



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 07:44 AM UTC

PDB ID : 7L4B / pdb_0000714b
Title : Dihydropicolinate synthase (DHDPS) from *C. jejuni* with pyruvate bound in the active site in P1211 space group
Authors : Saran, S.; Sanders, D.A.R.
Deposited on : 2020-12-18
Resolution : 2.42 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

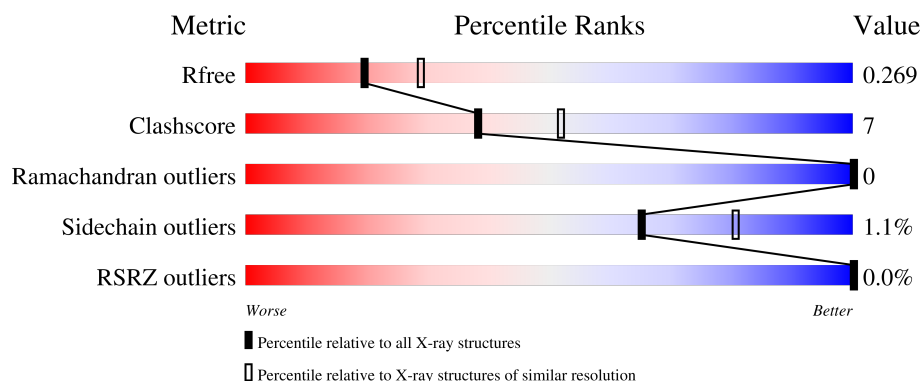
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	310	
1	B	310	
1	C	310	
1	D	310	
1	E	310	

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Mol	Chain	Length	Quality of chain
1	F	310	 84% 11% • 5%
1	G	310	 78% 16% • 5%
1	H	310	 81% 14% • 5%
1	I	310	 77% 19% 5%
1	J	310	 79% 16% • 5%
1	K	310	 76% 17% • 5%
1	L	310	 79% 16% • 5%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACT	C	305	-	-	X	-
4	ACT	L	302	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 27966 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 4-hydroxy-tetrahydrodipicolinate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	0	0
			2278	1449	379	437	13			
1	B	296	Total	C	N	O	S	0	0	0
			2275	1447	378	437	13			
1	C	296	Total	C	N	O	S	0	0	0
			2267	1442	375	437	13			
1	D	296	Total	C	N	O	S	0	0	0
			2257	1434	373	437	13			
1	E	294	Total	C	N	O	S	0	0	0
			2262	1440	375	434	13			
1	F	296	Total	C	N	O	S	0	0	0
			2271	1444	377	437	13			
1	G	296	Total	C	N	O	S	0	0	0
			2271	1444	377	437	13			
1	H	296	Total	C	N	O	S	0	0	0
			2279	1450	379	437	13			
1	I	296	Total	C	N	O	S	0	0	0
			2275	1447	378	437	13			
1	J	296	Total	C	N	O	S	0	0	0
			2279	1450	379	437	13			
1	K	296	Total	C	N	O	S	0	0	0
			2271	1444	378	436	13			
1	L	296	Total	C	N	O	S	0	0	0
			2271	1444	377	437	13			

There are 144 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP Q9PPB4
A	-10	ARG	-	expression tag	UNP Q9PPB4
A	-9	GLY	-	expression tag	UNP Q9PPB4
A	-8	SER	-	expression tag	UNP Q9PPB4
A	-7	HIS	-	expression tag	UNP Q9PPB4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	HIS	-	expression tag	UNP Q9PPB4
A	-5	HIS	-	expression tag	UNP Q9PPB4
A	-4	HIS	-	expression tag	UNP Q9PPB4
A	-3	HIS	-	expression tag	UNP Q9PPB4
A	-2	HIS	-	expression tag	UNP Q9PPB4
A	-1	GLY	-	expression tag	UNP Q9PPB4
A	0	SER	-	expression tag	UNP Q9PPB4
B	-11	MET	-	expression tag	UNP Q9PPB4
B	-10	ARG	-	expression tag	UNP Q9PPB4
B	-9	GLY	-	expression tag	UNP Q9PPB4
B	-8	SER	-	expression tag	UNP Q9PPB4
B	-7	HIS	-	expression tag	UNP Q9PPB4
B	-6	HIS	-	expression tag	UNP Q9PPB4
B	-5	HIS	-	expression tag	UNP Q9PPB4
B	-4	HIS	-	expression tag	UNP Q9PPB4
B	-3	HIS	-	expression tag	UNP Q9PPB4
B	-2	HIS	-	expression tag	UNP Q9PPB4
B	-1	GLY	-	expression tag	UNP Q9PPB4
B	0	SER	-	expression tag	UNP Q9PPB4
C	-11	MET	-	expression tag	UNP Q9PPB4
C	-10	ARG	-	expression tag	UNP Q9PPB4
C	-9	GLY	-	expression tag	UNP Q9PPB4
C	-8	SER	-	expression tag	UNP Q9PPB4
C	-7	HIS	-	expression tag	UNP Q9PPB4
C	-6	HIS	-	expression tag	UNP Q9PPB4
C	-5	HIS	-	expression tag	UNP Q9PPB4
C	-4	HIS	-	expression tag	UNP Q9PPB4
C	-3	HIS	-	expression tag	UNP Q9PPB4
C	-2	HIS	-	expression tag	UNP Q9PPB4
C	-1	GLY	-	expression tag	UNP Q9PPB4
C	0	SER	-	expression tag	UNP Q9PPB4
D	-11	MET	-	expression tag	UNP Q9PPB4
D	-10	ARG	-	expression tag	UNP Q9PPB4
D	-9	GLY	-	expression tag	UNP Q9PPB4
D	-8	SER	-	expression tag	UNP Q9PPB4
D	-7	HIS	-	expression tag	UNP Q9PPB4
D	-6	HIS	-	expression tag	UNP Q9PPB4
D	-5	HIS	-	expression tag	UNP Q9PPB4
D	-4	HIS	-	expression tag	UNP Q9PPB4
D	-3	HIS	-	expression tag	UNP Q9PPB4
D	-2	HIS	-	expression tag	UNP Q9PPB4
D	-1	GLY	-	expression tag	UNP Q9PPB4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	0	SER	-	expression tag	UNP Q9PPB4
E	-11	MET	-	expression tag	UNP Q9PPB4
E	-10	ARG	-	expression tag	UNP Q9PPB4
E	-9	GLY	-	expression tag	UNP Q9PPB4
E	-8	SER	-	expression tag	UNP Q9PPB4
E	-7	HIS	-	expression tag	UNP Q9PPB4
E	-6	HIS	-	expression tag	UNP Q9PPB4
E	-5	HIS	-	expression tag	UNP Q9PPB4
E	-4	HIS	-	expression tag	UNP Q9PPB4
E	-3	HIS	-	expression tag	UNP Q9PPB4
E	-2	HIS	-	expression tag	UNP Q9PPB4
E	-1	GLY	-	expression tag	UNP Q9PPB4
E	0	SER	-	expression tag	UNP Q9PPB4
F	-11	MET	-	expression tag	UNP Q9PPB4
F	-10	ARG	-	expression tag	UNP Q9PPB4
F	-9	GLY	-	expression tag	UNP Q9PPB4
F	-8	SER	-	expression tag	UNP Q9PPB4
F	-7	HIS	-	expression tag	UNP Q9PPB4
F	-6	HIS	-	expression tag	UNP Q9PPB4
F	-5	HIS	-	expression tag	UNP Q9PPB4
F	-4	HIS	-	expression tag	UNP Q9PPB4
F	-3	HIS	-	expression tag	UNP Q9PPB4
F	-2	HIS	-	expression tag	UNP Q9PPB4
F	-1	GLY	-	expression tag	UNP Q9PPB4
F	0	SER	-	expression tag	UNP Q9PPB4
G	-11	MET	-	expression tag	UNP Q9PPB4
G	-10	ARG	-	expression tag	UNP Q9PPB4
G	-9	GLY	-	expression tag	UNP Q9PPB4
G	-8	SER	-	expression tag	UNP Q9PPB4
G	-7	HIS	-	expression tag	UNP Q9PPB4
G	-6	HIS	-	expression tag	UNP Q9PPB4
G	-5	HIS	-	expression tag	UNP Q9PPB4
G	-4	HIS	-	expression tag	UNP Q9PPB4
G	-3	HIS	-	expression tag	UNP Q9PPB4
G	-2	HIS	-	expression tag	UNP Q9PPB4
G	-1	GLY	-	expression tag	UNP Q9PPB4
G	0	SER	-	expression tag	UNP Q9PPB4
H	-11	MET	-	expression tag	UNP Q9PPB4
H	-10	ARG	-	expression tag	UNP Q9PPB4
H	-9	GLY	-	expression tag	UNP Q9PPB4
H	-8	SER	-	expression tag	UNP Q9PPB4
H	-7	HIS	-	expression tag	UNP Q9PPB4

