE:HD12	1.98	0.46
1:D:177:MET:HB3	8:D:1131:HOH:O	2.16	0.46
1:E:543:ASP:CB	6:E:1009:1PE:H232	2.45	0.46
3:D:1002:CO3:O1	4:D:1004:1OT:O3	2.33	0.45
1:I:172:SER:O	1:I:173:LYS:HD2	2.15	0.45
1:G:321:LEU:HD11	1:G:411:TYR:HA	1.97	0.45
1:F:492:LEU:HD11	4:F:1004:1OT:C7	2.46	0.45
1:J:320:LYS:NZ	6:J:1007:1PE:OH5	2.49	0.45
1:H:135:PRO:HA	1:H:194:SER:O	2.17	0.45
1:E:492:LEU:HD11	4:E:1004:1OT:H12	1.97	0.45
1:A:106:PRO:HD2	1:A:247:MET:HE1	1.99	0.45
1:F:392:MET:HE2	4:F:1004:1OT:C8	2.47	0.45
1:H:122:ASN:HD22	7:H:1007:2PE:H171	1.81	0.45
1:F:292:TYR:O	1:F:296:GLN:HG3	2.17	0.45
1:L:518:LYS:HZ1	6:L:1008:1PE:H131	1.82	0.45
1:G:150:ASP:OD1	1:G:179:ASN:HB2	2.16	0.45
1:C:320:LYS:NZ	6:C:1006:1PE:OH6	2.50	0.45
1:C:150:ASP:OD1	1:C:179:ASN:HB2	2.17	0.44
1:D:411:TYR:CE1	6:D:1007:1PE:H232	2.51	0.44
1:L:321:LEU:HD11	1:L:411:TYR:HA	1.98	0.44
1:A:121:CYS:HA	1:A:270:TYR:CE2	2.53	0.44
1:D:451:LYS:HE3	1:D:564:GLU:O	2.17	0.44
1:F:392:MET:HE1	1:F:583:PHE:CE2	2.52	0.44
1:E:156:PHE:O	1:E:162:MET:HE3	2.17	0.44
1:K:364:ASP:O	1:K:420:ASN:HA	2.18	0.44
1:B:392:MET:HE2	4:B:1004:1OT:C8	2.48	0.44
1:D:451:LYS:HZ2	6:D:1009:1PE:H242	1.82	0.44
1:E:518:LYS:HB2	1:E:518:LYS:HE2	1.56	0.44
1:G:164:LYS:HD3	1:G:165:PHE:CE1	2.52	0.44
1:I:532:GLU:HG3	1:J:205:ARG:HH22	1.82	0.44
4:K:1004:1OT:H10	4:K:1004:1OT:O1	2.18	0.44
1:H:150:ASP:O	1:H:154:SER:HB2	2.17	0.44
1:I:392:MET:HE3	1:I:395:LEU:HD22	1.99	0.44
1:A:383:TYR:HE2	1:A:438:SER:HB2	1.83	0.44
1:B:173:LYS:HB2	1:B:189:TYR:CE1	2.52	0.44
1:B:287:TYR:CD2	1:B:594:ARG:HG2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:328:LEU:HB2	1:H:354:PHE:HB3	1.99	0.43
1:I:103:TYR:HB3	6:I:1008:1PE:H142	2.00	0.43
1:I:320:LYS:HD3	6:I:1007:1PE:H161	2.00	0.43
1:C:135:PRO:HA	1:C:194:SER:O	2.19	0.43
1:C:492:LEU:HD11	4:C:1004:1OT:H12	2.00	0.43
4:C:1004:1OT:H10	4:C:1004:1OT:O2	2.17	0.43
1:H:117:ILE:HD11	1:H:146:SER:OG	2.18	0.43
1:L:451:LYS:HZ3	6:L:1007:1PE:H251	1.83	0.43
1:D:320:LYS:HZ3	7:D:1010:2PE:H182	1.83	0.43
1:E:579:VAL:O	1:E:589:LYS:HD2	2.18	0.43
1:I:326:LYS:HE3	1:I:328:LEU:HD11	2.00	0.43
1:B:381:GLY:HA2	1:B:459:ASP:OD1	2.18	0.43
1:C:96:PRO:HA	7:C:1009:2PE:H172	2.00	0.43
1:D:127:LEU:HB2	1:D:219:LEU:HD22	1.99	0.43
1:E:536:THR:HG21	1:E:551:VAL:CG2	2.47	0.43
1:F:360:LYS:HZ2	1:F:361:SER:HB2	1.83	0.43
4:G:1004:1OT:O2	4:G:1004:1OT:H10	2.19	0.43
1:E:364:ASP:O	1:E:420:ASN:HA	2.19	0.43
1:I:492:LEU:HD11	4:I:1004:1OT:H12	2.01	0.43
1:L:360:LYS:HD2	1:L:361:SER:O	2.18	0.43
1:A:381:GLY:HA2	1:A:459:ASP:OD1	2.18	0.43
1:D:156:PHE:CG	1:D:177:MET:HE1	2.53	0.43
1:A:357:LEU:HB2	1:A:425:PHE:HB2	2.01	0.42
1:B:175:PHE:N	1:B:187:VAL:O	2.41	0.42
1:J:381:GLY:HA2	1:J:459:ASP:OD1	2.19	0.42
1:K:172:SER:OG	1:K:213:MET:HE3	2.19	0.42
1:L:311:SER:O	1:L:315:VAL:HG23	2.19	0.42
1:D:219:LEU:H	1:D:219:LEU:HG	1.30	0.42
1:D:398:PHE:CD2	1:D:398:PHE:C	2.95	0.42
1:J:219:LEU:H	1:J:219:LEU:HG	1.16	0.42
1:I:398:PHE:CD2	1:I:398:PHE:C	2.98	0.42
1:I:492:LEU:HD23	1:I:492:LEU:HA	1.82	0.42
1:D:198:LEU:HD22	1:D:202:ASP:HB3	2.02	0.42
1:L:498:SER:O	1:L:523:PRO:HG2	2.19	0.42
1:D:321:LEU:HD11	1:D:411:TYR:HA	2.01	0.42
1:F:162:MET:HE2	1:F:162:MET:HA	2.02	0.42
1:D:383:TYR:HE2	1:D:438:SER:HB2	1.85	0.42
1:F:386:LYS:NZ	4:F:1004:1OT:O2	2.41	0.42
4:I:1004:1OT:O2	4:I:1004:1OT:H10	2.20	0.42
1:C:173:LYS:HB2	1:C:189:TYR:CE1	2.55	0.42
1:D:156:PHE:O	1:D:162:MET:HE3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:176:TYR:OH	1:D:217:ASN:OD1	2.20	0.42
1:D:381:GLY:HA2	1:D:459:ASP:OD1	2.20	0.42
1:F:360:LYS:HA	1:F:360:LYS:HD2	1.91	0.42
1:G:411:TYR:CE1	6:G:1008:1PE:H132	2.50	0.42
1:I:364:ASP:O	1:I:420:ASN:HA	2.19	0.42
1:L:200:GLU:HG3	8:L:1158:HOH:O	2.19	0.42
1:C:230:VAL:HG12	1:C:234:LEU:HD23	2.02	0.42
1:D:451:LYS:CG	6:D:1009:1PE:H251	2.50	0.42
1:D:597:THR:HG22	1:D:601:LEU:HD22	2.02	0.42
1:H:174:HIS:HB3	1:L:175:PHE:CD1	2.55	0.42
1:J:392:MET:HE1	1:J:583:PHE:CE2	2.54	0.42
1:E:135:PRO:HA	1:E:194:SER:O	2.20	0.41
1:E:543:ASP:CG	6:E:1009:1PE:H232	2.44	0.41
3:L:1002:CO3:O1	4:L:1004:1OT:O3	2.38	0.41
1:F:341:TYR:CE1	1:F:428:ALA:HB1	2.55	0.41
1:G:488:THR:HG21	1:G:555:SER:HA	2.02	0.41
1:J:332:GLU:HG3	8:J:1209:HOH:O	2.20	0.41
1:D:132:VAL:HG21	1:D:144:ILE:HD13	2.03	0.41
1:F:360:LYS:HD2	1:F:361:SER:N	2.34	0.41
1:J:398:PHE:CD2	1:J:398:PHE:C	2.99	0.41
1:J:451:LYS:HE3	6:J:1008:1PE:H151	2.02	0.41
1:D:395:LEU:HD11	1:D:581:TRP:CG	2.56	0.41
1:K:514:LEU:O	1:K:518:LYS:HG2	2.20	0.41
1:L:198:LEU:HD12	1:L:198:LEU:HA	1.84	0.41
1:I:158:LYS:HB2	8:I:1103:HOH:O	2.20	0.41
1:E:152:GLN:HG2	1:E:180:ASP:OD1	2.20	0.41
1:E:340:ALA:HA	1:E:445:ILE:HD12	2.02	0.41
1:L:451:LYS:HZ3	6:L:1007:1PE:H141	1.86	0.41
1:L:451:LYS:HE2	6:L:1007:1PE:H251	2.03	0.41
1:A:372:VAL:O	1:A:483:ASP:HA	2.21	0.41
1:D:102:GLU:HG2	1:D:105:THR:HG22	2.03	0.41
1:G:357:LEU:HB2	1:G:425:PHE:HB2	2.03	0.41
1:K:102:GLU:HG2	1:K:105:THR:HG22	2.02	0.41
1:A:114:VAL:HG12	1:A:274:ALA:HB1	2.03	0.41
7:C:1009:2PE:H141	8:C:1146:HOH:O	2.20	0.41
1:L:127:LEU:HD11	1:L:129:ILE:HD11	2.03	0.41
1:B:551:VAL:HG12	1:B:553:ALA:H	1.87	0.40
1:D:316:GLU:HG2	1:D:320:LYS:HE2	2.02	0.40
1:I:236:ARG:HH11	1:I:236:ARG:HD3	1.75	0.40
1:B:165:PHE:CE2	1:B:173:LYS:HG2	2.56	0.40
1:F:300:ALA:HA	1:F:301:PRO:HD3	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:214:LEU:HD21	1:J:222:LEU:HD22	2.04	0.40
1:J:579:VAL:O	1:J:589:LYS:HD2	2.22	0.40
4:L:1004:1OT:O1	4:L:1004:1OT:H10	2.22	0.40
1:A:392:MET:HE1	1:A:583:PHE:HE2	1.86	0.40
1:C:198:LEU:HD22	1:C:202:ASP:HB3	2.04	0.40
1:F:311:SER:HB2	1:F:327:ILE:HD12	2.03	0.40
1:G:492:LEU:HD23	1:G:492:LEU:HA	1.89	0.40
1:H:488:THR:HG21	1:H:555:SER:HA	2.03	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	518/528 (98%)	507 (98%)	11 (2%)	0	100	100
1	B	516/528 (98%)	503 (98%)	11 (2%)	2 (0%)	30	27
1	C	517/528 (98%)	505 (98%)	12 (2%)	0	100	100
1	D	510/528 (97%)	498 (98%)	12 (2%)	0	100	100
1	E	506/528 (96%)	495 (98%)	11 (2%)	0	100	100
1	F	507/528 (96%)	497 (98%)	10 (2%)	0	100	100
1	G	517/528 (98%)	507 (98%)	10 (2%)	0	100	100
1	H	512/528 (97%)	502 (98%)	9 (2%)	1 (0%)	43	42
1	I	516/528 (98%)	505 (98%)	11 (2%)	0	100	100
1	J	509/528 (96%)	498 (98%)	11 (2%)	0	100	100
1	K	503/528 (95%)	494 (98%)	8 (2%)	1 (0%)	43	42
1	L	509/528 (96%)	500 (98%)	8 (2%)	1 (0%)	43	42
All	All	6140/6336 (97%)	6011 (98%)	124 (2%)	5 (0%)	48	46

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	138	GLU
1	L	137	LYS
1	B	257	LYS
1	B	258	ASN
1	K	363	GLY

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	429/455 (94%)	420 (98%)	9 (2%)	47	52
1	B	415/455 (91%)	402 (97%)	13 (3%)	35	37
1	C	421/455 (92%)	402 (96%)	19 (4%)	24	23
1	D	421/455 (92%)	414 (98%)	7 (2%)	53	60
1	E	416/455 (91%)	408 (98%)	8 (2%)	50	56
1	F	401/455 (88%)	390 (97%)	11 (3%)	39	42
1	G	428/455 (94%)	416 (97%)	12 (3%)	38	41
1	H	411/455 (90%)	394 (96%)	17 (4%)	27	26
1	I	423/455 (93%)	413 (98%)	10 (2%)	43	47
1	J	420/455 (92%)	411 (98%)	9 (2%)	47	52
1	K	417/455 (92%)	409 (98%)	8 (2%)	50	56
1	L	407/455 (90%)	392 (96%)	15 (4%)	30	30
All	All	5009/5460 (92%)	4871 (97%)	138 (3%)	38	41

All (138) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	169	LEU
1	A	200	GLU
1	A	204	LYS
1	A	208	LEU

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Mol	Chain	Res	Type
1	A	398	PHE
1	A	439	TYR
1	A	549	SER
1	A	552	LYS
1	A	603	ASP
1	B	125	GLU
1	B	169	LEU
1	B	200	GLU
1	B	208	LEU
1	B	215	HIS
1	B	255	THR
1	B	261	MET
1	B	328	LEU
1	B	357	LEU
1	B	398	PHE
1	B	421	VAL
1	B	439	TYR
1	B	603	ASP
1	C	86	SER
1	C	132	VAL
1	C	163	GLU
1	C	169	LEU
1	C	184	SER
1	C	185	VAL
1	C	200	GLU
1	C	208	LEU
1	C	288	TYR
1	C	322	ASN
1	C	357	LEU
1	C	398	PHE
1	C	421	VAL
1	C	436	LYS
1	C	439	TYR
1	C	483	ASP
1	C	549	SER
1	C	601	LEU
1	C	603	ASP
1	D	124	GLU
1	D	154	SER
1	D	219	LEU
1	D	322	ASN
1	D	398	PHE

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Mol	Chain	Res	Type
1	D	439	TYR
1	D	552	LYS
1	E	86	SER
1	E	169	LEU
1	E	195	VAL
1	E	208	LEU
1	E	398	PHE
1	E	436	LYS
1	E	439	TYR
1	E	567	GLN
1	F	86	SER
1	F	198	LEU
1	F	200	GLU
1	F	208	LEU
1	F	225	VAL
1	F	229	ASN
1	F	254	SER
1	F	326	LYS
1	F	364	ASP
1	F	398	PHE
1	F	439	TYR
1	G	117	ILE
1	G	164	LYS
1	G	169	LEU
1	G	200	GLU
1	G	204	LYS
1	G	208	LEU
1	G	272	ASN
1	G	367	LYS
1	G	398	PHE
1	G	436	LYS
1	G	439	TYR
1	G	603	ASP
1	H	86	SER
1	H	125	GLU
1	H	154	SER
1	H	164	LYS
1	H	173	LYS
1	H	181	ASN
1	H	200	GLU
1	H	208	LEU
1	H	215	HIS

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Mol	Chain	Res	Type
1	H	232	LYS
1	H	254	SER
1	H	357	LEU
1	H	398	PHE
1	H	421	VAL
1	H	439	TYR
1	H	601	LEU
1	H	603	ASP
1	I	86	SER
1	I	168	LYS
1	I	169	LEU
1	I	208	LEU
1	I	212	THR
1	I	322	ASN
1	I	398	PHE
1	I	421	VAL
1	I	439	TYR
1	I	549	SER
1	J	147	LYS
1	J	164	LYS
1	J	200	GLU
1	J	205	ARG
1	J	219	LEU
1	J	262	GLU
1	J	398	PHE
1	J	439	TYR
1	J	518	LYS
1	K	86	SER
1	K	132	VAL
1	K	145	SER
1	K	164	LYS
1	K	169	LEU
1	K	216	ASP
1	K	398	PHE
1	K	439	TYR
1	L	86	SER
1	L	116	ASP
1	L	169	LEU
1	L	198	LEU
1	L	200	GLU
1	L	254	SER
1	L	295	SER

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Mol	Chain	Res	Type
1	L	326	LYS
1	L	398	PHE
1	L	421	VAL
1	L	436	LYS
1	L	439	TYR
1	L	567	GLN
1	L	568	ASN
1	L	603	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	420	ASN
1	B	113	GLN
1	B	139	ASN
1	B	215	HIS
1	B	531	ASN
1	C	260	ASN
1	C	420	ASN
1	D	183	ASN
1	D	272	ASN
1	D	420	ASN
1	E	139	ASN
1	E	420	ASN
1	F	161	ASN
1	F	229	ASN
1	F	531	ASN
1	H	515	GLN
1	I	183	ASN
1	J	183	ASN
1	J	531	ASN
1	K	420	ASN
1	K	602	ASN