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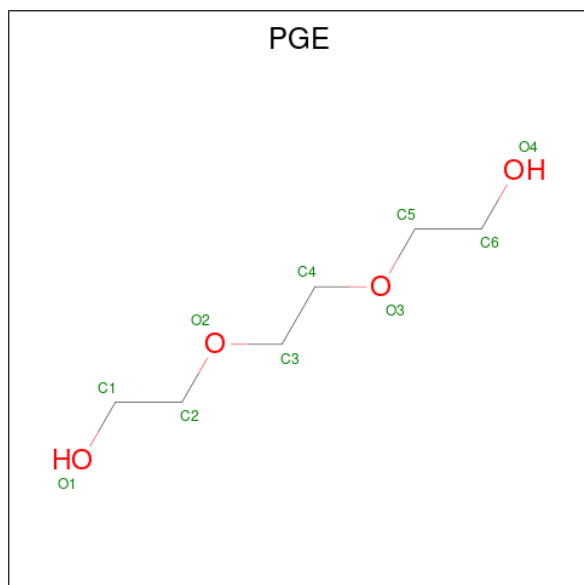
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H	-5	HIS	-	expression tag	UNP Q9PPB4
H	-4	HIS	-	expression tag	UNP Q9PPB4
H	-3	HIS	-	expression tag	UNP Q9PPB4
H	-2	HIS	-	expression tag	UNP Q9PPB4
H	-1	GLY	-	expression tag	UNP Q9PPB4
H	0	SER	-	expression tag	UNP Q9PPB4
I	-11	MET	-	expression tag	UNP Q9PPB4
I	-10	ARG	-	expression tag	UNP Q9PPB4
I	-9	GLY	-	expression tag	UNP Q9PPB4
I	-8	SER	-	expression tag	UNP Q9PPB4
I	-7	HIS	-	expression tag	UNP Q9PPB4
I	-6	HIS	-	expression tag	UNP Q9PPB4
I	-5	HIS	-	expression tag	UNP Q9PPB4
I	-4	HIS	-	expression tag	UNP Q9PPB4
I	-3	HIS	-	expression tag	UNP Q9PPB4
I	-2	HIS	-	expression tag	UNP Q9PPB4
I	-1	GLY	-	expression tag	UNP Q9PPB4
I	0	SER	-	expression tag	UNP Q9PPB4
J	-11	MET	-	expression tag	UNP Q9PPB4
J	-10	ARG	-	expression tag	UNP Q9PPB4
J	-9	GLY	-	expression tag	UNP Q9PPB4
J	-8	SER	-	expression tag	UNP Q9PPB4
J	-7	HIS	-	expression tag	UNP Q9PPB4
J	-6	HIS	-	expression tag	UNP Q9PPB4
J	-5	HIS	-	expression tag	UNP Q9PPB4
J	-4	HIS	-	expression tag	UNP Q9PPB4
J	-3	HIS	-	expression tag	UNP Q9PPB4
J	-2	HIS	-	expression tag	UNP Q9PPB4
J	-1	GLY	-	expression tag	UNP Q9PPB4
J	0	SER	-	expression tag	UNP Q9PPB4
K	-11	MET	-	expression tag	UNP Q9PPB4
K	-10	ARG	-	expression tag	UNP Q9PPB4
K	-9	GLY	-	expression tag	UNP Q9PPB4
K	-8	SER	-	expression tag	UNP Q9PPB4
K	-7	HIS	-	expression tag	UNP Q9PPB4
K	-6	HIS	-	expression tag	UNP Q9PPB4
K	-5	HIS	-	expression tag	UNP Q9PPB4
K	-4	HIS	-	expression tag	UNP Q9PPB4
K	-3	HIS	-	expression tag	UNP Q9PPB4
K	-2	HIS	-	expression tag	UNP Q9PPB4
K	-1	GLY	-	expression tag	UNP Q9PPB4

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Chain	Residue	Modelled	Actual	Comment	Reference
K	0	SER	-	expression tag	UNP Q9PPB4
L	-11	MET	-	expression tag	UNP Q9PPB4
L	-10	ARG	-	expression tag	UNP Q9PPB4
L	-9	GLY	-	expression tag	UNP Q9PPB4
L	-8	SER	-	expression tag	UNP Q9PPB4
L	-7	HIS	-	expression tag	UNP Q9PPB4
L	-6	HIS	-	expression tag	UNP Q9PPB4
L	-5	HIS	-	expression tag	UNP Q9PPB4
L	-4	HIS	-	expression tag	UNP Q9PPB4
L	-3	HIS	-	expression tag	UNP Q9PPB4
L	-2	HIS	-	expression tag	UNP Q9PPB4
L	-1	GLY	-	expression tag	UNP Q9PPB4
L	0	SER	-	expression tag	UNP Q9PPB4

- Molecule 2 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	6	4		
2	B	1	Total	C	O	0	0
			10	6	4		
2	C	1	Total	C	O	0	0
			10	6	4		
2	C	1	Total	C	O	0	0
			10	6	4		

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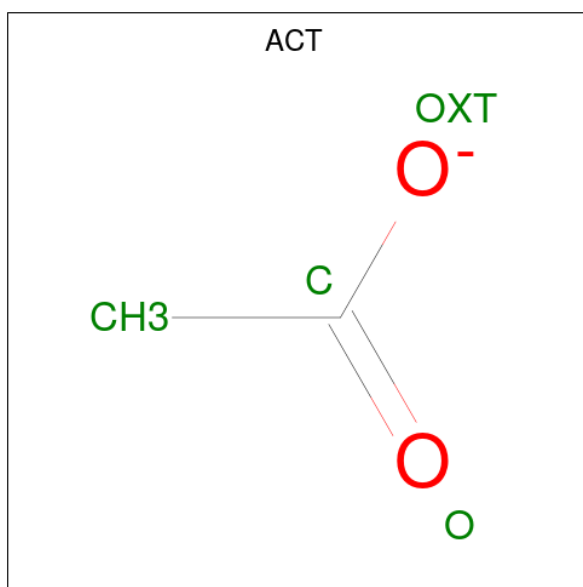
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total C O 10 6 4	0	0
2	E	1	Total C O 10 6 4	0	0
2	E	1	Total C O 10 6 4	0	0
2	F	1	Total C O 10 6 4	0	0
2	G	1	Total C O 10 6 4	0	0
2	H	1	Total C O 10 6 4	0	0
2	I	1	Total C O 10 6 4	0	0
2	J	1	Total C O 10 6 4	0	0
2	K	1	Total C O 10 6 4	0	0
2	L	1	Total C O 10 6 4	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

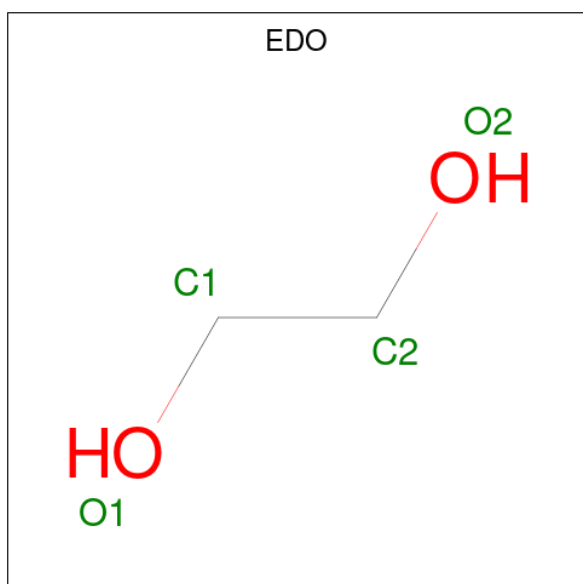
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	C	2	Total Mg 2 2	0	0
3	D	1	Total Mg 1 1	0	0
3	E	1	Total Mg 1 1	0	0
3	F	1	Total Mg 1 1	0	0
3	J	1	Total Mg 1 1	0	0

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		
4	L	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	J	1	Total	C	O	0	0
			4	2	2		

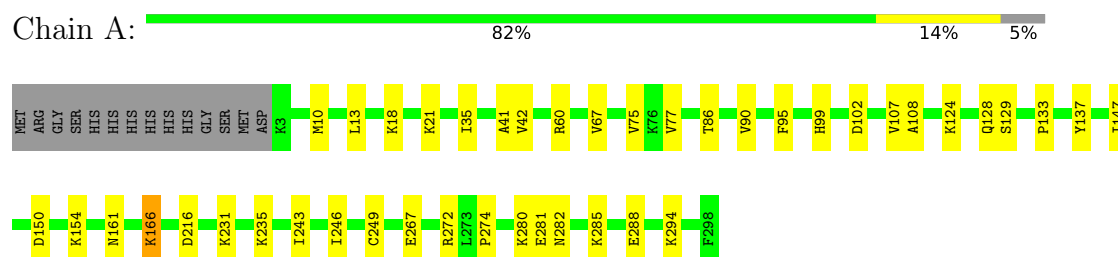
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	54	Total	O	0	0
			54	54		
6	B	64	Total	O	0	0
			64	64		
6	C	43	Total	O	0	0
			43	43		
6	D	54	Total	O	0	0
			54	54		
6	E	38	Total	O	0	0
			38	38		
6	F	35	Total	O	0	0
			35	35		
6	G	53	Total	O	0	0
			53	53		
6	H	44	Total	O	0	0
			44	44		
6	I	47	Total	O	0	0
			47	47		
6	J	32	Total	O	0	0
			32	32		
6	K	46	Total	O	0	0
			46	46		
6	L	37	Total	O	0	0
			37	37		

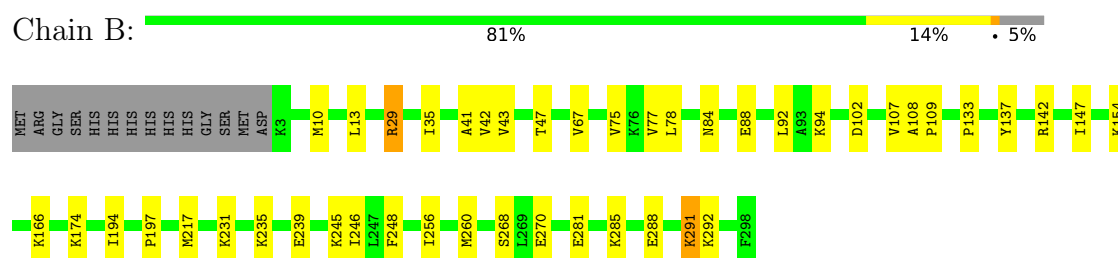
3 Residue-property plots

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.