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 115 ligands modelled in this entry, 24 are monoatomic - leaving 91 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	1PE	F	1008	-	9,9,15	0.47	0	8,8,14	0.18	0
4	1OT	F	1004	2	17,18,18	2.83	7 (41%)	23,26,26	2.20	7 (30%)
7	2PE	B	1006	-	25,25,27	0.18	0	24,24,26	0.17	0
6	1PE	G	1008	-	8,8,15	0.41	0	7,7,14	0.36	0
3	CO3	D	1002	-	3,3,3	0.84	0	2,3,3	0.29	0
4	1OT	A	1004	2	17,18,18	2.83	5 (29%)	23,26,26	2.38	9 (39%)
3	CO3	E	1002	-	3,3,3	0.68	0	2,3,3	0.08	0
3	CO3	C	1002	-	3,3,3	0.75	0	2,3,3	0.17	0
6	1PE	K	1007	-	11,11,15	0.52	0	10,10,14	0.44	0
6	1PE	J	1008	-	10,10,15	0.55	0	9,9,14	0.26	0
6	1PE	K	1009	-	10,10,15	0.48	0	9,9,14	0.15	0
7	2PE	B	1007	-	9,9,27	0.06	0	8,8,26	0.23	0
6	1PE	K	1010	-	5,5,15	0.47	0	4,4,14	0.16	0
6	1PE	G	1011	-	14,14,15	0.41	0	13,13,14	0.56	0
5	SO4	G	1007	-	4,4,4	0.25	0	6,6,6	0.19	0
6	1PE	L	1007	-	11,11,15	0.52	0	10,10,14	0.19	0
6	1PE	C	1008	-	6,6,15	0.50	0	5,5,14	0.27	0
5	SO4	I	1006	-	4,4,4	0.29	0	6,6,6	0.12	0
5	SO4	H	1005	-	4,4,4	0.24	0	6,6,6	0.21	0
6	1PE	J	1006	-	10,10,15	0.43	0	9,9,14	0.43	0
5	SO4	D	1006	-	4,4,4	0.22	0	6,6,6	0.08	0
6	1PE	I	1009	-	4,4,15	0.39	0	3,3,14	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	1PE	C	1006	-	12,12,15	0.55	0	11,11,14	0.19	0
6	1PE	E	1008	-	11,11,15	0.49	0	10,10,14	0.31	0
3	CO3	F	1002	-	3,3,3	0.79	0	2,3,3	0.17	0
6	1PE	D	1009	-	10,10,15	0.53	0	9,9,14	0.40	0
6	1PE	E	1009	-	7,7,15	0.44	0	6,6,14	0.36	0
5	SO4	E	1005	-	4,4,4	0.21	0	6,6,6	0.12	0
5	SO4	F	1005	-	4,4,4	0.21	0	6,6,6	0.09	0
6	1PE	A	1010	-	11,11,15	0.51	0	10,10,14	0.15	0
3	CO3	G	1002	-	3,3,3	0.93	0	2,3,3	0.31	0
3	CO3	J	1002	-	3,3,3	0.70	0	2,3,3	0.36	0
4	1OT	H	1004	2	17,18,18	2.85	5 (29%)	23,26,26	2.09	7 (30%)
5	SO4	A	1008	-	4,4,4	0.28	0	6,6,6	0.27	0
5	SO4	G	1006	-	4,4,4	0.26	0	6,6,6	0.28	0
5	SO4	A	1005	-	4,4,4	0.21	0	6,6,6	0.24	0
6	1PE	J	1007	-	9,9,15	0.52	0	8,8,14	0.32	0
3	CO3	H	1002	-	3,3,3	0.71	0	2,3,3	0.45	0
4	1OT	K	1004	2	17,18,18	2.72	5 (29%)	23,26,26	2.21	9 (39%)
4	1OT	D	1004	2	17,18,18	2.90	5 (29%)	23,26,26	2.50	8 (34%)
5	SO4	L	1005	-	4,4,4	0.21	0	6,6,6	0.13	0
7	2PE	D	1012	-	8,8,27	0.16	0	7,7,26	0.16	0
7	2PE	J	1010	-	9,9,27	0.20	0	8,8,26	0.09	0
7	2PE	H	1006	-	24,24,27	0.20	0	23,23,26	0.15	0
6	1PE	I	1007	-	14,14,15	0.39	0	13,13,14	0.49	0
5	SO4	J	1005	-	4,4,4	0.21	0	6,6,6	0.07	0
7	2PE	D	1011	-	4,4,27	0.24	0	3,3,26	0.37	0
5	SO4	C	1005	-	4,4,4	0.20	0	6,6,6	0.20	0
6	1PE	F	1007	-	9,9,15	0.48	0	8,8,14	0.14	0
5	SO4	K	1005	-	4,4,4	0.34	0	6,6,6	0.16	0
4	1OT	E	1004	2	17,18,18	2.86	5 (29%)	23,26,26	2.35	7 (30%)
6	1PE	J	1009	-	8,8,15	0.55	0	7,7,14	0.39	0
7	2PE	H	1007	-	9,9,27	0.19	0	8,8,26	0.08	0
3	CO3	K	1002	-	3,3,3	0.76	0	2,3,3	0.13	0
6	1PE	K	1008	-	11,11,15	0.46	0	10,10,14	0.20	0
6	1PE	D	1008	-	8,8,15	0.50	0	7,7,14	0.22	0
7	2PE	F	1010	-	8,8,27	0.19	0	7,7,26	0.22	0
6	1PE	G	1009	-	5,5,15	0.45	0	4,4,14	0.49	0
6	1PE	D	1007	-	9,9,15	0.44	0	8,8,14	0.44	0
6	1PE	E	1007	-	11,11,15	0.48	0	10,10,14	0.32	0
3	CO3	B	1002	-	3,3,3	0.81	0	2,3,3	0.28	0
4	1OT	C	1004	2	17,18,18	2.76	6 (35%)	23,26,26	2.15	7 (30%)
5	SO4	I	1005	-	4,4,4	0.26	0	6,6,6	0.10	0
6	1PE	G	1010	-	5,5,15	0.49	0	4,4,14	0.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	1PE	F	1009	-	10,10,15	0.48	0	9,9,14	0.21	0
5	SO4	D	1005	-	4,4,4	0.39	0	6,6,6	0.15	0
7	2PE	E	1010	-	8,8,27	0.12	0	7,7,26	0.18	0
5	SO4	A	1006	-	4,4,4	0.23	0	6,6,6	0.20	0
6	1PE	L	1008	-	7,7,15	0.46	0	6,6,14	0.38	0
5	SO4	A	1007	-	4,4,4	0.25	0	6,6,6	0.29	0
6	1PE	L	1009	-	10,10,15	0.50	0	9,9,14	0.49	0
3	CO3	I	1002	-	3,3,3	0.63	0	2,3,3	0.34	0
7	2PE	I	1010	-	7,7,27	0.14	0	6,6,26	0.11	0
4	1OT	L	1004	2	17,18,18	2.81	5 (29%)	23,26,26	2.34	8 (34%)
6	1PE	F	1006	-	9,9,15	0.44	0	8,8,14	0.24	0
5	SO4	K	1006	-	4,4,4	0.27	0	6,6,6	0.09	0
3	CO3	L	1002	-	3,3,3	0.91	0	2,3,3	0.21	0
6	1PE	C	1007	-	8,8,15	0.42	0	7,7,14	0.22	0
5	SO4	B	1005	-	4,4,4	0.26	0	6,6,6	0.06	0
5	SO4	E	1006	-	4,4,4	0.28	0	6,6,6	0.27	0
4	1OT	G	1004	2	17,18,18	2.79	5 (29%)	23,26,26	2.14	6 (26%)
4	1OT	J	1004	2	17,18,18	2.81	6 (35%)	23,26,26	2.23	7 (30%)
6	1PE	L	1006	-	9,9,15	0.37	0	8,8,14	0.68	0
7	2PE	C	1009	-	8,8,27	0.14	0	7,7,26	0.24	0
7	2PE	D	1010	-	12,12,27	0.18	0	11,11,26	0.14	0
4	1OT	B	1004	2	17,18,18	2.84	5 (29%)	23,26,26	2.22	7 (30%)
3	CO3	A	1002	-	3,3,3	0.88	0	2,3,3	0.25	0
6	1PE	A	1009	-	8,8,15	0.43	0	7,7,14	0.26	0
5	SO4	G	1005	-	4,4,4	0.28	0	6,6,6	0.27	0
6	1PE	I	1008	-	10,10,15	0.46	0	9,9,14	0.33	0
4	1OT	I	1004	2	17,18,18	2.70	6 (35%)	23,26,26	2.24	10 (43%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	1PE	I	1009	-	-	1/2/2/13	-
6	1PE	L	1008	-	-	3/5/5/13	-
6	1PE	F	1007	-	-	4/7/7/13	-
6	1PE	C	1006	-	-	1/10/10/13	-
6	1PE	F	1008	-	-	2/7/7/13	-
6	1PE	E	1008	-	-	3/9/9/13	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	1OT	F	1004	2	-	2/13/14/14	0/2/2/2
6	1PE	L	1009	-	-	4/8/8/13	-
7	2PE	B	1006	-	-	9/23/23/25	-
4	1OT	E	1004	2	-	2/13/14/14	0/2/2/2
6	1PE	D	1009	-	-	7/8/8/13	-
7	2PE	I	1010	-	-	1/5/5/25	-
4	1OT	L	1004	2	-	2/13/14/14	0/2/2/2
6	1PE	F	1006	-	-	1/7/7/13	-
6	1PE	J	1009	-	-	3/6/6/13	-
6	1PE	E	1009	-	-	3/5/5/13	-
6	1PE	A	1010	-	-	5/9/9/13	-
6	1PE	G	1008	-	-	4/6/6/13	-
7	2PE	H	1007	-	-	3/7/7/25	-
6	1PE	K	1008	-	-	2/9/9/13	-
4	1OT	A	1004	2	-	3/13/14/14	0/2/2/2
4	1OT	H	1004	2	-	3/13/14/14	0/2/2/2
6	1PE	D	1008	-	-	3/6/6/13	-
6	1PE	K	1007	-	-	3/9/9/13	-
6	1PE	C	1007	-	-	4/6/6/13	-
7	2PE	F	1010	-	-	3/6/6/25	-
6	1PE	J	1008	-	-	3/8/8/13	-
6	1PE	G	1009	-	-	1/3/3/13	-
6	1PE	D	1007	-	-	4/7/7/13	-
6	1PE	K	1009	-	-	1/8/8/13	-
6	1PE	J	1007	-	-	0/7/7/13	-
6	1PE	E	1007	-	-	4/9/9/13	-
4	1OT	C	1004	2	-	3/13/14/14	0/2/2/2
4	1OT	G	1004	2	-	2/13/14/14	0/2/2/2
4	1OT	J	1004	2	-	2/13/14/14	0/2/2/2
4	1OT	K	1004	2	-	3/13/14/14	0/2/2/2
6	1PE	L	1006	-	-	3/7/7/13	-
7	2PE	C	1009	-	-	3/6/6/25	-
4	1OT	D	1004	2	-	2/13/14/14	0/2/2/2
7	2PE	D	1012	-	-	3/6/6/25	-
7	2PE	B	1007	-	-	4/7/7/25	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	1PE	G	1010	-	-	3/3/3/13	-
7	2PE	J	1010	-	-	5/7/7/25	-
6	1PE	K	1010	-	-	2/3/3/13	-
7	2PE	D	1010	-	-	7/10/10/25	-
4	1OT	B	1004	2	-	3/13/14/14	0/2/2/2
6	1PE	F	1009	-	-	2/8/8/13	-
7	2PE	H	1006	-	-	7/22/22/25	-
6	1PE	I	1007	-	-	5/12/12/13	-
6	1PE	G	1011	-	-	6/12/12/13	-
6	1PE	L	1007	-	-	6/9/9/13	-
6	1PE	C	1008	-	-	2/4/4/13	-
7	2PE	D	1011	-	-	2/2/2/25	-
6	1PE	A	1009	-	-	0/6/6/13	-
7	2PE	E	1010	-	-	2/6/6/25	-
6	1PE	I	1008	-	-	3/8/8/13	-
4	1OT	I	1004	2	-	3/13/14/14	0/2/2/2
6	1PE	J	1006	-	-	6/8/8/13	-

All (65) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1004	1OT	P-O3	7.77	1.62	1.49
4	F	1004	1OT	P-O3	7.58	1.62	1.49
4	J	1004	1OT	P-O3	7.56	1.62	1.49
4	H	1004	1OT	P-O3	7.55	1.62	1.49
4	E	1004	1OT	P-O3	7.53	1.62	1.49
4	C	1004	1OT	P-O3	7.41	1.62	1.49
4	B	1004	1OT	P-O3	7.41	1.62	1.49
4	G	1004	1OT	P-O3	7.32	1.61	1.49
4	A	1004	1OT	P-O3	7.26	1.61	1.49
4	L	1004	1OT	P-O3	7.06	1.61	1.49
4	K	1004	1OT	P-O3	6.95	1.61	1.49
4	I	1004	1OT	P-O3	6.80	1.61	1.49
4	L	1004	1OT	C5-N2	-5.27	1.33	1.43
4	E	1004	1OT	C5-N2	-5.24	1.33	1.43
4	A	1004	1OT	C5-N2	-5.07	1.33	1.43
4	B	1004	1OT	C5-N2	-5.03	1.33	1.43
4	J	1004	1OT	C5-N2	-5.03	1.33	1.43
4	D	1004	1OT	C5-N2	-5.02	1.33	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	1004	1OT	C5-N2	-5.01	1.33	1.43
4	F	1004	1OT	C5-N2	-5.01	1.33	1.43
4	K	1004	1OT	C5-N2	-5.01	1.33	1.43
4	H	1004	1OT	C5-N2	-4.99	1.33	1.43
4	C	1004	1OT	C5-N2	-4.85	1.33	1.43
4	D	1004	1OT	P-O2	4.78	1.62	1.54
4	I	1004	1OT	C5-N2	-4.73	1.34	1.43
4	L	1004	1OT	P-O2	4.59	1.62	1.54
4	K	1004	1OT	P-O2	4.44	1.61	1.54
4	F	1004	1OT	P-O2	4.33	1.61	1.54
4	E	1004	1OT	P-O2	4.31	1.61	1.54
4	G	1004	1OT	P-O2	4.30	1.61	1.54
4	A	1004	1OT	P-O2	4.30	1.61	1.54
4	B	1004	1OT	P-O2	4.27	1.61	1.54
4	I	1004	1OT	P-O2	4.24	1.61	1.54
4	C	1004	1OT	P-O2	4.22	1.61	1.54
4	B	1004	1OT	P-O1	-4.15	1.48	1.54
4	H	1004	1OT	P-O1	-4.11	1.48	1.54
4	H	1004	1OT	P-O2	3.98	1.61	1.54
4	I	1004	1OT	P-O1	-3.95	1.48	1.54
4	E	1004	1OT	P-O1	-3.94	1.48	1.54
4	J	1004	1OT	P-O2	3.93	1.61	1.54
4	A	1004	1OT	P-O1	-3.86	1.48	1.54
4	C	1004	1OT	P-O1	-3.70	1.49	1.54
4	L	1004	1OT	P-O1	-3.69	1.49	1.54
4	A	1004	1OT	C2-C1	-3.64	1.47	1.52
4	K	1004	1OT	C2-C1	-3.63	1.47	1.52
4	H	1004	1OT	C2-C1	-3.60	1.47	1.52
4	G	1004	1OT	C2-C1	-3.57	1.47	1.52
4	F	1004	1OT	P-O1	-3.54	1.49	1.54
4	J	1004	1OT	C2-C1	-3.46	1.47	1.52
4	I	1004	1OT	C2-C1	-3.44	1.47	1.52
4	G	1004	1OT	P-O1	-3.43	1.49	1.54
4	L	1004	1OT	C2-C1	-3.39	1.48	1.52
4	D	1004	1OT	P-O1	-3.31	1.49	1.54
4	J	1004	1OT	P-O1	-3.27	1.49	1.54
4	D	1004	1OT	C2-C1	-3.26	1.48	1.52
4	E	1004	1OT	C2-C1	-3.23	1.48	1.52
4	B	1004	1OT	C2-C1	-3.22	1.48	1.52
4	K	1004	1OT	P-O1	-3.07	1.50	1.54
4	F	1004	1OT	C2-C1	-3.00	1.48	1.52
4	C	1004	1OT	C2-C1	-2.93	1.48	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	1004	1OT	N2-N3	2.27	1.40	1.36
4	J	1004	1OT	N2-N3	2.15	1.39	1.36
4	C	1004	1OT	N2-N3	2.11	1.39	1.36
4	I	1004	1OT	N2-N3	2.04	1.39	1.36
4	F	1004	1OT	C6-N2	-2.01	1.32	1.35

All (92) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	1004	1OT	C6-N2-N3	5.07	116.06	111.83
4	L	1004	1OT	C7-C6-N2	-5.01	102.63	106.76
4	E	1004	1OT	C6-N2-N3	4.99	115.99	111.83
4	K	1004	1OT	C7-C6-N2	-4.93	102.70	106.76
4	F	1004	1OT	C7-C6-N2	-4.92	102.71	106.76
4	G	1004	1OT	C7-C6-N2	-4.90	102.72	106.76
4	J	1004	1OT	C6-N2-N3	4.81	115.84	111.83
4	B	1004	1OT	C7-C6-N2	-4.79	102.82	106.76
4	A	1004	1OT	C7-C6-N2	-4.79	102.82	106.76
4	J	1004	1OT	C7-C6-N2	-4.78	102.82	106.76
4	H	1004	1OT	C7-C6-N2	-4.75	102.85	106.76
4	D	1004	1OT	C4-C5-N2	4.74	125.38	119.58
4	C	1004	1OT	C7-C6-N2	-4.73	102.86	106.76
4	I	1004	1OT	C7-C6-N2	-4.73	102.87	106.76
4	E	1004	1OT	C7-C6-N2	-4.68	102.91	106.76
4	K	1004	1OT	C6-N2-N3	4.65	115.71	111.83
4	A	1004	1OT	C6-N2-N3	4.61	115.67	111.83
4	F	1004	1OT	C6-N2-N3	4.59	115.66	111.83
4	D	1004	1OT	C10-C2-C1	-4.50	115.65	120.76
4	C	1004	1OT	C6-N2-N3	4.48	115.57	111.83
4	G	1004	1OT	C6-N2-N3	4.44	115.53	111.83
4	B	1004	1OT	C6-N2-N3	4.43	115.52	111.83
4	H	1004	1OT	C6-N2-N3	4.40	115.50	111.83
4	D	1004	1OT	C6-N2-N3	4.21	115.34	111.83
4	I	1004	1OT	C6-N2-N3	4.11	115.26	111.83
4	D	1004	1OT	C7-C6-N2	-4.08	103.40	106.76
4	A	1004	1OT	C4-C5-N2	4.04	124.53	119.58
4	A	1004	1OT	C10-C2-C1	-3.90	116.34	120.76
4	I	1004	1OT	C4-C5-N2	3.67	124.07	119.58
4	E	1004	1OT	C4-C5-N2	3.67	124.07	119.58
4	E	1004	1OT	C10-C2-C1	-3.56	116.72	120.76
4	D	1004	1OT	C5-N2-C6	-3.52	120.72	128.39
4	D	1004	1OT	C3-C2-C1	3.50	124.73	120.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1004	1OT	C4-C5-N2	3.49	123.85	119.58
4	E	1004	1OT	C5-N2-C6	-3.48	120.80	128.39
4	J	1004	1OT	C10-C2-C1	-3.45	116.85	120.76
4	A	1004	1OT	C5-N2-C6	-3.43	120.92	128.39
4	L	1004	1OT	C10-C2-C1	-3.43	116.87	120.76
4	B	1004	1OT	C5-N2-C6	-3.32	121.14	128.39
4	L	1004	1OT	C4-C5-N2	3.31	123.63	119.58
4	F	1004	1OT	C5-N2-C6	-3.28	121.24	128.39
4	J	1004	1OT	C5-N2-C6	-3.27	121.27	128.39
4	J	1004	1OT	C4-C5-N2	3.21	123.51	119.58
4	I	1004	1OT	C5-N2-C6	-3.18	121.45	128.39
4	L	1004	1OT	C5-N2-C6	-3.17	121.47	128.39
4	C	1004	1OT	C5-N2-C6	-3.16	121.50	128.39
4	I	1004	1OT	C10-C2-C1	-3.15	117.18	120.76
4	K	1004	1OT	C5-N2-C6	-3.11	121.61	128.39
4	H	1004	1OT	C5-N2-C6	-3.07	121.70	128.39
4	C	1004	1OT	C4-C5-N2	3.04	123.30	119.58
4	F	1004	1OT	C10-C2-C1	-3.00	117.35	120.76
4	B	1004	1OT	C10-C2-C1	-2.95	117.41	120.76
4	F	1004	1OT	C4-C5-N2	2.90	123.13	119.58
4	G	1004	1OT	C5-N2-C6	-2.86	122.16	128.39
4	G	1004	1OT	C10-C2-C1	-2.82	117.56	120.76
4	H	1004	1OT	C4-C5-N2	2.73	122.93	119.58
4	C	1004	1OT	O1-P-O3	-2.65	106.81	113.45
4	G	1004	1OT	C4-C5-N2	2.56	122.72	119.58
4	K	1004	1OT	C4-C5-N2	2.56	122.71	119.58
4	K	1004	1OT	C10-C2-C1	-2.52	117.89	120.76
4	H	1004	1OT	C10-C2-C1	-2.48	117.94	120.76
4	C	1004	1OT	C10-C2-C1	-2.46	117.97	120.76
4	D	1004	1OT	C4-C3-C2	-2.45	118.74	121.18
4	E	1004	1OT	C3-C2-C1	2.33	123.40	120.76
4	F	1004	1OT	C6-C7-C8	2.31	109.96	105.44
4	G	1004	1OT	C6-C7-C8	2.27	109.89	105.44
4	K	1004	1OT	C6-C7-C8	2.22	109.78	105.44
4	I	1004	1OT	O1-P-O3	-2.21	107.90	113.45
4	B	1004	1OT	C3-C2-C1	2.21	123.27	120.76
4	B	1004	1OT	C6-C7-C8	2.21	109.76	105.44
4	A	1004	1OT	C3-C2-C1	2.18	123.23	120.76
4	H	1004	1OT	C6-C7-C8	2.18	109.70	105.44
4	K	1004	1OT	O2-P-O3	-2.16	108.04	113.45
4	I	1004	1OT	C3-C2-C1	2.16	123.20	120.76
4	H	1004	1OT	C10-C9-C5	2.13	123.00	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	1004	1OT	C3-C2-C1	2.13	123.17	120.76
4	E	1004	1OT	O1-P-O3	-2.13	108.11	113.45
4	A	1004	1OT	C6-C7-C8	2.12	109.59	105.44
4	L	1004	1OT	O1-P-O3	-2.11	108.15	113.45
4	J	1004	1OT	C3-C2-C1	2.11	123.15	120.76
4	I	1004	1OT	P-C1-C2	2.11	115.63	111.63
4	I	1004	1OT	C6-C7-C8	2.11	109.56	105.44
4	I	1004	1OT	C4-C3-C2	-2.11	119.08	121.18
4	J	1004	1OT	C6-C7-C8	2.10	109.55	105.44
4	A	1004	1OT	C4-C3-C2	-2.10	119.09	121.18
4	K	1004	1OT	P-C1-C2	2.08	115.57	111.63
4	K	1004	1OT	C4-C3-C2	-2.08	119.11	121.18
4	C	1004	1OT	C6-C7-C8	2.07	109.48	105.44
4	L	1004	1OT	C6-C7-C8	2.06	109.47	105.44
4	L	1004	1OT	C3-C2-C1	2.05	123.08	120.76
4	D	1004	1OT	P-C1-C2	2.04	115.49	111.63
4	A	1004	1OT	C10-C9-C5	2.03	122.87	120.30

There are no chirality outliers.

All (183) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	1004	1OT	P-C1-C2-C10
4	D	1004	1OT	P-C1-C2-C3
4	E	1004	1OT	P-C1-C2-C10
4	F	1004	1OT	P-C1-C2-C10
4	F	1004	1OT	P-C1-C2-C3
4	J	1004	1OT	P-C1-C2-C10
4	J	1004	1OT	P-C1-C2-C3
4	L	1004	1OT	P-C1-C2-C10
6	K	1010	1PE	OH5-C14-C24-OH4
7	H	1006	2PE	O16-C17-C18-O19
6	F	1007	1PE	OH6-C15-C25-OH5
6	G	1011	1PE	OH6-C15-C25-OH5
6	L	1009	1PE	OH6-C15-C25-OH5
6	G	1011	1PE	OH5-C14-C24-OH4
6	I	1007	1PE	OH4-C13-C23-OH3
6	I	1008	1PE	OH4-C13-C23-OH3
4	A	1004	1OT	P-C1-C2-C10
4	A	1004	1OT	P-C1-C2-C3
4	B	1004	1OT	P-C1-C2-C10
4	B	1004	1OT	P-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
4	C	1004	1OT	P-C1-C2-C10
4	C	1004	1OT	P-C1-C2-C3
4	E	1004	1OT	P-C1-C2-C3
4	G	1004	1OT	P-C1-C2-C10
4	G	1004	1OT	P-C1-C2-C3
4	H	1004	1OT	P-C1-C2-C10
4	H	1004	1OT	P-C1-C2-C3
4	I	1004	1OT	P-C1-C2-C10
4	I	1004	1OT	P-C1-C2-C3
4	K	1004	1OT	P-C1-C2-C10
4	K	1004	1OT	P-C1-C2-C3
4	L	1004	1OT	P-C1-C2-C3
6	I	1007	1PE	OH6-C15-C25-OH5
7	C	1009	2PE	C14-C15-O16-C17
7	F	1010	2PE	C11-C12-O13-C14
6	D	1007	1PE	OH5-C14-C24-OH4
6	L	1007	1PE	OH5-C14-C24-OH4
7	D	1011	2PE	C14-C15-O16-C17
6	I	1009	1PE	OH5-C14-C24-OH4
7	J	1010	2PE	O13-C14-C15-O16
6	L	1006	1PE	OH6-C15-C25-OH5
6	D	1007	1PE	OH4-C13-C23-OH3
6	E	1009	1PE	OH4-C13-C23-OH3
6	F	1007	1PE	OH5-C14-C24-OH4
6	L	1007	1PE	OH6-C15-C25-OH5
7	D	1011	2PE	O13-C14-C15-O16
7	D	1010	2PE	O10-C11-C12-O13
6	F	1008	1PE	OH5-C14-C24-OH4
6	G	1010	1PE	OH6-C15-C25-OH5
6	C	1007	1PE	OH4-C13-C23-OH3
6	C	1008	1PE	OH5-C14-C24-OH4
6	J	1008	1PE	OH6-C15-C25-OH5
6	L	1006	1PE	OH7-C16-C26-OH6
6	D	1009	1PE	OH6-C15-C25-OH5
6	F	1009	1PE	OH4-C13-C23-OH3
6	K	1007	1PE	OH6-C15-C25-OH5
6	J	1006	1PE	OH5-C14-C24-OH4
6	J	1008	1PE	OH5-C14-C24-OH4
6	G	1010	1PE	C24-C14-OH5-C25
7	D	1012	2PE	C18-C17-O16-C15
6	J	1006	1PE	OH6-C15-C25-OH5
6	E	1009	1PE	OH5-C14-C24-OH4

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Mol	Chain	Res	Type	Atoms
6	D	1007	1PE	C12-C22-OH3-C23
6	D	1009	1PE	C23-C13-OH4-C24
6	A	1010	1PE	C13-C23-OH3-C22
6	I	1007	1PE	C12-C22-OH3-C23
7	B	1007	2PE	O13-C14-C15-O16
6	J	1006	1PE	C14-C24-OH4-C13
6	K	1010	1PE	C14-C24-OH4-C13
6	L	1008	1PE	C24-C14-OH5-C25
7	H	1006	2PE	C14-C15-O16-C17
7	H	1007	2PE	C14-C15-O16-C17
6	E	1009	1PE	C23-C13-OH4-C24
7	H	1006	2PE	C17-C18-O19-C20
6	J	1006	1PE	C16-C26-OH6-C15
6	C	1008	1PE	C15-C25-OH5-C14
7	H	1006	2PE	C24-C23-O22-C21
6	L	1007	1PE	C24-C14-OH5-C25
7	B	1006	2PE	C2-C3-O4-C5
6	J	1006	1PE	C15-C25-OH5-C14
6	L	1009	1PE	C15-C25-OH5-C14
6	L	1008	1PE	C23-C13-OH4-C24
6	L	1009	1PE	C25-C15-OH6-C26
6	D	1008	1PE	C14-C24-OH4-C13
6	F	1007	1PE	C15-C25-OH5-C14
7	D	1010	2PE	C11-C12-O13-C14
6	C	1007	1PE	C12-C22-OH3-C23
7	B	1006	2PE	C17-C18-O19-C20
6	F	1008	1PE	C16-C26-OH6-C15
7	B	1006	2PE	C20-C21-O22-C23
6	E	1007	1PE	OH6-C15-C25-OH5
7	B	1007	2PE	O16-C17-C18-O19
7	D	1012	2PE	C14-C15-O16-C17
6	D	1009	1PE	C16-C26-OH6-C15
7	H	1006	2PE	C21-C20-O19-C18
6	D	1008	1PE	OH5-C14-C24-OH4
6	K	1007	1PE	C14-C24-OH4-C13
6	L	1009	1PE	C14-C24-OH4-C13
7	B	1006	2PE	C23-C24-O25-C26
6	G	1008	1PE	C24-C14-OH5-C25
6	I	1008	1PE	C13-C23-OH3-C22
7	D	1010	2PE	O13-C14-C15-O16
7	B	1006	2PE	O1-C2-C3-O4
6	G	1011	1PE	C14-C24-OH4-C13

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Mol	Chain	Res	Type	Atoms
6	I	1007	1PE	C25-C15-OH6-C26
6	F	1009	1PE	C25-C15-OH6-C26
6	J	1008	1PE	C25-C15-OH6-C26
7	H	1007	2PE	C18-C17-O16-C15
6	I	1008	1PE	C14-C24-OH4-C13
6	J	1006	1PE	OH7-C16-C26-OH6
6	A	1010	1PE	OH4-C13-C23-OH3
7	D	1010	2PE	O16-C17-C18-O19
6	L	1007	1PE	C15-C25-OH5-C14
6	G	1011	1PE	C16-C26-OH6-C15
6	K	1009	1PE	C25-C15-OH6-C26
7	D	1010	2PE	C21-C20-O19-C18
7	B	1006	2PE	O22-C23-C24-O25
6	G	1008	1PE	C13-C23-OH3-C22
6	E	1007	1PE	C12-C22-OH3-C23
7	I	1010	2PE	C14-C15-O16-C17
6	J	1009	1PE	C23-C13-OH4-C24
6	G	1011	1PE	OH7-C16-C26-OH6
6	E	1008	1PE	OH5-C14-C24-OH4
6	G	1008	1PE	OH5-C14-C24-OH4
7	H	1006	2PE	C23-C24-O25-C26
6	K	1008	1PE	OH6-C15-C25-OH5
6	G	1011	1PE	C24-C14-OH5-C25
6	E	1008	1PE	C12-C22-OH3-C23
7	E	1010	2PE	C21-C20-O19-C18
7	B	1007	2PE	C14-C15-O16-C17
6	D	1009	1PE	C24-C14-OH5-C25
6	L	1006	1PE	C16-C26-OH6-C15
7	F	1010	2PE	C8-C9-O10-C11
7	B	1006	2PE	C15-C14-O13-C12
7	J	1010	2PE	C18-C17-O16-C15
6	A	1010	1PE	C25-C15-OH6-C26
6	D	1008	1PE	C13-C23-OH3-C22
6	C	1007	1PE	OH5-C14-C24-OH4
7	D	1010	2PE	C18-C17-O16-C15
6	D	1009	1PE	C15-C25-OH5-C14
6	J	1009	1PE	C14-C24-OH4-C13
6	I	1007	1PE	OH5-C14-C24-OH4
7	B	1006	2PE	C18-C17-O16-C15
7	C	1009	2PE	C18-C17-O16-C15
7	J	1010	2PE	C15-C14-O13-C12
6	K	1008	1PE	C15-C25-OH5-C14

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Mol	Chain	Res	Type	Atoms
6	G	1008	1PE	C23-C13-OH4-C24
6	D	1009	1PE	OH5-C14-C24-OH4
6	F	1006	1PE	C24-C14-OH5-C25
7	H	1007	2PE	C12-C11-O10-C9
7	F	1010	2PE	C15-C14-O13-C12
6	A	1010	1PE	OH6-C15-C25-OH5
6	D	1007	1PE	C23-C13-OH4-C24
7	C	1009	2PE	C11-C12-O13-C14
6	E	1007	1PE	OH5-C14-C24-OH4
6	J	1009	1PE	C15-C25-OH5-C14
6	A	1010	1PE	C15-C25-OH5-C14
6	E	1008	1PE	C23-C13-OH4-C24
6	K	1007	1PE	C12-C22-OH3-C23
6	L	1007	1PE	C14-C24-OH4-C13
7	D	1012	2PE	C15-C14-O13-C12
6	D	1009	1PE	C14-C24-OH4-C13
7	D	1010	2PE	C15-C14-O13-C12
6	L	1007	1PE	C12-C22-OH3-C23
7	J	1010	2PE	C17-C18-O19-C20
7	J	1010	2PE	C21-C20-O19-C18
6	F	1007	1PE	C25-C15-OH6-C26
6	G	1010	1PE	C15-C25-OH5-C14
6	L	1008	1PE	C14-C24-OH4-C13
7	B	1007	2PE	C11-C12-O13-C14
6	E	1007	1PE	OH4-C13-C23-OH3
6	G	1009	1PE	C15-C25-OH5-C14
7	B	1006	2PE	O16-C17-C18-O19
6	C	1006	1PE	C12-C22-OH3-C23
4	A	1004	1OT	N1-C1-C2-C3
4	B	1004	1OT	N1-C1-C2-C3
4	C	1004	1OT	N1-C1-C2-C3
4	H	1004	1OT	N1-C1-C2-C3
4	I	1004	1OT	N1-C1-C2-C3
4	K	1004	1OT	N1-C1-C2-C3
7	H	1006	2PE	O22-C23-C24-O25
6	C	1007	1PE	C23-C13-OH4-C24
7	E	1010	2PE	C18-C17-O16-C15

There are no ring outliers.

44 monomers are involved in 100 short contacts:

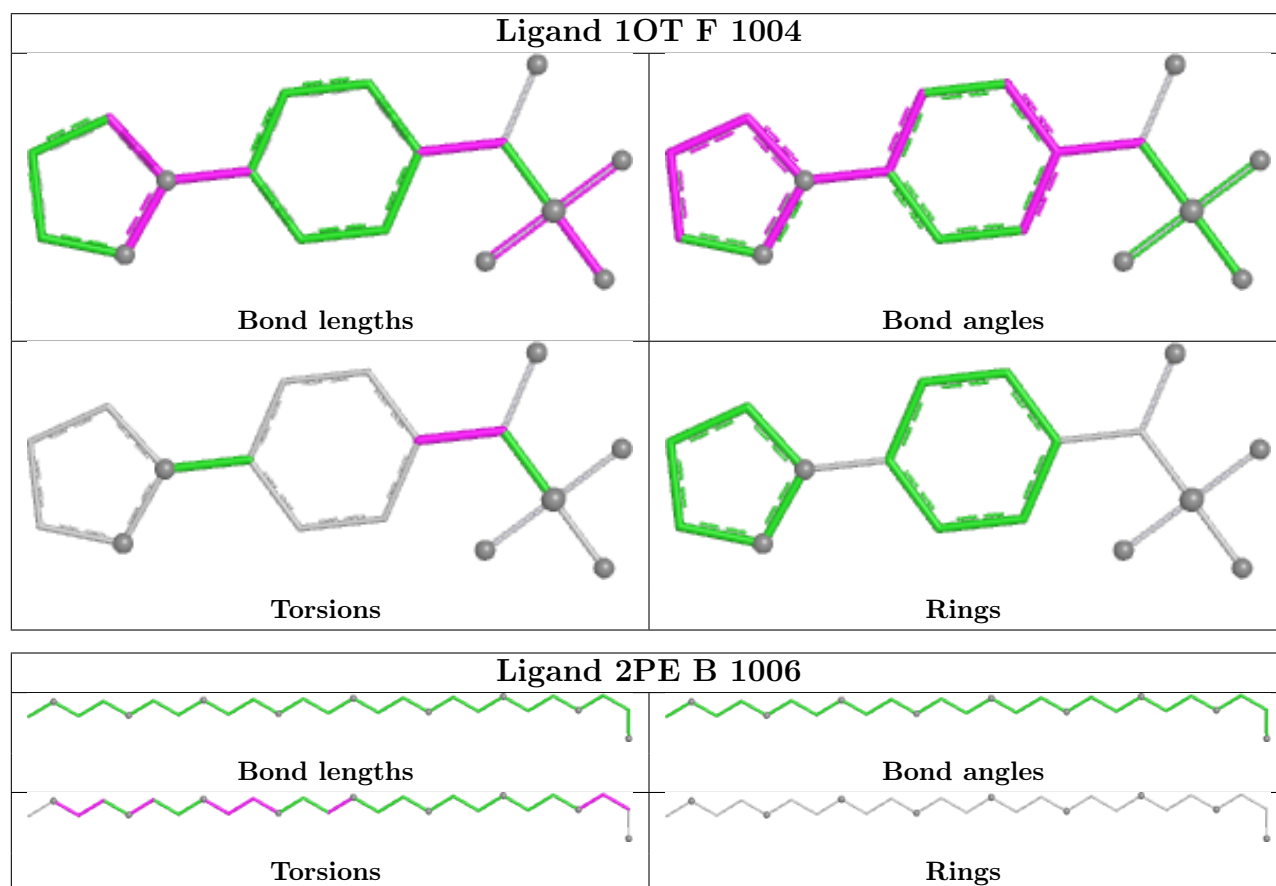
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	F	1008	1PE	1	0
4	F	1004	1OT	4	0
6	G	1008	1PE	2	0
3	D	1002	CO3	1	0
4	A	1004	1OT	4	0
6	K	1007	1PE	2	0
6	J	1008	1PE	2	0
6	K	1010	1PE	1	0
6	G	1011	1PE	6	0
6	L	1007	1PE	4	0
6	J	1006	1PE	3	0
6	I	1009	1PE	1	0
6	C	1006	1PE	1	0
6	D	1009	1PE	5	0
6	E	1009	1PE	4	0
6	A	1010	1PE	1	0
4	H	1004	1OT	2	0
6	J	1007	1PE	1	0
4	K	1004	1OT	3	0
4	D	1004	1OT	2	0
7	D	1012	2PE	1	0
7	H	1006	2PE	1	0
6	I	1007	1PE	1	0
5	J	1005	SO4	1	0
4	E	1004	1OT	4	0
6	J	1009	1PE	2	0
7	H	1007	2PE	1	0
7	F	1010	2PE	1	0
6	D	1007	1PE	2	0
4	C	1004	1OT	3	0
6	F	1009	1PE	1	0
7	E	1010	2PE	1	0
6	L	1008	1PE	2	0
6	L	1009	1PE	4	0
4	L	1004	1OT	4	0
3	L	1002	CO3	1	0
6	C	1007	1PE	1	0
4	G	1004	1OT	4	0
4	J	1004	1OT	3	0
7	C	1009	2PE	2	0
7	D	1010	2PE	4	0
4	B	1004	1OT	4	0
6	I	1008	1PE	1	0