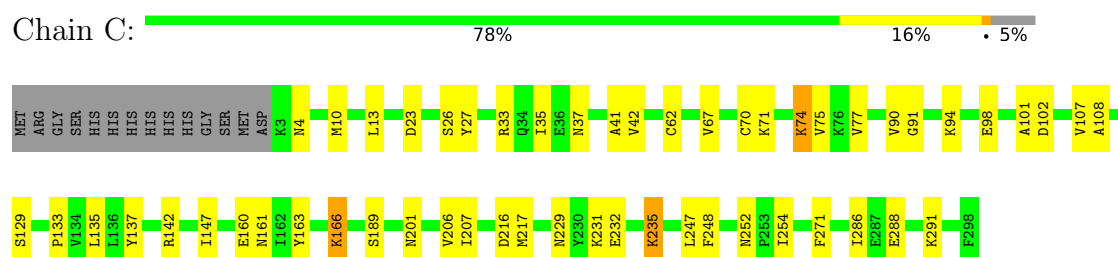
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase



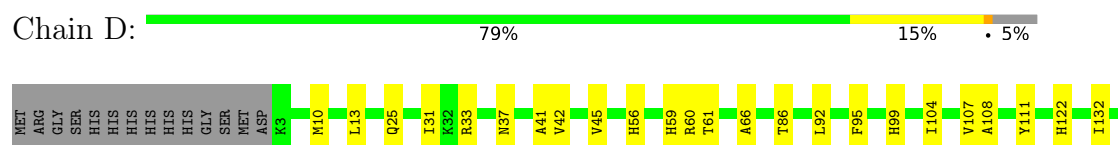
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

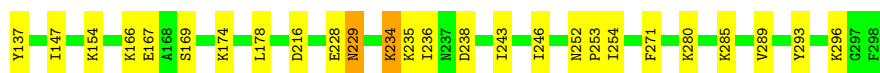


- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase



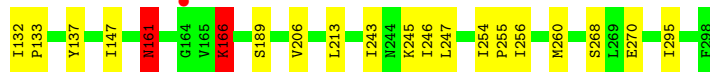
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase





- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain E: 81% 13% • 5%



- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain F: 84% 11% • 5%



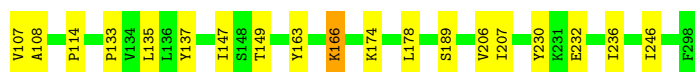
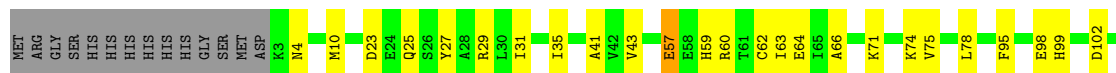
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain G: 78% 16% • 5%



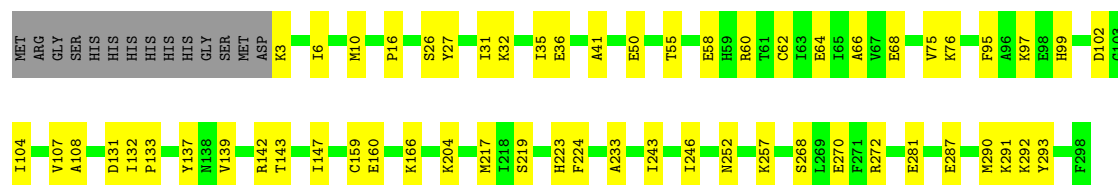
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain H: 81% 14% • 5%



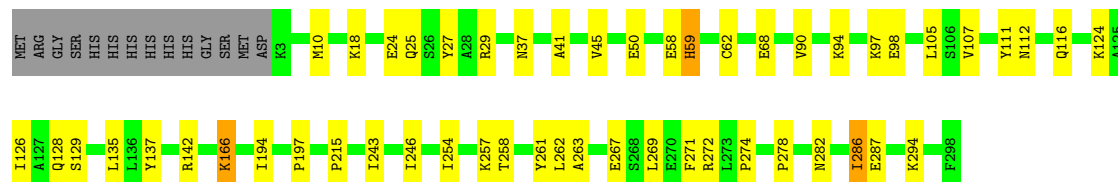
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain I: 77% 19% 5%



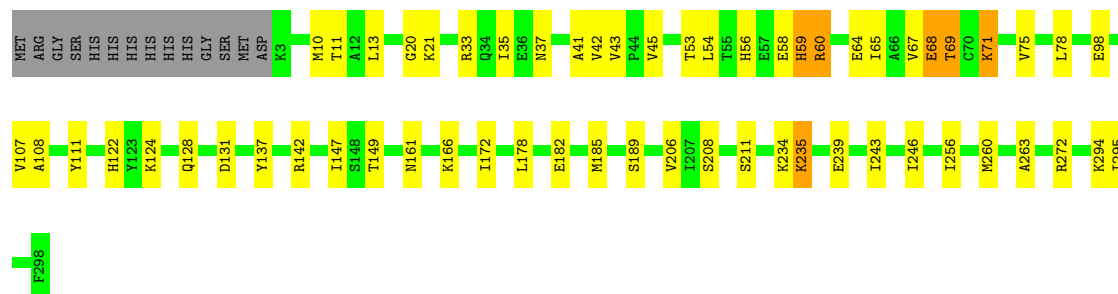
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain J: 79% 16% • 5%



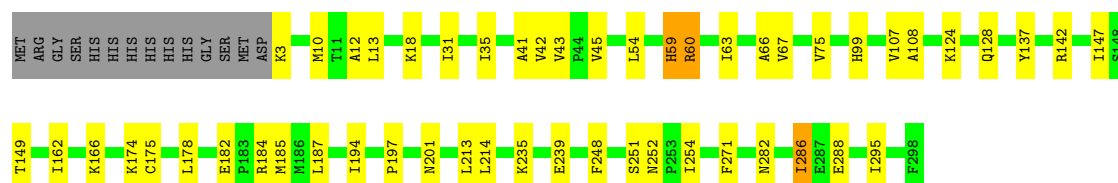
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain K: 76% 17% • 5%



- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain L: 79% 16% • 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.51Å 200.36Å 122.24Å 90.00° 109.83° 90.00°	Depositor
Resolution (Å)	49.15 – 2.42 49.15 – 2.42	Depositor EDS
% Data completeness (in resolution range)	98.5 (49.15-2.42) 88.1 (49.15-2.42)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 2.42Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.225 , 0.268 0.229 , 0.269	Depositor DCC
R_{free} test set	7072 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	41.7	Xtriage
Anisotropy	0.515	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.369 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	27966	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% 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: PGE, MG, ACT, EDO, KPI

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	0.25	0/2300	0.55	2/3109 (0.1%)
1	B	0.41	0/2297	0.54	1/3106 (0.0%)
1	C	0.60	1/2291 (0.0%)	0.60	4/3101 (0.1%)
1	D	0.58	4/2281 (0.2%)	0.57	2/3091 (0.1%)
1	E	0.48	1/2286 (0.0%)	0.60	2/3091 (0.1%)
1	F	0.38	0/2295	0.54	2/3105 (0.1%)
1	G	0.62	2/2295 (0.1%)	0.59	5/3105 (0.2%)
1	H	0.44	0/2303	0.54	0/3113
1	I	0.52	0/2299	0.56	4/3109 (0.1%)
1	J	0.48	0/2303	0.65	7/3113 (0.2%)
1	K	0.57	0/2295	0.57	1/3103 (0.0%)
1	L	0.62	0/2295	0.66	3/3105 (0.1%)
All	All	0.51	8/27540 (0.0%)	0.58	33/37251 (0.1%)

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	A	0	1
1	E	0	1
1	G	0	1
All	All	0	3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	295	ILE	C-O	-5.73	1.17	1.24
1	G	77	VAL	C-O	-5.72	1.18	1.24
1	E	254	ILE	CA-CB	-5.46	1.51	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	286	ILE	C-O	-5.37	1.17	1.24
1	D	61	THR	C-O	-5.24	1.17	1.24
1	D	235	LYS	C-O	-5.17	1.17	1.24
1	D	234	LYS	C-O	-5.07	1.18	1.24
1	D	33	ARG	C-O	-5.00	1.18	1.24

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	286	ILE	N-CA-C	-10.56	100.60	110.53
1	J	286	ILE	N-CA-C	-8.60	102.03	110.72
1	E	161	ASN	N-CA-C	8.14	122.78	113.01
1	J	263	ALA	N-CA-C	-7.70	103.47	113.17
1	G	57	GLU	N-CA-C	-7.63	101.84	112.45
1	B	291	LYS	N-CA-C	7.47	121.31	111.75
1	F	71	LYS	CG-CD-CE	6.34	125.88	111.30
1	L	59	HIS	N-CA-C	-6.11	104.53	111.07
1	I	159	CYS	CA-C-N	6.09	128.93	120.29
1	I	159	CYS	C-N-CA	6.09	128.93	120.29
1	G	277	SER	CA-C-N	-5.87	114.70	120.98
1	G	277	SER	C-N-CA	-5.87	114.70	120.98
1	L	288	GLU	N-CA-C	5.84	117.65	111.28
1	D	236	ILE	N-CA-C	-5.70	104.97	110.72
1	A	281	GLU	CB-CA-C	-5.69	101.87	110.92
1	E	73	THR	N-CA-C	-5.63	101.19	109.81
1	I	160	GLU	CB-CG-CD	-5.59	103.10	112.60
1	C	160	GLU	N-CA-C	5.58	118.12	111.71
1	I	281	GLU	CB-CG-CD	5.45	121.86	112.60
1	G	278	PRO	N-CA-C	5.42	119.13	110.40
1	J	287	GLU	CA-C-N	-5.40	112.15	121.66
1	J	287	GLU	C-N-CA	-5.40	112.15	121.66
1	K	71	LYS	CG-CD-CE	5.25	123.38	111.30
1	F	285	LYS	CB-CG-CD	5.21	123.29	111.30
1	J	24	GLU	N-CA-C	5.16	118.68	112.38
1	A	281	GLU	N-CA-CB	5.12	117.60	109.82
1	G	285	LYS	CB-CG-CD	-5.07	99.64	111.30
1	J	29	ARG	CB-CG-CD	5.06	122.94	111.30
1	C	4	ASN	CB-CA-C	5.05	118.08	109.80
1	J	68	GLU	CB-CA-C	-5.04	102.28	110.85
1	D	229	ASN	N-CA-C	-5.04	102.61	109.71
1	C	74	LYS	CA-CB-CG	5.03	124.16	114.10
1	C	4	ASN	N-CA-CB	-5.02	101.88	110.16

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	166	KPI	Mainchain
1	E	166	KPI	Mainchain
1	G	166	KPI	Mainchain

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	2278	0	2317	33	0
1	B	2275	0	2308	30	0
1	C	2267	0	2283	39	0
1	D	2257	0	2255	32	0
1	E	2262	0	2299	32	0
1	F	2271	0	2297	23	0
1	G	2271	0	2297	32	0
1	H	2279	0	2319	33	0
1	I	2275	0	2308	34	0
1	J	2279	0	2319	34	0
1	K	2271	0	2297	42	0
1	L	2271	0	2297	35	0
2	A	10	0	14	3	0
2	B	10	0	14	1	0
2	C	20	0	28	0	0
2	D	10	0	14	1	0
2	E	20	0	28	0	0
2	F	10	0	14	0	0
2	G	10	0	14	1	0
2	H	10	0	14	0	0
2	I	10	0	14	0	0
2	J	10	0	14	1	0
2	K	10	0	14	0	0
2	L	10	0	14	0	0
3	A	1	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	J	1	0	0	0	0
4	C	4	0	3	5	0
4	F	4	0	3	0	0
4	L	4	0	3	4	0
5	J	4	0	6	0	0
6	A	54	0	0	0	0
6	B	64	0	0	2	0
6	C	43	0	0	0	0
6	D	54	0	0	1	0
6	E	38	0	0	0	0
6	F	35	0	0	2	0
6	G	53	0	0	1	0
6	H	44	0	0	1	0
6	I	47	0	0	3	0
6	J	32	0	0	0	0
6	K	46	0	0	1	0
6	L	37	0	0	0	0
All	All	27966	0	27807	382	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 7.