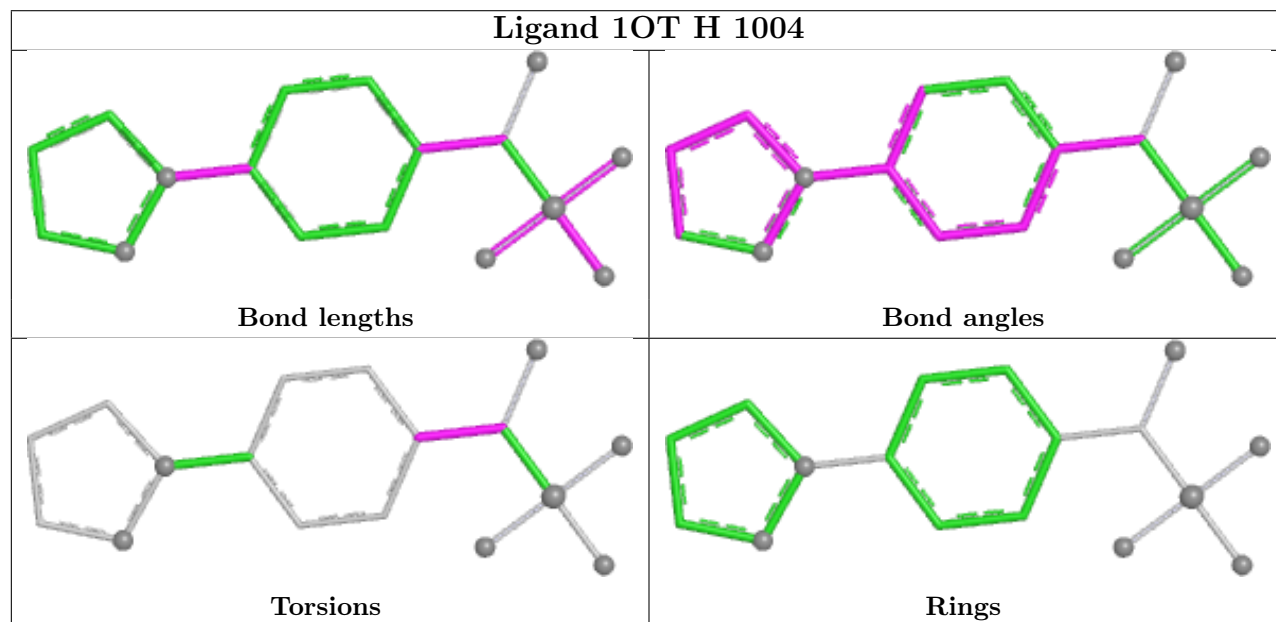
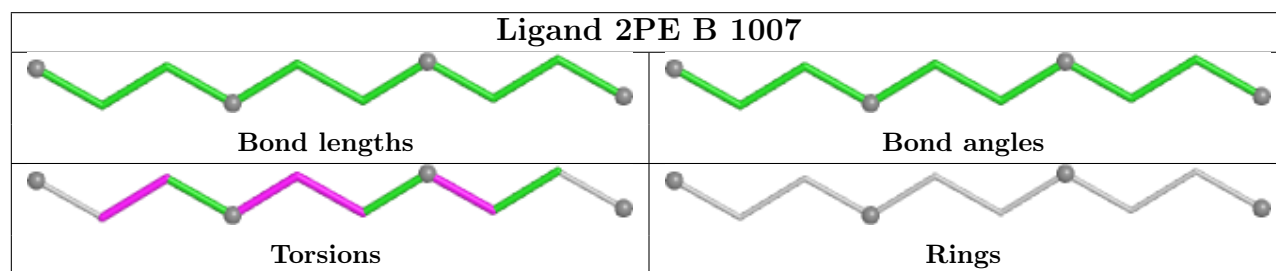
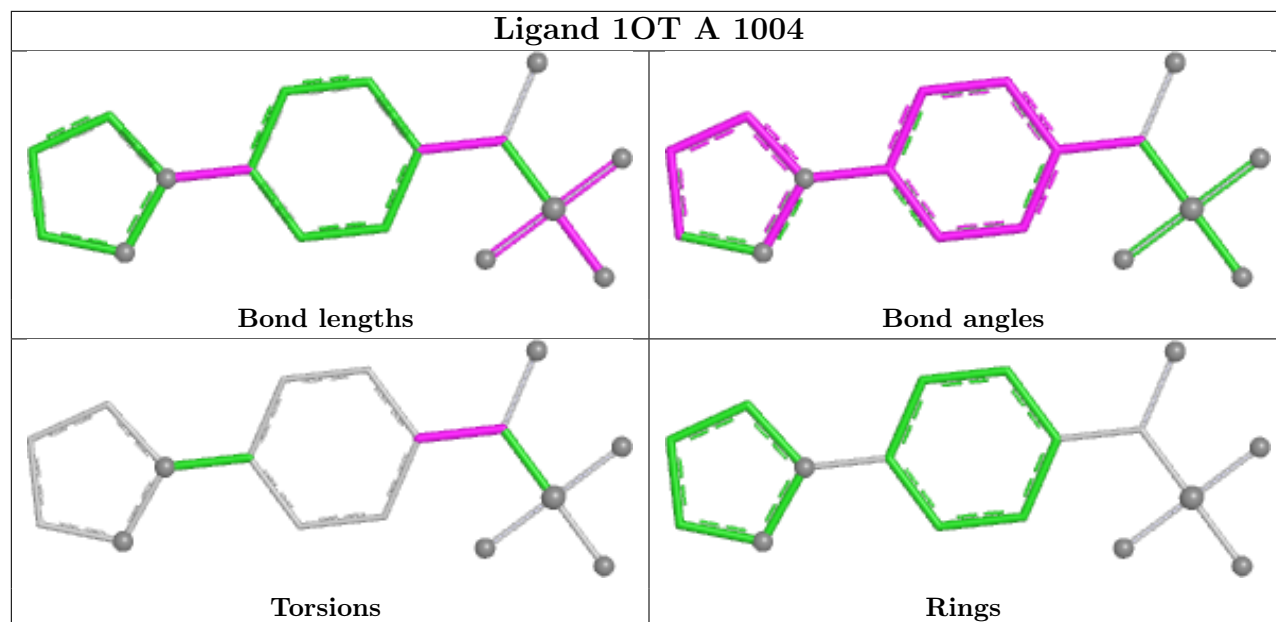
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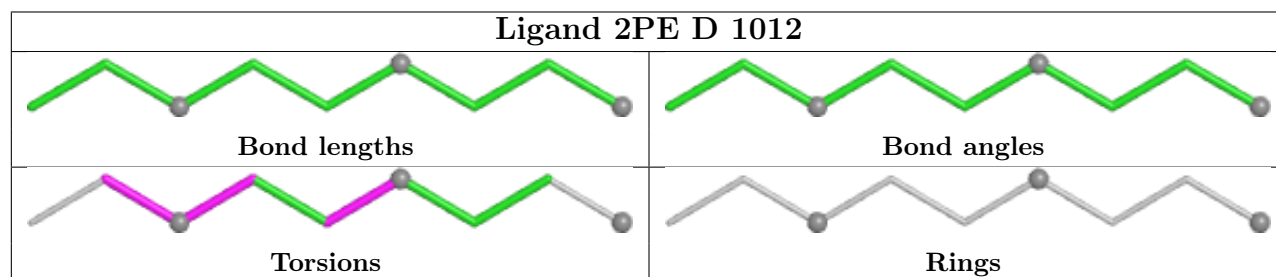
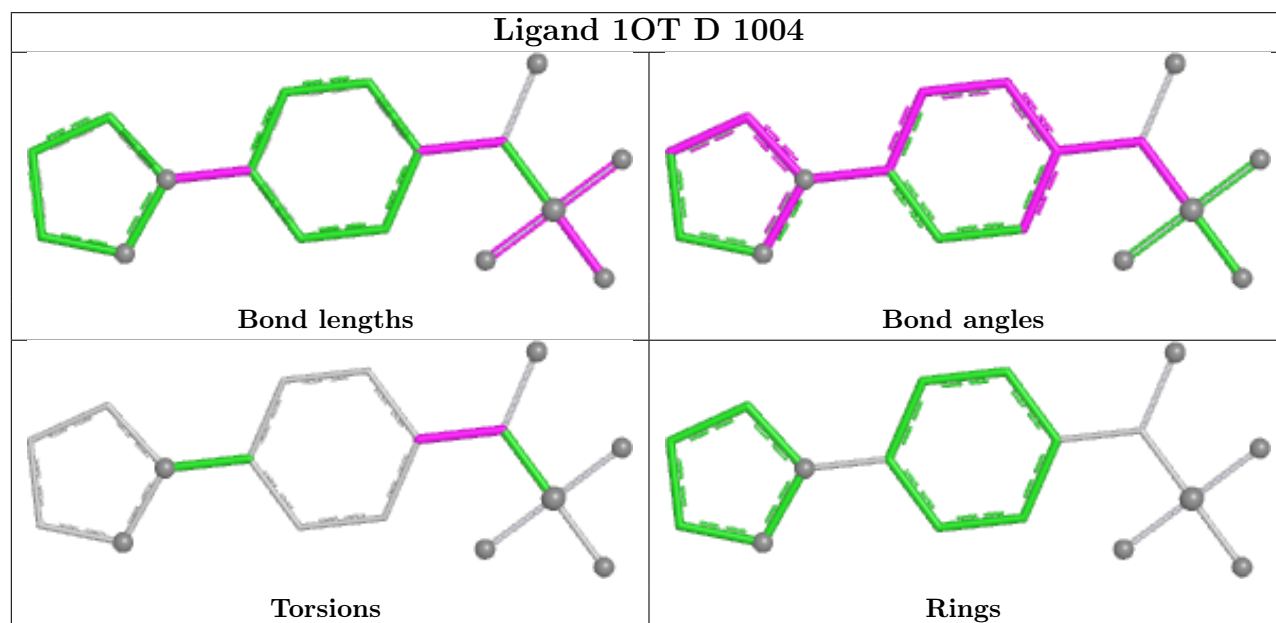
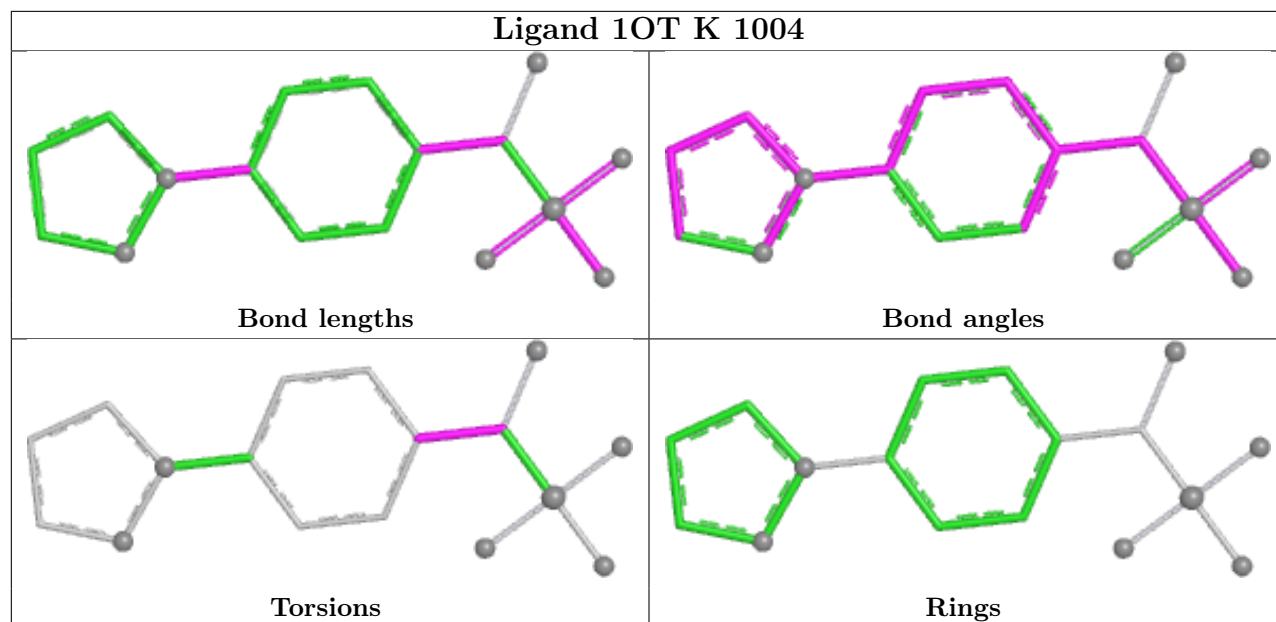
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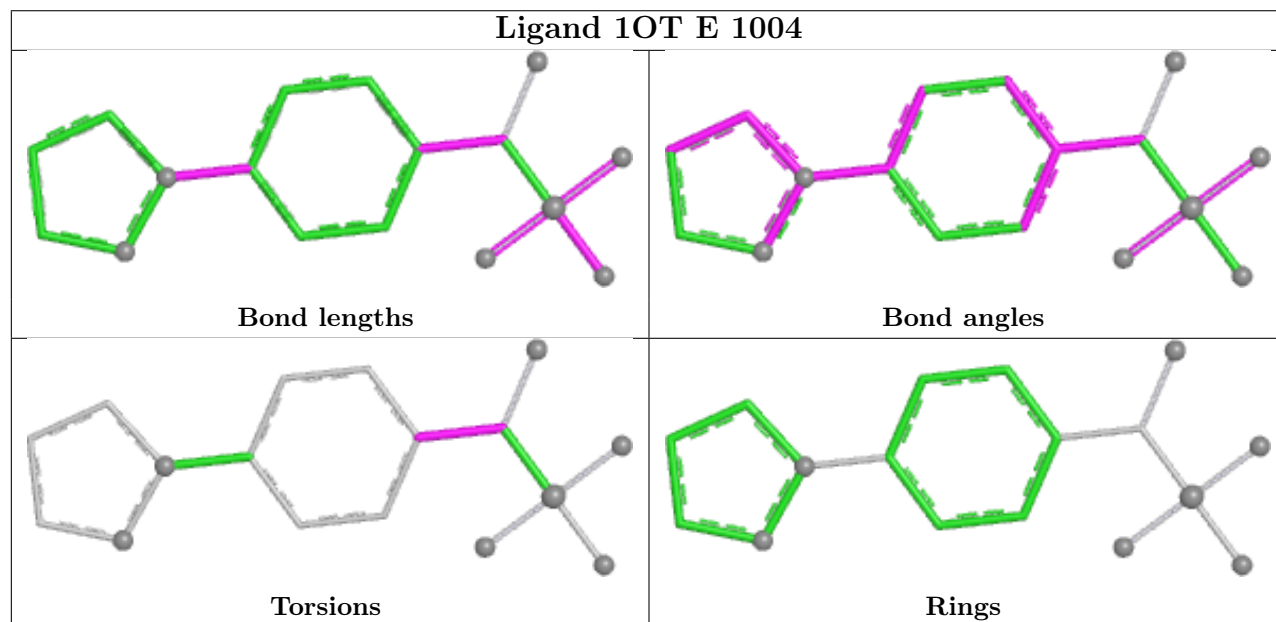
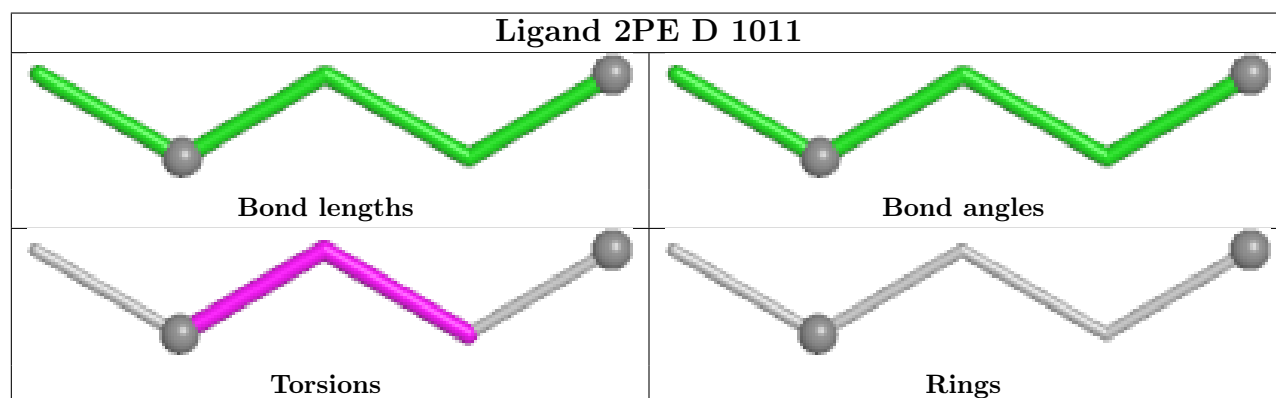
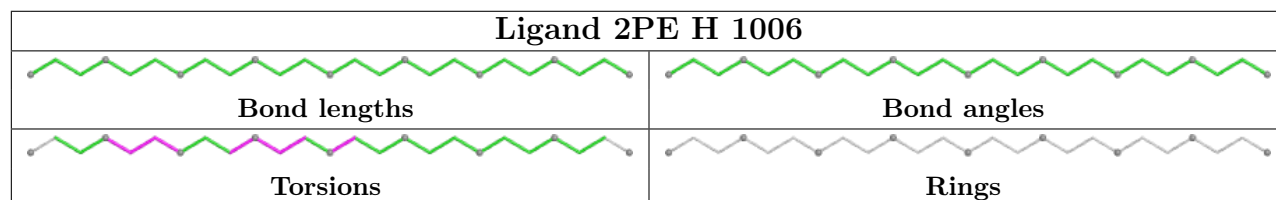
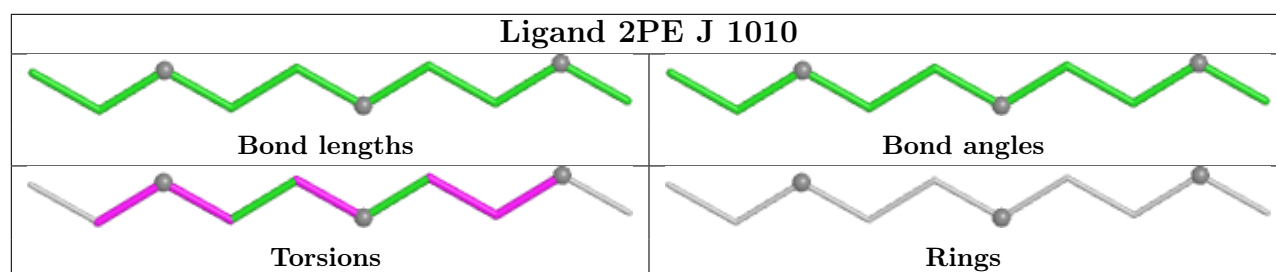
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	I	1004	1OT	3	0