All (382) 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:E:60:ARG:HB2	1:E:95:PHE:CZ	1.86	1.10
1:E:60:ARG:HB2	1:E:95:PHE:HZ	1.16	1.06
1:J:45:VAL:HG11	1:J:59:HIS:CE1	2.07	0.90
1:K:65:ILE:O	1:K:69:THR:OG1	1.90	0.88
1:K:124:LYS:O	1:K:128:GLN:HG3	1.74	0.87
1:I:64:GLU:O	1:I:68:GLU:HG3	1.79	0.81
1:H:60:ARG:HB2	1:H:95:PHE:HZ	1.46	0.79
1:H:107:VAL:HA	1:H:137:TYR:HB3	1.68	0.76
1:L:60:ARG:HG2	1:L:99:HIS:CE1	2.19	0.76
1:I:107:VAL:HA	1:I:137:TYR:HB3	1.68	0.76
1:J:258:THR:HG22	1:J:269:LEU:HD11	1.68	0.75
1:K:56:HIS:O	1:K:60:ARG:HD3	1.86	0.75
1:E:60:ARG:CB	1:E:95:PHE:HZ	1.97	0.74
1:J:282:ASN:O	1:J:286:ILE:HD12	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:107:VAL:HA	1:C:137:TYR:HB3	1.70	0.74
1:G:288:GLU:HA	1:G:291:LYS:HD2	1.69	0.74
1:F:107:VAL:HA	1:F:137:TYR:HB3	1.69	0.73
1:L:251:SER:HA	4:L:302:ACT:H3	1.69	0.73
1:J:107:VAL:HA	1:J:137:TYR:HB3	1.70	0.73
1:K:67:VAL:O	1:K:71:LYS:HG3	1.88	0.72
1:K:208:SER:HB3	1:K:211:SER:HB2	1.72	0.72
1:L:235:LYS:O	1:L:239:GLU:HG3	1.89	0.72
1:D:60:ARG:HB2	1:D:95:PHE:CZ	2.24	0.72
1:C:67:VAL:HG22	1:C:77:VAL:HG21	1.73	0.71
1:K:107:VAL:HA	1:K:137:TYR:HB3	1.73	0.71
1:H:35:ILE:HG12	1:H:75:VAL:HG21	1.72	0.71
1:G:142:ARG:NH2	6:G:401:HOH:O	2.23	0.71
1:L:107:VAL:HA	1:L:137:TYR:HB3	1.73	0.70
1:L:60:ARG:CD	1:L:99:HIS:CE1	2.74	0.70
1:E:107:VAL:HA	1:E:137:TYR:HB3	1.73	0.69
1:H:60:ARG:HB2	1:H:95:PHE:CZ	2.27	0.69
1:L:60:ARG:CG	1:L:99:HIS:CE1	2.75	0.69
1:A:107:VAL:HA	1:A:137:TYR:HB3	1.75	0.69
1:B:288:GLU:O	1:B:291:LYS:HB2	1.92	0.69
1:I:35:ILE:HG12	1:I:75:VAL:HG21	1.74	0.69
1:B:107:VAL:HA	1:B:137:TYR:HB3	1.74	0.68
1:C:252:ASN:H	4:C:305:ACT:H3	1.59	0.67
1:D:107:VAL:HA	1:D:137:TYR:HB3	1.76	0.67
1:D:60:ARG:HB2	1:D:95:PHE:HZ	1.57	0.67
1:K:33:ARG:O	1:K:37:ASN:ND2	2.27	0.67
1:G:288:GLU:OE1	1:G:291:LYS:NZ	2.29	0.66
1:G:107:VAL:HA	1:G:137:TYR:HB3	1.76	0.66
1:A:108:ALA:HB2	1:A:147:ILE:HD11	1.79	0.65
1:A:67:VAL:HA	1:A:77:VAL:HG21	1.79	0.65
1:K:235:LYS:O	1:K:239:GLU:HG3	1.98	0.64
1:K:263:ALA:HB1	1:K:295:ILE:HG21	1.79	0.64
1:D:154:LYS:HD2	2:D:301:PGE:H4	1.80	0.64
1:K:35:ILE:HG12	1:K:75:VAL:HG21	1.79	0.64
1:I:55:THR:OG1	1:I:58:GLU:HG3	1.98	0.63
1:F:138:ASN:ND2	1:F:167:GLU:OE1	2.32	0.63
1:E:60:ARG:HB2	1:E:95:PHE:CE1	2.32	0.63
1:C:189:SER:HB3	1:C:206:VAL:HG12	1.81	0.62
1:A:60:ARG:HB2	1:A:95:PHE:CZ	2.33	0.62
1:B:142:ARG:NH2	6:B:405:HOH:O	2.32	0.62
1:A:124:LYS:HG2	1:A:128:GLN:NE2	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:33:ARG:O	1:C:37:ASN:ND2	2.33	0.62
1:A:95:PHE:O	1:A:99:HIS:ND1	2.33	0.62
1:C:235:LYS:O	1:C:235:LYS:HD3	2.00	0.62
1:K:108:ALA:HB2	1:K:147:ILE:HD11	1.81	0.61
1:F:246:ILE:HD11	1:F:285:LYS:HG2	1.82	0.61
1:K:58:GLU:OE1	1:K:272:ARG:NH2	2.33	0.61
1:G:154:LYS:HD3	2:G:301:PGE:H62	1.83	0.61
1:D:108:ALA:HB2	1:D:147:ILE:HD11	1.82	0.61
1:D:296:LYS:NZ	6:D:401:HOH:O	2.33	0.61
1:G:288:GLU:CD	1:G:291:LYS:HZ3	2.09	0.61
1:J:45:VAL:CG1	1:J:59:HIS:CE1	2.83	0.61
1:F:31:ILE:O	1:F:35:ILE:HG12	2.01	0.60
1:E:21:LYS:NZ	1:E:24:GLU:OE2	2.33	0.60
1:K:45:VAL:HG11	1:K:59:HIS:CE1	2.36	0.60
1:B:13:LEU:HD11	1:B:42:VAL:HB	1.84	0.60
1:C:77:VAL:HG23	1:C:102:ASP:H	1.67	0.60
1:E:268:SER:OG	1:E:270:GLU:HG3	2.02	0.60
1:K:149:THR:HG23	1:K:178:LEU:HD23	1.82	0.59
1:K:294:LYS:NZ	6:K:404:HOH:O	2.34	0.59
1:A:13:LEU:HD11	1:A:42:VAL:HB	1.83	0.59
1:F:32:LYS:O	1:F:36:GLU:HG3	2.03	0.59