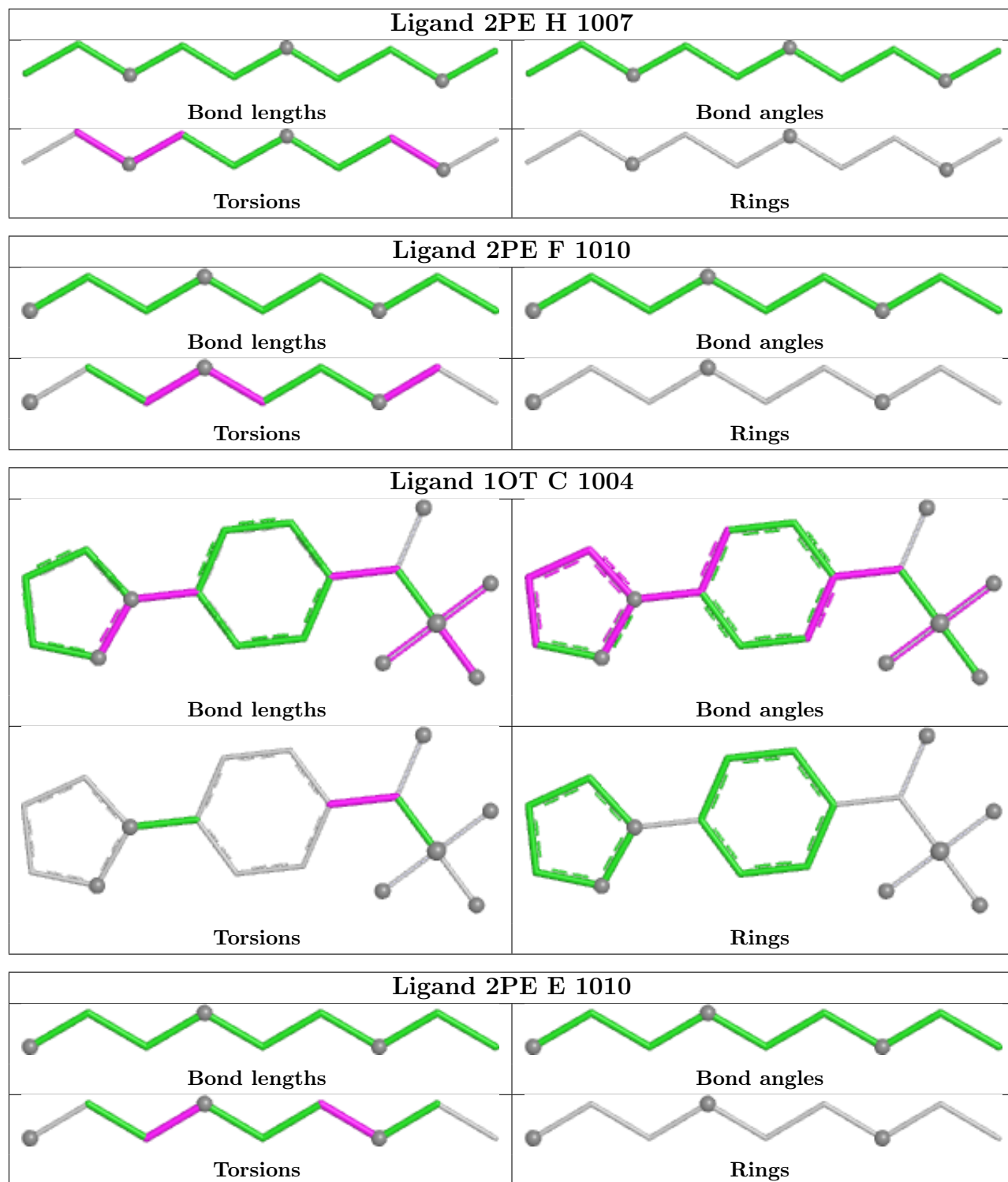
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

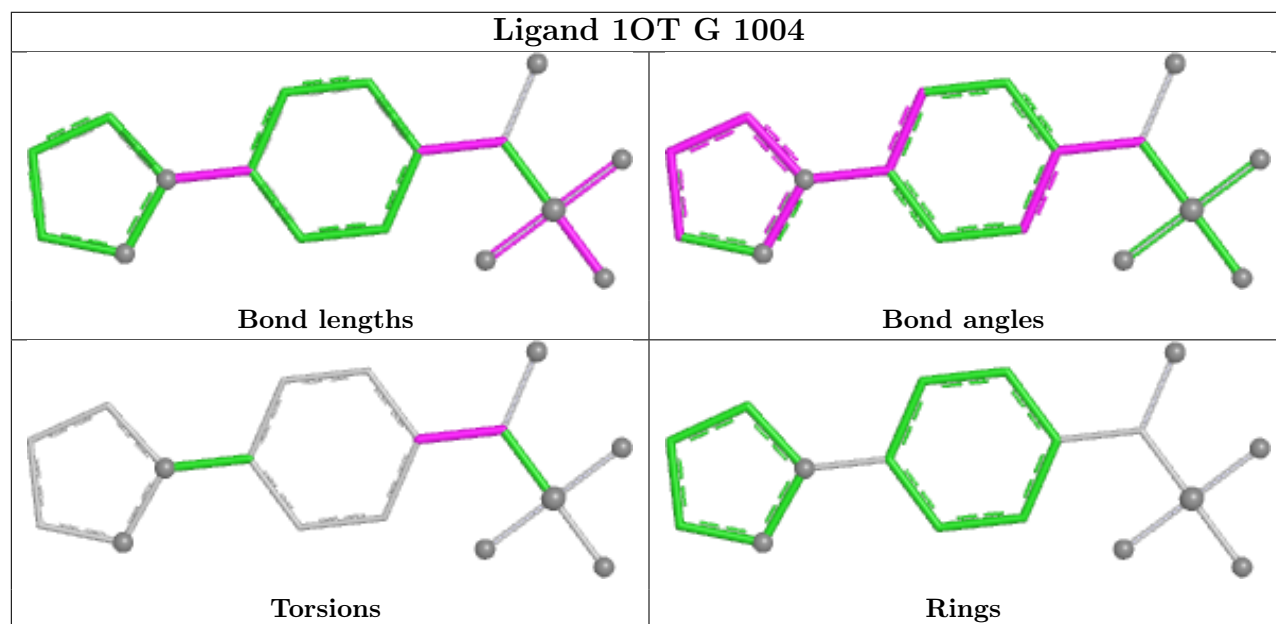
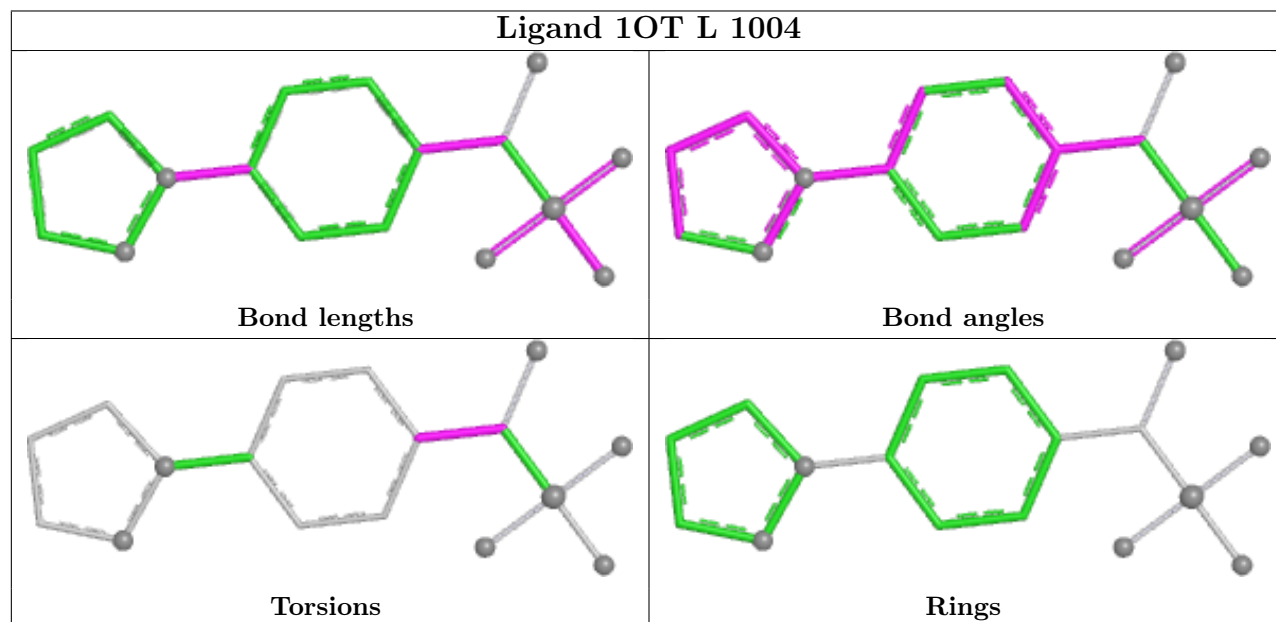
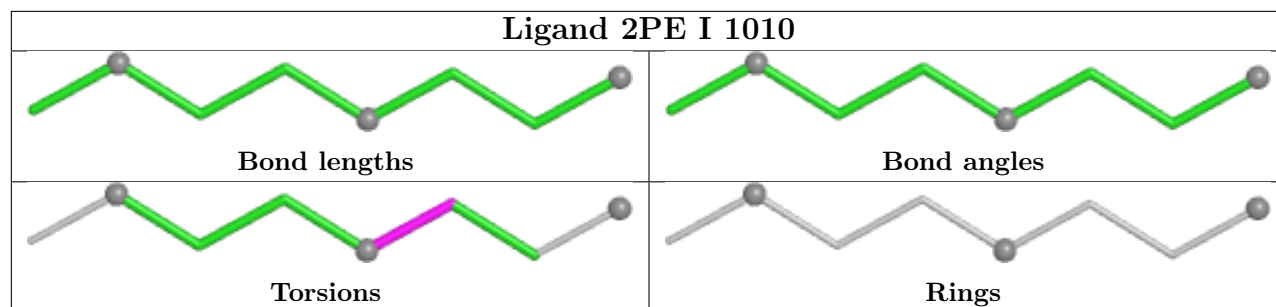


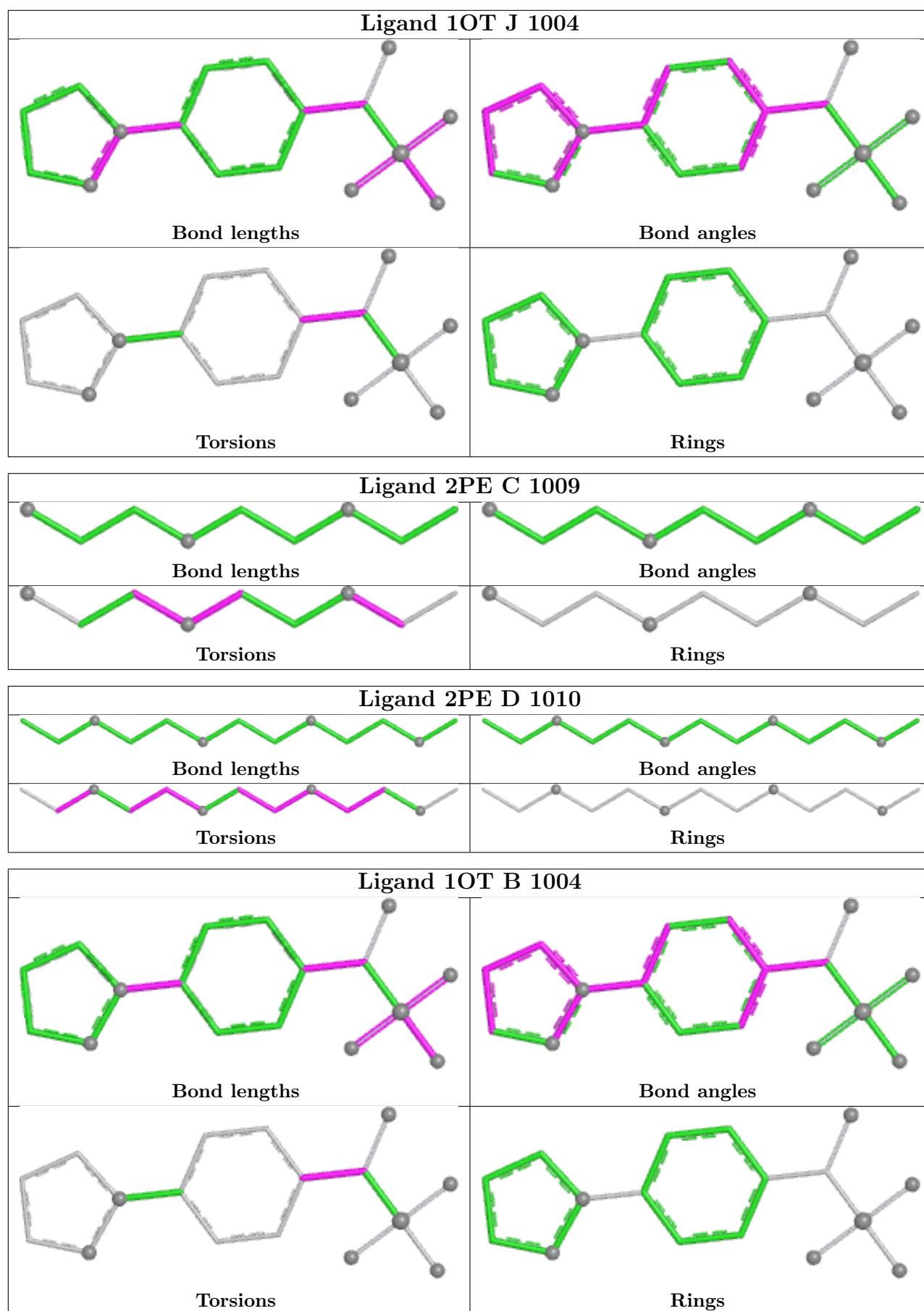


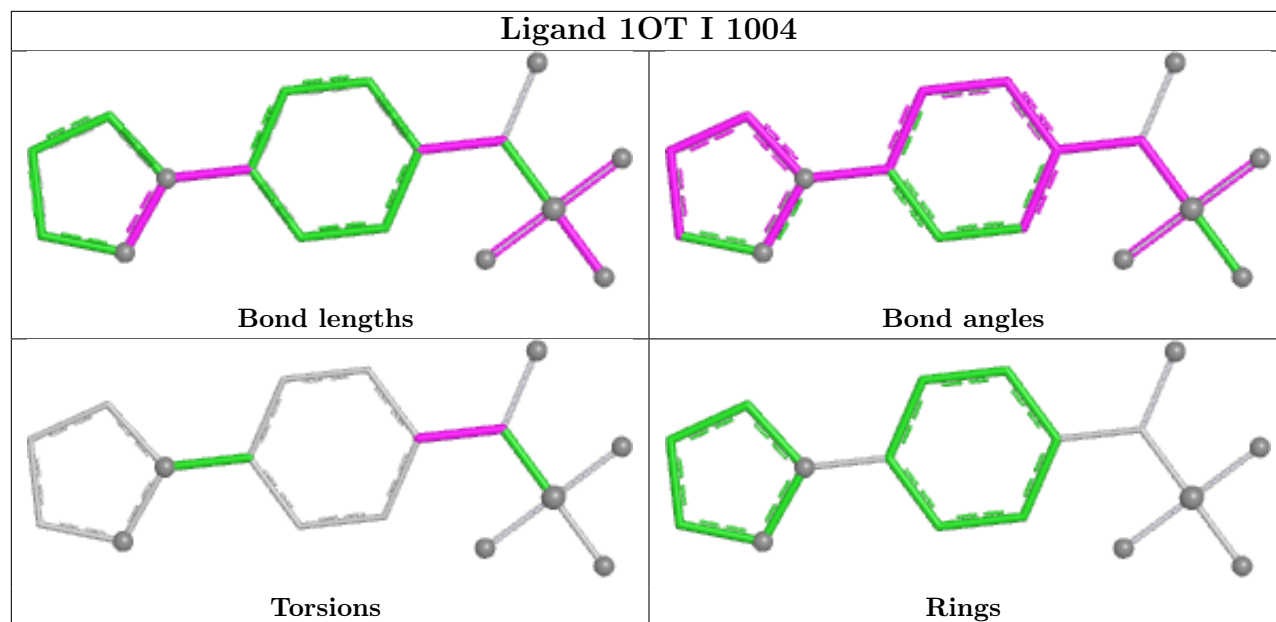












5.7 Other polymers [i](#)

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	519/528 (98%)	-0.49	6 (1%) 76 76	6, 13, 30, 50	1 (0%)
1	B	518/528 (98%)	-0.27	11 (2%) 63 63	5, 14, 43, 65	1 (0%)
1	C	518/528 (98%)	-0.48	5 (0%) 79 79	4, 12, 33, 48	1 (0%)
1	D	514/528 (97%)	-0.48	10 (1%) 66 66	5, 12, 31, 54	0
1	E	510/528 (96%)	-0.61	4 (0%) 82 82	4, 11, 25, 50	0
1	F	511/528 (96%)	-0.27	9 (1%) 67 67	6, 16, 37, 69	0
1	G	519/528 (98%)	-0.50	5 (0%) 79 79	5, 12, 31, 52	0
1	H	516/528 (97%)	-0.30	11 (2%) 63 63	5, 14, 41, 63	1 (0%)
1	I	518/528 (98%)	-0.47	7 (1%) 73 73	4, 13, 33, 56	0
1	J	513/528 (97%)	-0.49	10 (1%) 66 66	4, 13, 32, 57	0
1	K	509/528 (96%)	-0.58	3 (0%) 85 85	5, 12, 26, 52	0
1	L	513/528 (97%)	-0.34	8 (1%) 70 70	5, 14, 35, 52	0
All	All	6178/6336 (97%)	-0.44	89 (1%) 73 73	4, 13, 34, 69	4 (0%)