1:I:60:ARG:HB2	1:I:95:PHE:HZ	1.68	0.59
1:K:20:GLY:C	1:K:21:LYS:HD2	2.28	0.59
1:B:217:MET:HE1	1:B:239:GLU:OE2	2.03	0.59
1:C:288:GLU:OE2	1:C:291:LYS:CE	2.50	0.58
1:J:254:ILE:O	1:J:258:THR:HG23	2.04	0.58
1:I:97:LYS:HE2	1:I:131:ASP:OD1	2.02	0.58
1:J:37:ASN:HD22	1:J:215:PRO:HG2	1.69	0.58
1:J:258:THR:HG21	1:J:278:PRO:HG3	1.85	0.57
1:L:282:ASN:O	1:L:286:ILE:HG13	2.05	0.57
1:B:94:LYS:NZ	6:B:408:HOH:O	2.37	0.57
1:L:182:GLU:OE2	1:L:184:ARG:NH1	2.38	0.57
1:C:248:PHE:HD2	4:C:305:ACT:H2	1.70	0.57
1:I:60:ARG:HB2	1:I:95:PHE:CZ	2.40	0.57
1:A:35:ILE:HG12	1:A:75:VAL:HG21	1.86	0.56
1:G:137:TYR:CZ	1:G:166:KPI:HEA	2.40	0.56
1:H:60:ARG:NE	1:H:99:HIS:CE1	2.73	0.56
1:I:10:MET:HG2	1:I:41:ALA:HB3	1.88	0.56
1:J:18:LYS:HE3	1:J:267:GLU:OE1	2.06	0.56
1:B:47:THR:OG1	1:D:111:TYR:OH	2.23	0.56
1:C:201:ASN:O	1:D:234:LYS:CE	2.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:166:KPI:HDA	1:H:207:ILE:HB	1.87	0.56
1:J:58:GLU:OE2	1:J:272:ARG:NH2	2.34	0.56
1:B:281:GLU:O	1:B:285:LYS:HG3	2.06	0.56
1:K:45:VAL:HG13	1:K:54:LEU:HD12	1.88	0.56
1:F:76:LYS:NZ	6:F:401:HOH:O	2.36	0.55
1:G:230:TYR:CD2	1:H:230:TYR:HD2	2.24	0.55
1:J:126:ILE:O	1:J:129:SER:OG	2.24	0.55
1:C:90:VAL:HG13	1:C:129:SER:HB2	1.88	0.55
1:C:201:ASN:O	1:D:234:LYS:HE2	2.06	0.55
1:F:37:ASN:OD1	1:F:296:LYS:HD2	2.06	0.55
1:G:146:GLU:OE2	1:G:174:LYS:NZ	2.34	0.55
1:K:13:LEU:HD11	1:K:42:VAL:HB	1.87	0.55
1:B:231:LYS:H	1:B:231:LYS:HD2	1.70	0.55
1:E:108:ALA:HB2	1:E:147:ILE:HD11	1.88	0.55
1:J:112:ASN:HA	1:K:142:ARG:HH12	1.72	0.55
1:L:248:PHE:HD2	4:L:302:ACT:H2	1.72	0.55
1:H:74:LYS:NZ	6:H:401:HOH:O	2.26	0.55
1:C:35:ILE:HG12	1:C:75:VAL:HG21	1.90	0.54
1:C:27:TYR:CE2	1:C:62:CYS:HB3	2.43	0.54
1:G:10:MET:HG2	1:G:41:ALA:HB3	1.89	0.54
1:H:43:VAL:HG22	1:H:78:LEU:HB3	1.90	0.54
1:E:29:ARG:HA	1:E:32:LYS:HG3	1.89	0.54
1:K:10:MET:HG2	1:K:41:ALA:HB3	1.89	0.54
1:F:124:LYS:O	1:F:128:GLN:HG3	2.08	0.54
1:H:108:ALA:HB2	1:H:147:ILE:HD11	1.90	0.54
1:A:150:ASP:HB3	2:A:301:PGE:H3	1.89	0.53
1:E:60:ARG:HD3	1:E:99:HIS:CD2	2.43	0.53
1:G:288:GLU:CD	1:G:291:LYS:NZ	2.66	0.53
1:I:243:ILE:HA	1:I:246:ILE:HG22	1.91	0.53
1:E:29:ARG:O	1:E:32:LYS:HB2	2.08	0.53
1:J:50:GLU:OE2	1:J:257:LYS:NZ	2.39	0.53
1:E:10:MET:HG2	1:E:41:ALA:HB3	1.91	0.53
1:J:10:MET:HG2	1:J:41:ALA:HB3	1.90	0.52
1:K:53:THR:HB	1:K:272:ARG:HG2	1.91	0.52
1:B:108:ALA:HB2	1:B:147:ILE:HD11	1.91	0.52
1:L:35:ILE:HG12	1:L:75:VAL:HG21	1.92	0.52
1:K:234:LYS:HE2	1:L:201:ASN:O	2.09	0.52
1:C:248:PHE:CD2	4:C:305:ACT:H2	2.44	0.52
1:L:175:CYS:HB3	1:L:187:LEU:HD21	1.90	0.52
1:H:10:MET:HG2	1:H:41:ALA:HB3	1.91	0.52
1:D:174:LYS:HD3	1:D:178:LEU:HD11	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:149:THR:HG23	1:L:178:LEU:HD23	1.92	0.52
1:E:161:ASN:OD1	1:E:161:ASN:N	2.33	0.52
1:L:13:LEU:HD11	1:L:42:VAL:HB	1.92	0.52
1:H:31:ILE:HD13	1:H:66:ALA:HA	1.91	0.51
1:J:105:LEU:HD13	1:J:135:LEU:HD23	1.91	0.51
1:A:235:LYS:O	1:A:235:LYS:HD3	2.10	0.51
1:D:167:GLU:OE1	1:D:169:SER:OG	2.23	0.51
1:I:217:MET:HE1	1:I:293:TYR:HD1	1.74	0.51
1:B:29:ARG:HH11	1:B:29:ARG:HG3	1.75	0.51
1:G:58:GLU:OE1	1:G:272:ARG:NH1	2.36	0.51
1:A:60:ARG:NE	1:A:99:HIS:NE2	2.58	0.51
1:L:252:ASN:H	4:L:302:ACT:CH3	2.24	0.51
1:J:90:VAL:HG13	1:J:129:SER:HB3	1.93	0.51
1:G:256:ILE:O	1:G:260:MET:HG2	2.11	0.51
1:E:60:ARG:HD3	1:E:99:HIS:NE2	2.26	0.50