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	259	VAL	6.1
1	D	603	ASP	5.9
1	F	137	LYS	5.0
1	K	550	SER	4.8
1	F	138	GLU	4.5
1	B	257	LYS	4.3
1	D	219	LEU	3.9
1	E	550	SER	3.6
1	F	360	LYS	3.6
1	J	219	LEU	3.4
1	B	256	ASP	3.3

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Mol	Chain	Res	Type	RSRZ
1	J	549	SER	3.3
1	D	123	VAL	3.3
1	J	603	ASP	3.3
1	F	361	SER	3.2
1	F	136	GLY	3.2
1	B	136	GLY	3.2
1	H	195	VAL	3.2
1	A	196	ALA	3.1
1	D	85	ALA	3.1
1	D	137	LYS	2.9
1	H	136	GLY	2.9
1	H	273	ASN	2.9
1	L	363	GLY	2.9
1	J	216	ASP	2.9
1	F	549	SER	2.9
1	D	136	GLY	2.8
1	H	196	ALA	2.8
1	L	136	GLY	2.8
1	H	261	MET	2.7
1	D	124	GLU	2.7
1	I	138	GLU	2.7
1	D	260	ASN	2.6
1	H	137	LYS	2.6
1	B	205	ARG	2.6
1	H	258	ASN	2.6
1	I	196	ALA	2.6
1	F	119	GLY	2.5
1	A	602	ASN	2.5
1	J	85	ALA	2.5
1	G	195	VAL	2.5
1	J	136	GLY	2.5
1	G	549	SER	2.5
1	B	258	ASN	2.5
1	A	197	ASP	2.4
1	A	362	LYS	2.4
1	G	196	ALA	2.4
1	L	605	LEU	2.4
1	H	138	GLU	2.4
1	G	137	LYS	2.4
1	B	135	PRO	2.4
1	J	123	VAL	2.3
1	B	198	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	549	SER	2.3
1	K	549	SER	2.3
1	D	217	ASN	2.3
1	E	136	GLY	2.3
1	L	549	SER	2.2
1	I	195	VAL	2.2
1	J	137	LYS	2.2
1	C	196	ALA	2.2
1	C	124	GLU	2.2
1	H	257	LYS	2.2
1	E	86	SER	2.2
1	E	549	SER	2.2
1	B	137	LYS	2.2
1	H	135	PRO	2.2
1	C	136	GLY	2.2
1	G	362	LYS	2.2
1	C	195	VAL	2.1
1	A	85	ALA	2.1
1	A	137	LYS	2.1
1	J	262	GLU	2.1
1	B	195	VAL	2.1
1	I	136	GLY	2.1
1	I	124	GLU	2.1
1	K	363	GLY	2.1
1	I	137	LYS	2.1
1	L	137	LYS	2.1
1	F	550	SER	2.1
1	L	550	SER	2.1
1	F	121	CYS	2.1
1	I	549	SER	2.0
1	B	260	ASN	2.0
1	J	217	ASN	2.0
1	C	603	ASP	2.0
1	L	216	ASP	2.0
1	H	118	LYS	2.0
1	L	362	LYS	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($\text{\AA}^2$)	Q<0.9
6	1PE	D	1008	9/16	0.64	0.25	46,49,60,60	0
6	1PE	K	1010	6/16	0.67	0.19	23,28,37,39	0
6	1PE	J	1008	11/16	0.69	0.20	26,32,42,42	0
6	1PE	A	1010	12/16	0.70	0.21	32,38,46,48	0
6	1PE	C	1008	7/16	0.71	0.24	42,43,44,45	0
6	1PE	L	1008	8/16	0.71	0.20	48,51,53,53	0
6	1PE	C	1006	13/16	0.74	0.15	12,34,44,45	0
6	1PE	J	1009	9/16	0.74	0.19	26,32,38,39	0
6	1PE	G	1011	15/16	0.75	0.19	26,38,47,51	0
6	1PE	L	1009	11/16	0.76	0.16	36,39,43,45	0
7	2PE	H	1007	10/28	0.76	0.17	36,45,53,53	0
6	1PE	F	1009	11/16	0.77	0.16	46,51,55,56	0
7	2PE	D	1012	9/28	0.77	0.16	36,42,45,47	0
5	SO4	B	1005	5/5	0.77	0.13	67,68,68,69	0
6	1PE	D	1009	11/16	0.78	0.17	23,35,40,43	0
6	1PE	L	1007	12/16	0.78	0.15	28,35,39,41	0
7	2PE	I	1010	8/28	0.78	0.16	32,39,48,49	0
7	2PE	J	1010	10/28	0.78	0.20	18,47,48,48	0
7	2PE	E	1010	9/28	0.79	0.18	36,39,42,46	0
6	1PE	K	1007	12/16	0.79	0.15	31,37,40,41	0
6	1PE	G	1008	9/16	0.80	0.14	17,25,29,30	0
6	1PE	F	1007	10/16	0.80	0.16	40,42,49,51	0
7	2PE	D	1010	13/28	0.81	0.17	28,38,49,51	0
6	1PE	J	1007	10/16	0.81	0.15	28,32,39,41	0
6	1PE	E	1007	12/16	0.82	0.13	29,32,34,34	0
7	2PE	F	1010	9/28	0.82	0.15	35,38,39,42	0
7	2PE	H	1006	25/28	0.82	0.13	21,31,39,44	0
7	2PE	B	1007	10/28	0.82	0.17	45,49,53,55	0
6	1PE	E	1009	8/16	0.82	0.16	29,42,43,43	0
6	1PE	K	1009	11/16	0.82	0.13	30,33,36,38	0
7	2PE	B	1006	26/28	0.84	0.15	17,43,51,51	0
6	1PE	I	1007	15/16	0.84	0.14	20,31,40,40	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	2PE	C	1009	9/28	0.85	0.13	29,32,38,39	0
6	1PE	I	1008	11/16	0.85	0.13	18,24,31,33	0
5	SO4	I	1005	5/5	0.85	0.12	50,51,54,55	0
6	1PE	C	1007	9/16	0.86	0.12	13,22,29,31	0
6	1PE	I	1009	5/16	0.87	0.12	14,20,20,21	0
5	SO4	C	1005	5/5	0.87	0.18	41,47,51,54	0
6	1PE	F	1006	10/16	0.87	0.12	29,32,36,37	0
6	1PE	G	1010	6/16	0.88	0.14	30,32,32,33	0
6	1PE	E	1008	12/16	0.88	0.12	19,23,39,39	0
6	1PE	F	1008	10/16	0.88	0.11	26,32,35,36	0
5	SO4	A	1008	5/5	0.88	0.14	26,35,41,48	0
5	SO4	E	1006	5/5	0.88	0.15	46,53,54,62	0
6	1PE	D	1007	10/16	0.89	0.10	17,24,28,32	0
5	SO4	G	1007	5/5	0.89	0.14	43,48,51,56	0
5	SO4	D	1006	5/5	0.89	0.19	44,51,54,55	0
5	SO4	A	1007	5/5	0.89	0.17	38,42,45,49	0
6	1PE	J	1006	11/16	0.89	0.11	20,25,35,37	0
5	SO4	I	1006	5/5	0.90	0.13	53,55,56,60	0
6	1PE	A	1009	9/16	0.90	0.10	18,21,25,34	0
7	2PE	D	1011	5/28	0.90	0.12	21,21,28,28	0
6	1PE	L	1006	10/16	0.90	0.11	19,25,29,35	0
5	SO4	E	1005	5/5	0.90	0.18	42,45,49,50	0
5	SO4	A	1006	5/5	0.91	0.15	45,46,50,50	0
5	SO4	F	1005	5/5	0.91	0.20	44,46,49,51	0
5	SO4	G	1005	5/5	0.91	0.17	32,32,33,40	0
5	SO4	L	1005	5/5	0.91	0.18	33,40,45,48	0
5	SO4	K	1006	5/5	0.92	0.18	42,45,47,53	0
5	SO4	G	1006	5/5	0.92	0.16	40,41,45,49	0
5	SO4	J	1005	5/5	0.92	0.10	51,51,55,56	0
6	1PE	G	1009	6/16	0.92	0.10	21,22,26,33	0
4	1OT	F	1004	17/17	0.93	0.11	10,24,32,32	0
4	1OT	J	1004	17/17	0.94	0.10	15,21,43,44	0
4	1OT	L	1004	17/17	0.94	0.10	7,25,34,34	0
6	1PE	K	1008	12/16	0.94	0.10	16,25,38,39	0
4	1OT	C	1004	17/17	0.95	0.09	9,18,36,37	0
4	1OT	K	1004	17/17	0.95	0.10	12,20,40,40	0
2	ZN	J	1003	1/1	0.95	0.21	53,53,53,53	0
4	1OT	G	1004	17/17	0.95	0.09	8,20,32,35	0
4	1OT	H	1004	17/17	0.95	0.10	10,24,37,37	0
4	1OT	I	1004	17/17	0.95	0.10	9,24,37,37	0
2	ZN	K	1003	1/1	0.96	0.20	46,46,46,46	0
4	1OT	B	1004	17/17	0.96	0.10	9,18,39,41	0

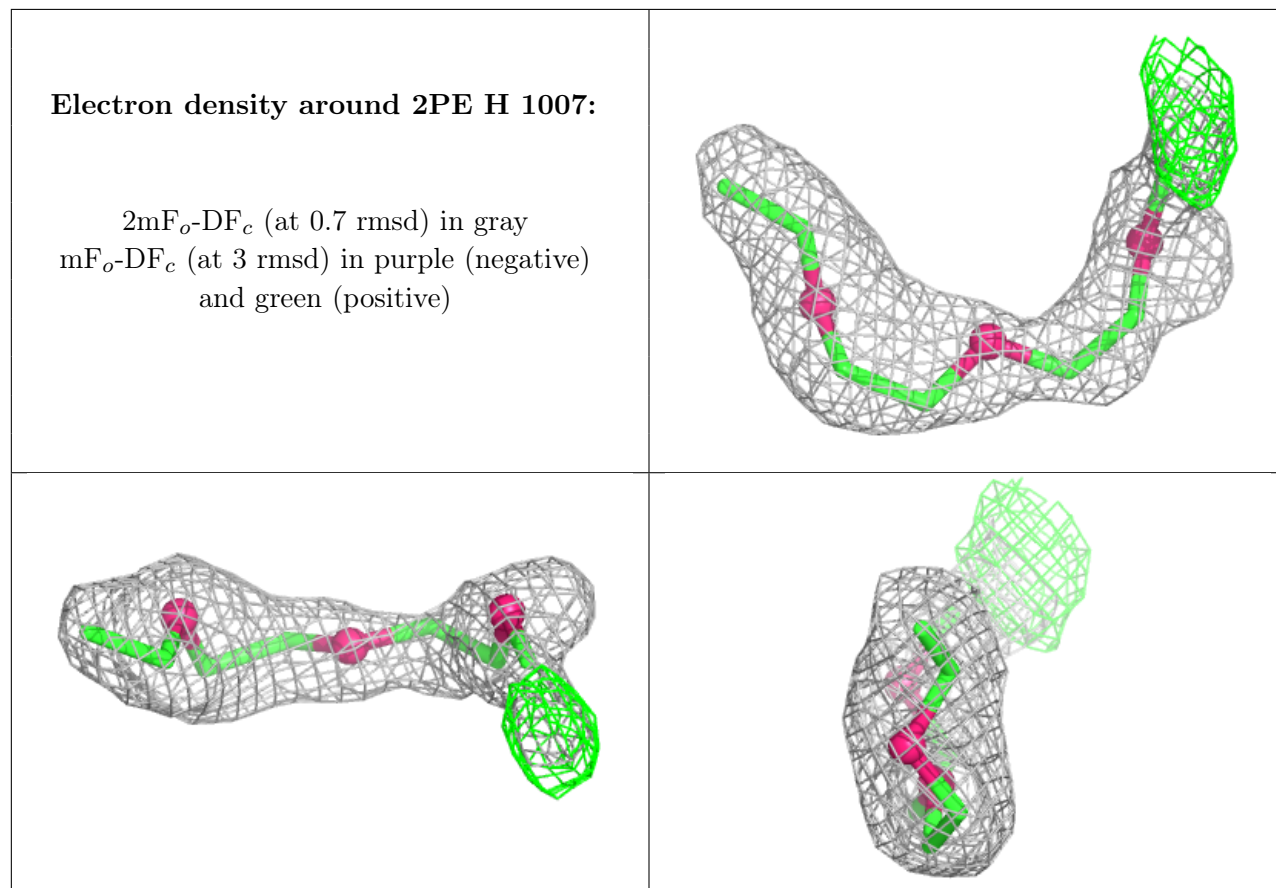
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	H	1003	1/1	0.96	0.20	46,46,46,46	0
4	1OT	D	1004	17/17	0.96	0.09	12,21,39,39	0
4	1OT	E	1004	17/17	0.96	0.09	14,20,41,42	0
3	CO3	F	1002	4/4	0.97	0.05	7,12,14,15	0
3	CO3	J	1002	4/4	0.97	0.04	5,6,11,15	0
4	1OT	A	1004	17/17	0.97	0.08	5,20,33,34	0
2	ZN	A	1003	1/1	0.97	0.16	46,46,46,46	0
2	ZN	D	1003	1/1	0.97	0.20	45,45,45,45	0
2	ZN	F	1001	1/1	0.97	0.13	39,39,39,39	0
3	CO3	D	1002	4/4	0.97	0.04	6,7,10,12	0
2	ZN	H	1001	1/1	0.98	0.17	36,36,36,36	0
3	CO3	K	1002	4/4	0.98	0.04	8,12,13,18	0
3	CO3	L	1002	4/4	0.98	0.04	6,8,10,12	0
2	ZN	B	1003	1/1	0.98	0.19	38,38,38,38	0
2	ZN	I	1001	1/1	0.98	0.17	40,40,40,40	0
2	ZN	E	1003	1/1	0.98	0.22	49,49,49,49	0
2	ZN	C	1003	1/1	0.98	0.21	41,41,41,41	0
2	ZN	L	1003	1/1	0.98	0.18	45,45,45,45	0
3	CO3	B	1002	4/4	0.98	0.04	5,6,7,11	0
2	ZN	F	1003	1/1	0.98	0.18	51,51,51,51	0
3	CO3	E	1002	4/4	0.98	0.04	8,10,11,17	0
2	ZN	G	1003	1/1	0.98	0.15	39,39,39,39	0
3	CO3	H	1002	4/4	0.98	0.04	4,10,10,12	0
2	ZN	J	1001	1/1	0.99	0.14	36,36,36,36	0
5	SO4	H	1005	5/5	0.99	0.03	7,8,10,11	0
3	CO3	I	1002	4/4	0.99	0.03	7,8,8,15	0
2	ZN	D	1001	1/1	0.99	0.16	36,36,36,36	0
5	SO4	A	1005	5/5	0.99	0.03	6,7,11,13	0
5	SO4	K	1005	5/5	0.99	0.04	9,11,11,12	0
2	ZN	K	1001	1/1	0.99	0.15	35,35,35,35	0
2	ZN	G	1001	1/1	0.99	0.11	39,39,39,39	0
2	ZN	L	1001	1/1	0.99	0.14	41,41,41,41	0
2	ZN	A	1001	1/1	0.99	0.13	37,37,37,37	0
3	CO3	A	1002	4/4	0.99	0.04	7,10,12,18	0
5	SO4	D	1005	5/5	0.99	0.03	9,9,11,12	0
2	ZN	E	1001	1/1	0.99	0.14	36,36,36,36	0
3	CO3	C	1002	4/4	0.99	0.04	8,10,12,17	0
2	ZN	C	1001	1/1	0.99	0.17	41,41,41,41	0
2	ZN	B	1001	1/1	0.99	0.14	36,36,36,36	0
2	ZN	I	1003	1/1	0.99	0.21	42,42,42,42	0
3	CO3	G	1002	4/4	0.99	0.03	7,8,8,17	0

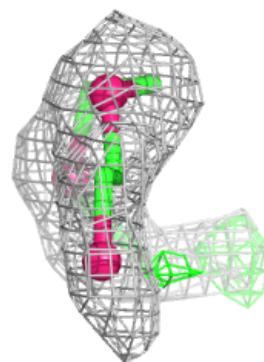
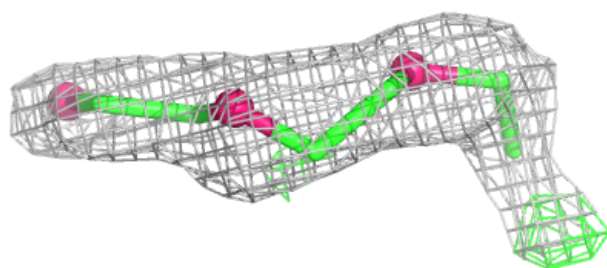
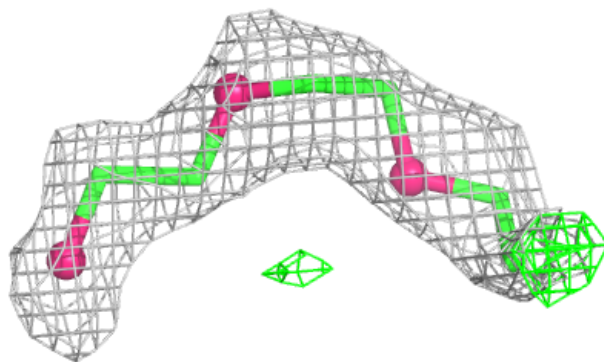
The following is a graphical depiction of the model fit to experimental electron density of all

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.

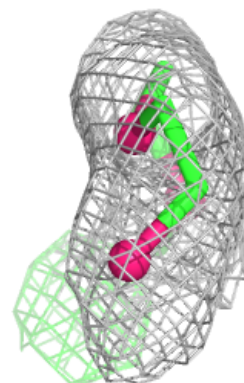
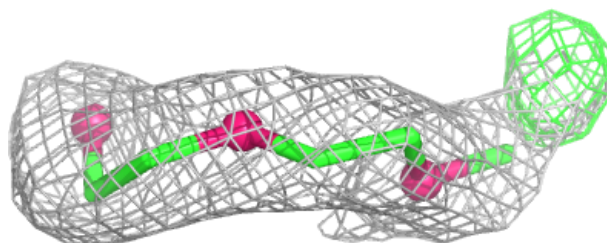
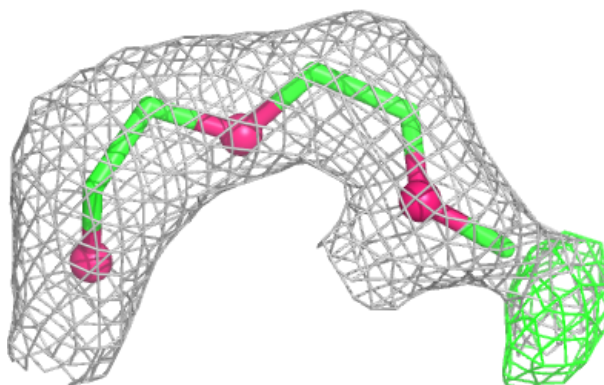


Electron density around 2PE D 1012:

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

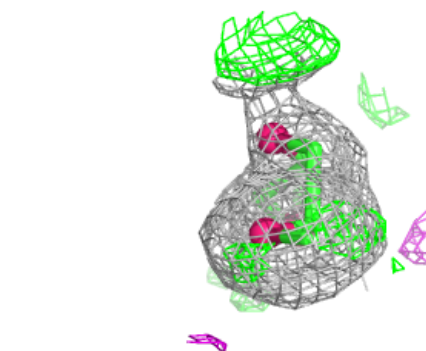
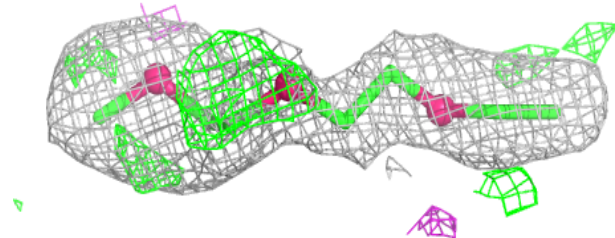
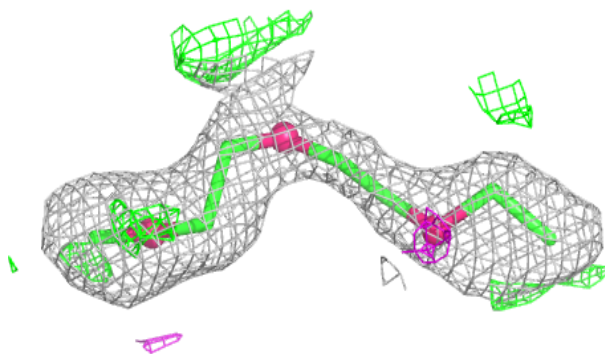
**Electron density around 2PE I 1010:**

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

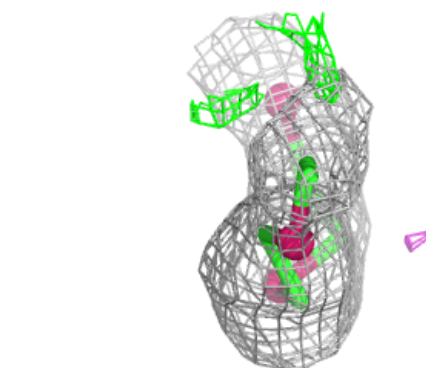
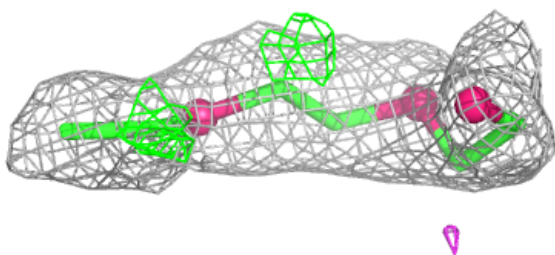
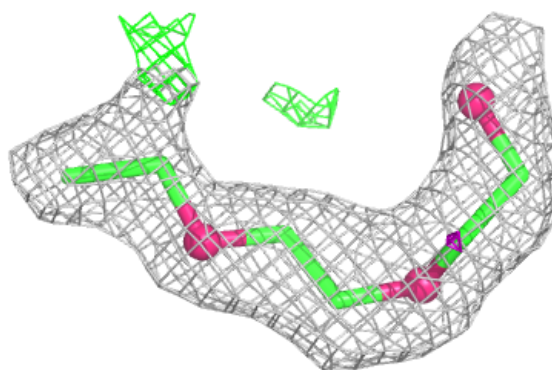


Electron density around 2PE J 1010:

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

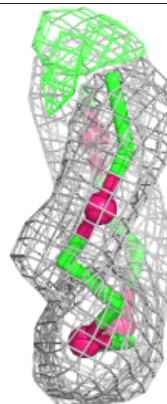
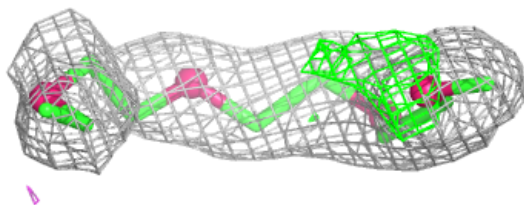
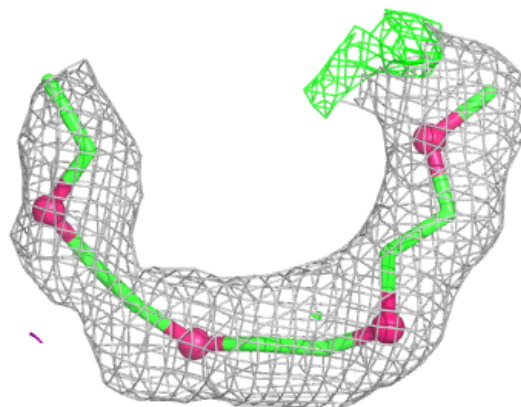
**Electron density around 2PE E 1010:**

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

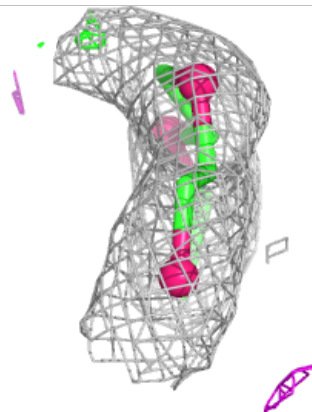
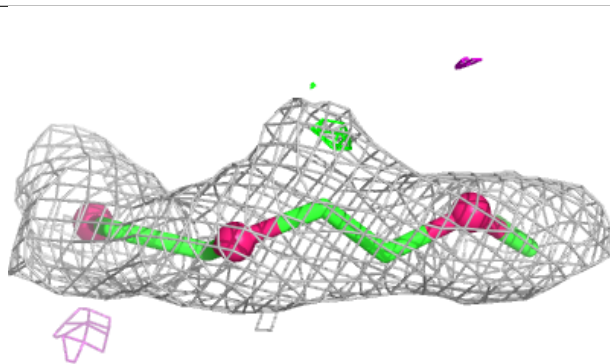
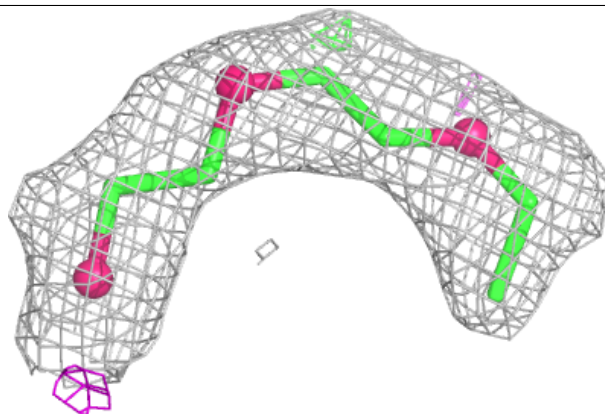


Electron density around 2PE D 1010:

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

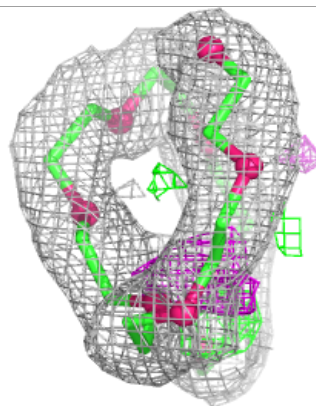
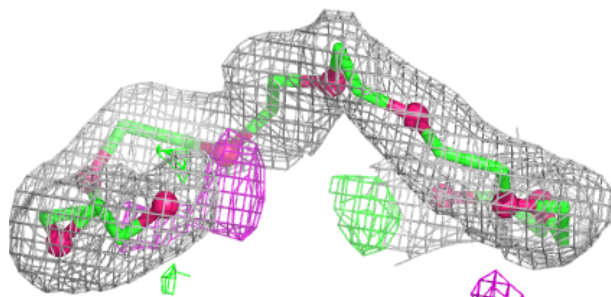
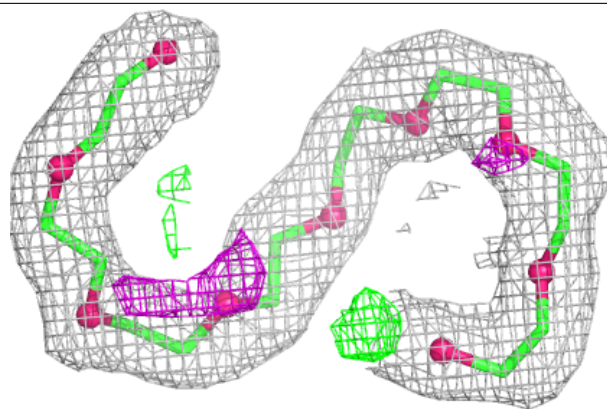
**Electron density around 2PE F 1010:**

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

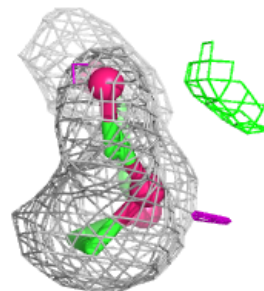
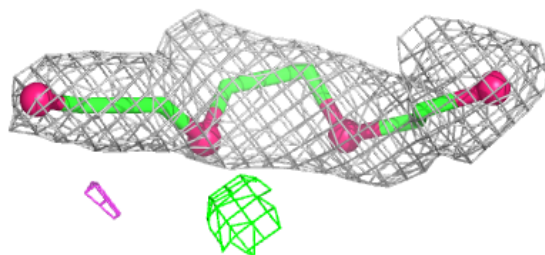
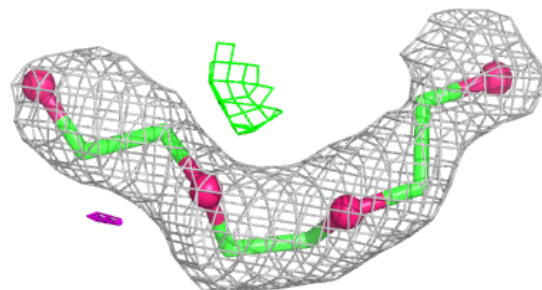


Electron density around 2PE H 1006:

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

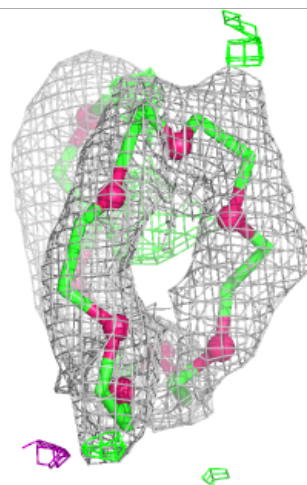
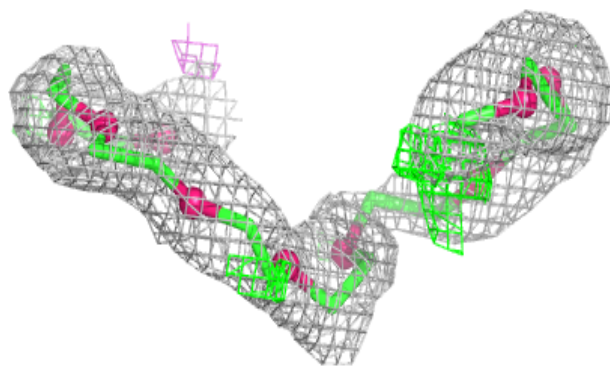
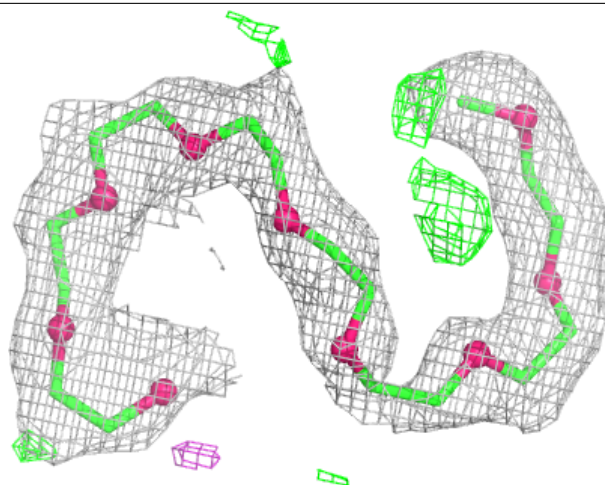
**Electron density around 2PE B 1007:**

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



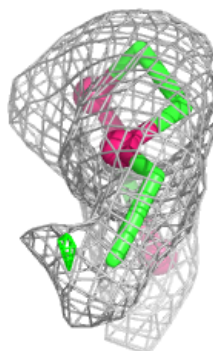
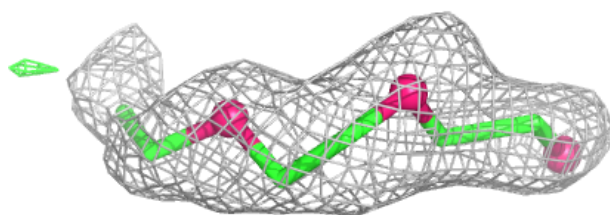
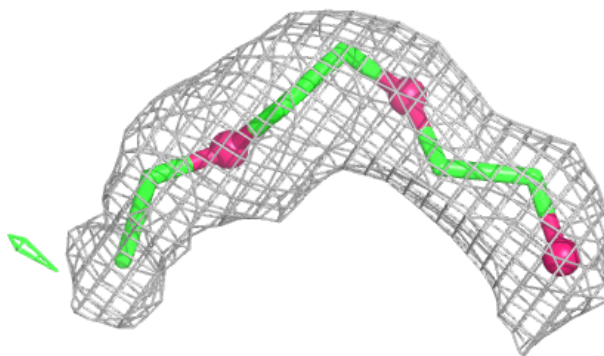
Electron density around 2PE B 1006:

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

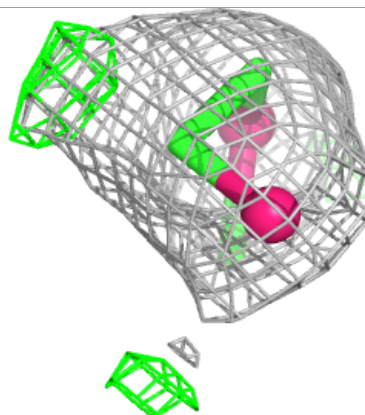
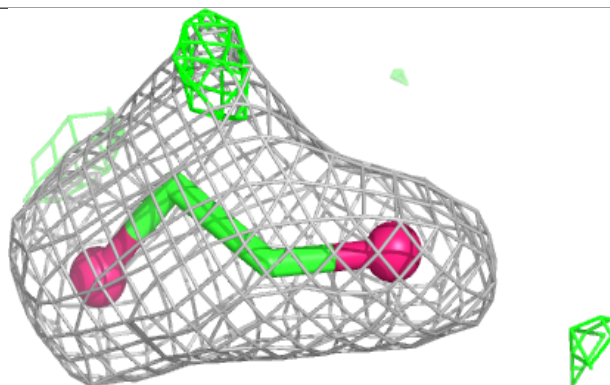
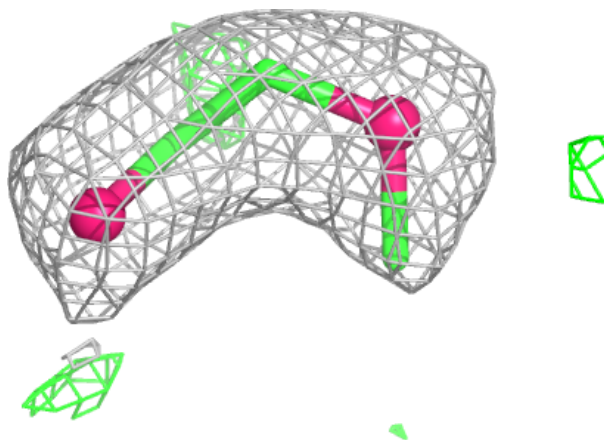


Electron density around 2PE C 1009:

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

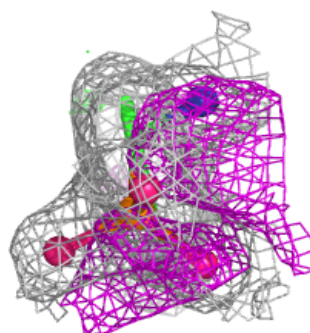
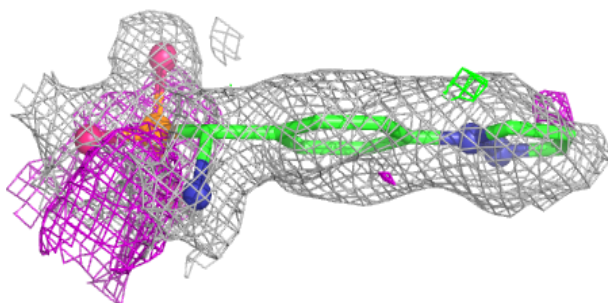
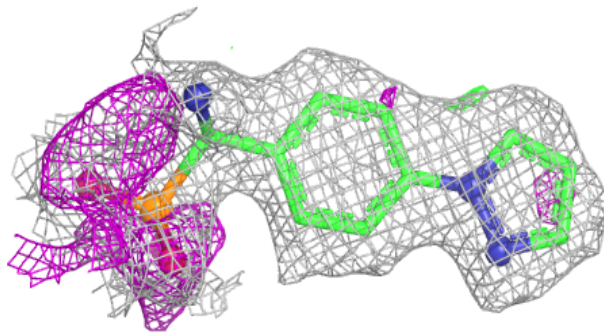
**Electron density around 2PE D 1011:**

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

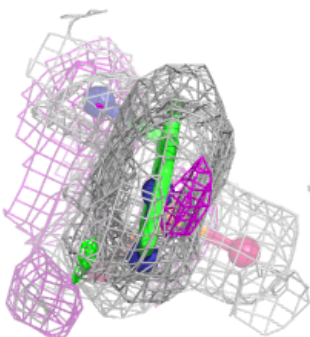
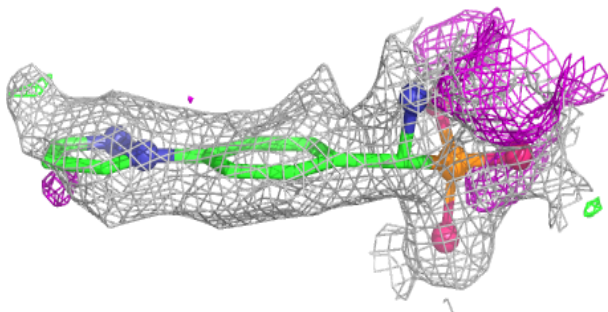
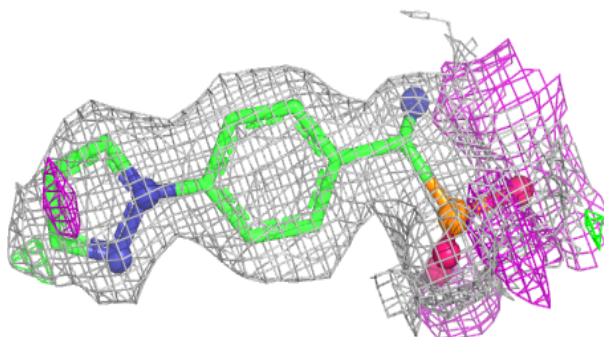


Electron density around 1OT F 1004:

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

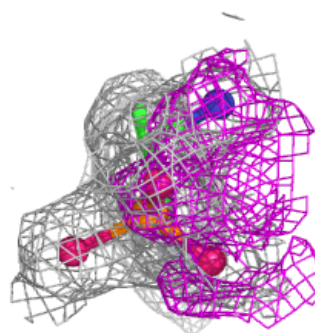
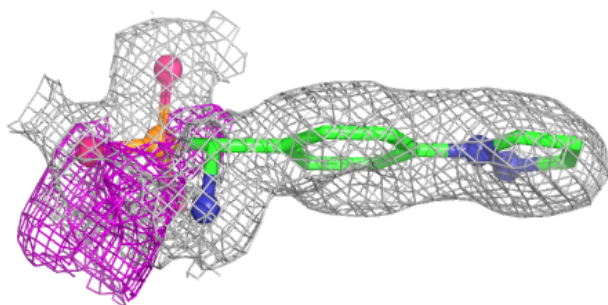
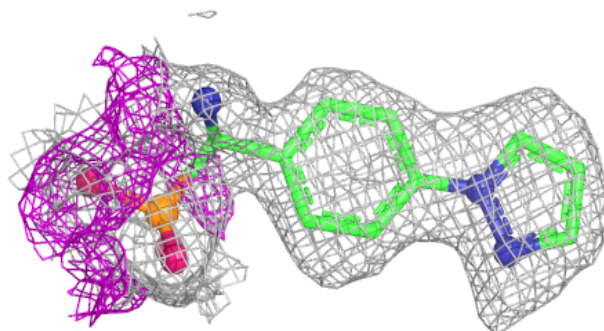
**Electron density around 1OT J 1004:**

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

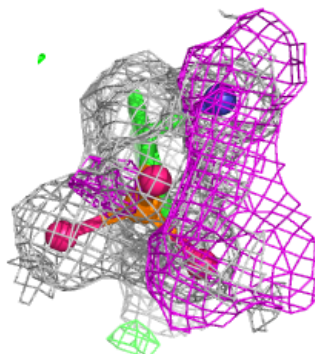
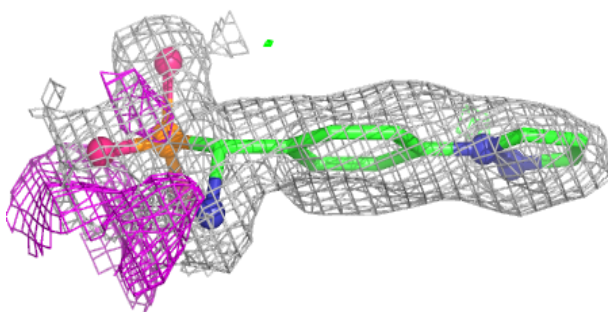
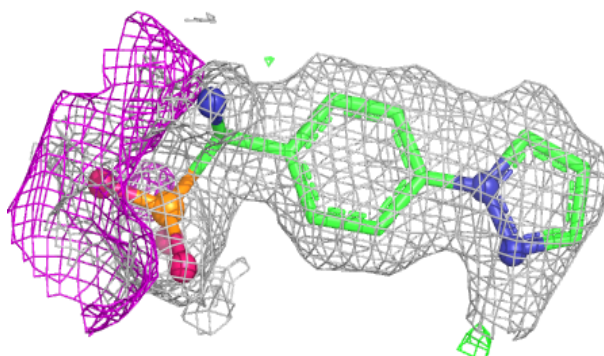


Electron density around 1OT L 1004:

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

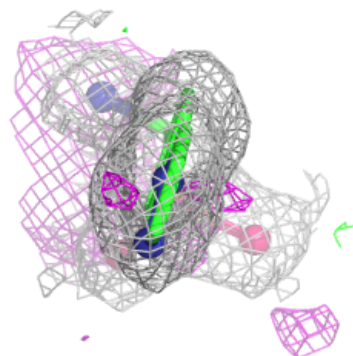
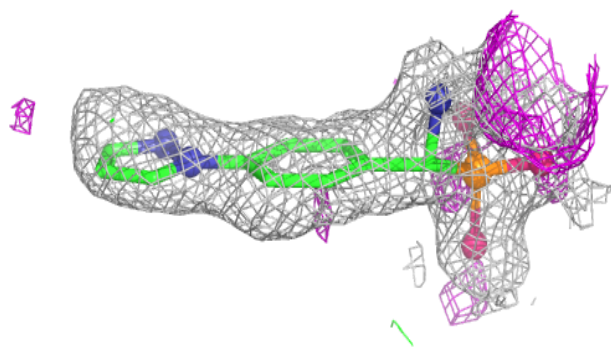
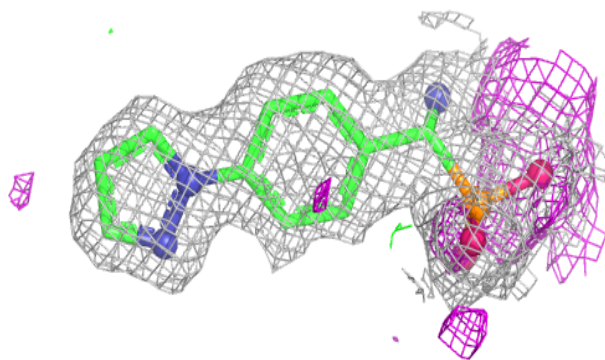
**Electron density around 1OT C 1004:**

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

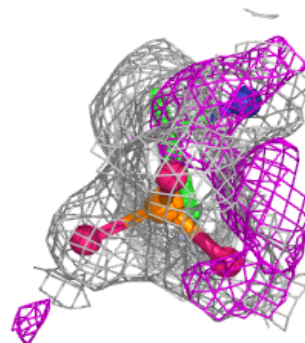
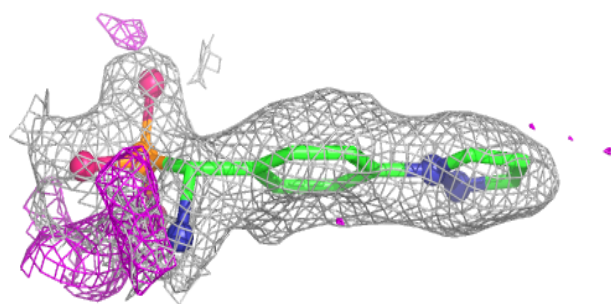
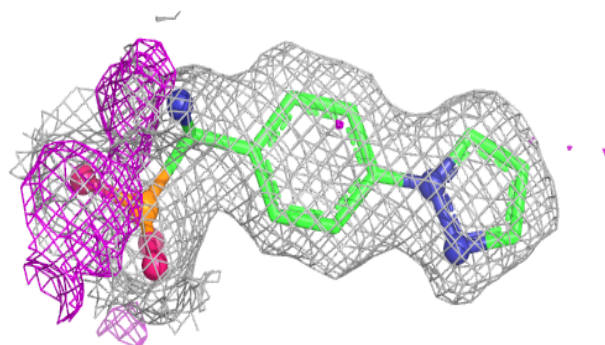


Electron density around 1OT K 1004:

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

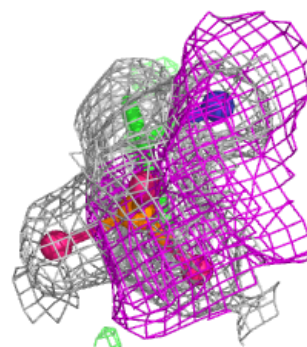
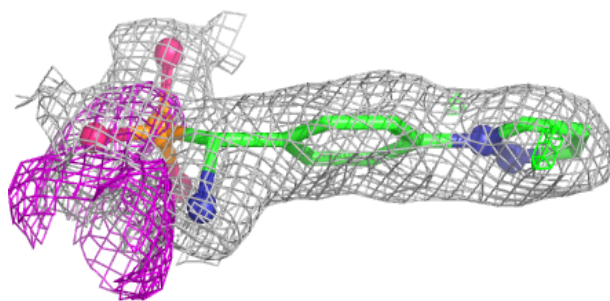
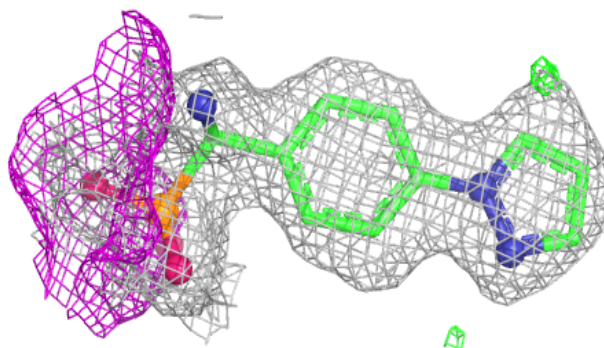
**Electron density around 1OT G 1004:**

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

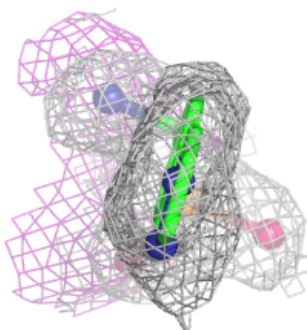
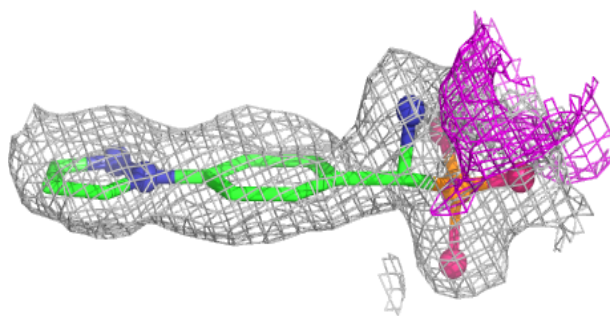
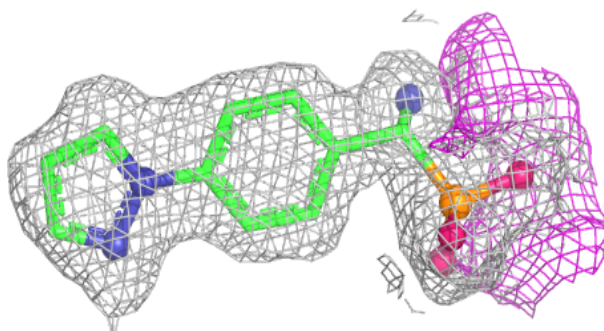


Electron density around 1OT H 1004:

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

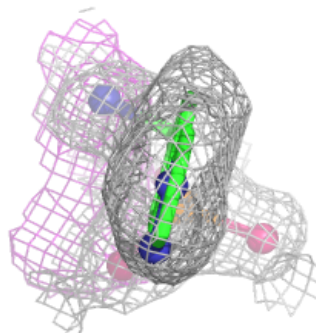
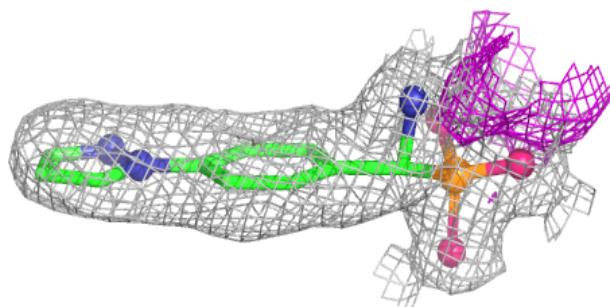
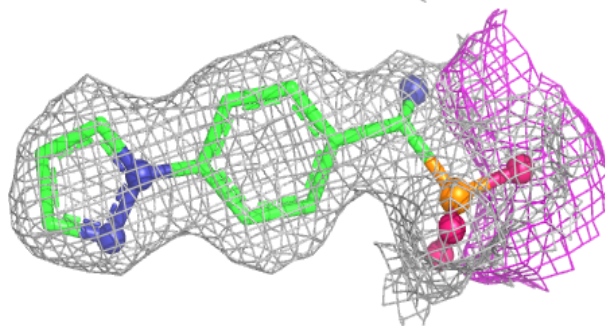
**Electron density around 1OT I 1004:**

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

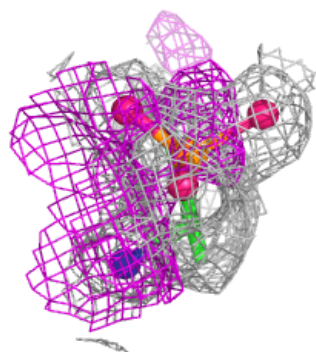
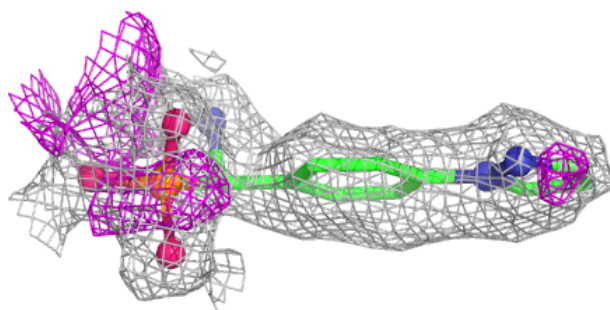
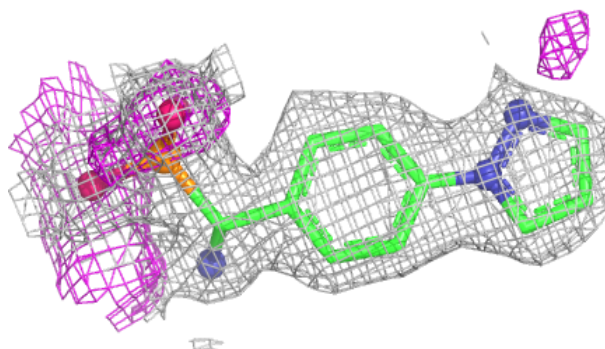


Electron density around 1OT B 1004:

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

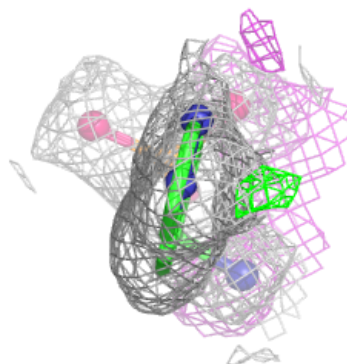
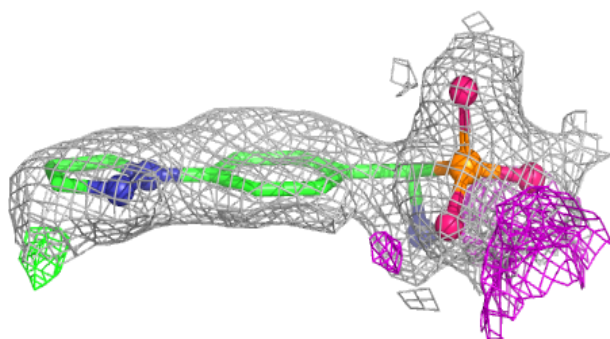
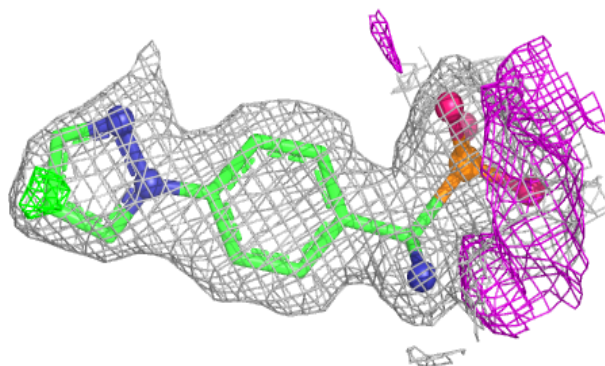
**Electron density around 1OT D 1004:**

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

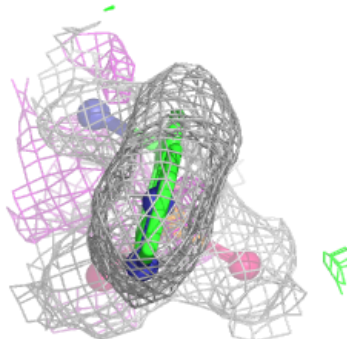
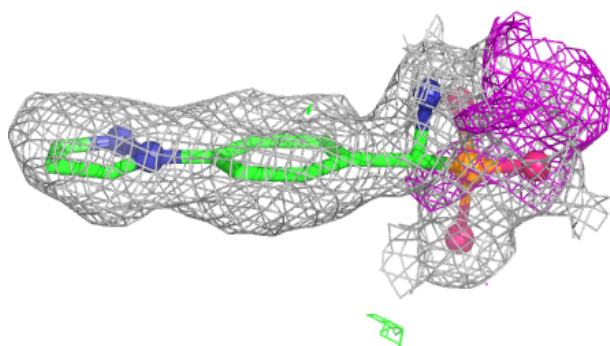
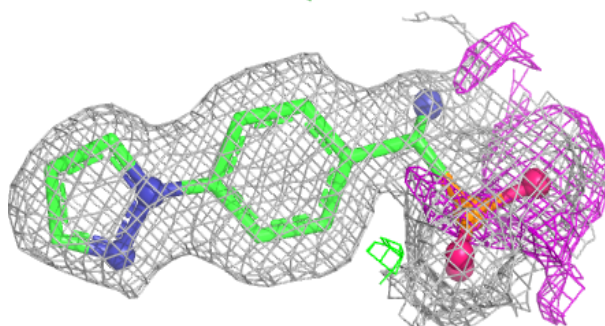


Electron density around 1OT E 1004:

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

**Electron density around 1OT A 1004:**

$2mF_o-DF_c$ (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.