1:F:108:ALA:HB2	1:F:147:ILE:HD11	1.93	0.50
1:G:60:ARG:HG3	1:G:99:HIS:CE1	2.46	0.50
1:L:10:MET:HG2	1:L:41:ALA:HB3	1.94	0.50
1:I:108:ALA:HB2	1:I:147:ILE:HD11	1.93	0.50
1:L:213:LEU:HG	1:L:214:LEU:HG	1.92	0.50
1:A:10:MET:HG2	1:A:41:ALA:HB3	1.93	0.50
1:G:137:TYR:CE1	1:G:166:KPI:HEA	2.47	0.50
1:K:11:THR:OG1	1:K:211:SER:OG	2.28	0.50
1:B:35:ILE:HG12	1:B:75:VAL:HG21	1.94	0.50
1:B:10:MET:HG2	1:B:41:ALA:HB3	1.94	0.50
1:H:137:TYR:CE1	1:H:166:KPI:HEA	2.47	0.50
1:A:243:ILE:HA	1:A:246:ILE:HG22	1.92	0.49
1:E:243:ILE:HA	1:E:246:ILE:HG22	1.93	0.49
1:G:281:GLU:O	1:G:285:LYS:HB2	2.12	0.49
1:I:58:GLU:OE1	1:I:272:ARG:NH1	2.23	0.49
1:I:50:GLU:OE2	1:I:257:LYS:NZ	2.29	0.49
1:C:254:ILE:HA	1:C:271:PHE:CE2	2.47	0.49
1:K:208:SER:HB3	1:K:211:SER:CB	2.41	0.49
1:L:182:GLU:HB3	1:L:185:MET:HE3	1.95	0.49
1:G:175:CYS:HB3	1:G:187:LEU:HD21	1.95	0.49
1:E:213:LEU:HD11	1:E:295:ILE:HD13	1.94	0.49
1:L:108:ALA:HB2	1:L:147:ILE:HD11	1.93	0.49
1:C:161:ASN:OD1	1:C:161:ASN:N	2.34	0.49
1:K:58:GLU:CD	1:K:272:ARG:HH22	2.21	0.49
1:C:77:VAL:HG23	1:C:101:ALA:HA	1.95	0.49
1:A:18:LYS:HD3	1:A:267:GLU:OE2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:31:ILE:HD13	1:I:66:ALA:HA	1.94	0.48
1:C:13:LEU:HD11	1:C:42:VAL:HB	1.94	0.48
1:I:99:HIS:ND1	6:I:405:HOH:O	2.34	0.48
1:B:268:SER:HB3	1:B:270:GLU:HG3	1.96	0.48
1:I:104:ILE:HG23	1:I:132:ILE:HD11	1.95	0.48
1:K:21:LYS:HD2	1:K:21:LYS:N	2.28	0.48
1:A:154:LYS:HD2	2:A:301:PGE:H22	1.96	0.48
1:A:21:LYS:HD3	1:A:272:ARG:NH2	2.29	0.48
1:H:60:ARG:NH1	1:H:99:HIS:HE1	2.11	0.48
1:L:60:ARG:HD2	1:L:99:HIS:CE1	2.47	0.48
1:B:88:GLU:HG2	1:B:92:LEU:HD11	1.96	0.48
1:A:231:LYS:HE2	1:A:231:LYS:HB2	1.62	0.48
1:C:229:ASN:HB3	1:C:232:GLU:OE1	2.14	0.48
1:K:208:SER:CB	1:K:211:SER:HB2	2.42	0.48
1:F:293:TYR:C	1:F:294:LYS:HD2	2.38	0.48
1:D:10:MET:HG2	1:D:41:ALA:HB3	1.95	0.47
1:D:45:VAL:HG11	1:D:59:HIS:CE1	2.48	0.47
1:A:60:ARG:HB2	1:A:95:PHE:CE2	2.49	0.47
1:F:10:MET:HG2	1:F:41:ALA:HB3	1.97	0.47
1:K:64:GLU:O	1:K:68:GLU:CG	2.63	0.47
1:D:254:ILE:HA	1:D:271:PHE:CE2	2.49	0.47
1:E:245:LYS:HA	1:E:245:LYS:HD2	1.77	0.47
1:F:35:ILE:HG23	1:F:75:VAL:HG21	1.95	0.47
1:J:137:TYR:CE1	1:J:166:KPI:HEA	2.49	0.47
1:D:56:HIS:CE1	1:D:92:LEU:HD21	2.49	0.47
1:E:28:ALA:O	1:E:32:LYS:HG2	2.14	0.47
1:E:189:SER:HB3	1:E:206:VAL:HG12	1.96	0.47
1:J:94:LYS:HG3	1:J:129:SER:HB2	1.96	0.47
1:K:243:ILE:HA	1:K:246:ILE:HG22	1.97	0.47
1:D:86:THR:OG1	1:D:122:HIS:ND1	2.39	0.47
1:D:234:LYS:NZ	1:D:238:ASP:OD2	2.47	0.47
1:D:285:LYS:O	1:D:289:VAL:HG23	2.14	0.47
1:I:287:GLU:HA	1:I:290:MET:HE2	1.96	0.47
1:L:3:LYS:NZ	1:L:162:ILE:O	2.41	0.47
1:L:45:VAL:HG13	1:L:54:LEU:HD12	1.97	0.47
1:K:56:HIS:O	1:K:60:ARG:CD	2.59	0.47
1:L:12:ALA:HA	1:L:43:VAL:HB	1.96	0.47
1:E:74:LYS:H	1:E:74:LYS:HG3	1.55	0.47
1:A:90:VAL:HG13	1:A:129:SER:HB3	1.97	0.47
1:L:147:ILE:O	1:L:174:LYS:NZ	2.39	0.47
1:A:216:ASP:OD1	1:A:216:ASP:N	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:VAL:HA	1:B:77:VAL:HG21	1.96	0.46
1:B:102:ASP:O	1:B:133:PRO:HD2	2.14	0.46
1:F:243:ILE:HA	1:F:246:ILE:HG22	1.97	0.46
1:A:60:ARG:CZ	1:A:99:HIS:NE2	2.78	0.46
1:B:235:LYS:O	1:B:239:GLU:HB2	2.15	0.46
1:H:4:ASN:HB3	1:H:133:PRO:HG3	1.96	0.46
1:D:95:PHE:O	1:D:99:HIS:CD2	2.68	0.46
1:H:137:TYR:CZ	1:H:166:KPI:HEA	2.51	0.46
1:L:60:ARG:CG	1:L:99:HIS:HE1	2.27	0.46
1:C:90:VAL:O	1:C:94:LYS:HG3	2.15	0.46
1:H:60:ARG:CZ	1:H:99:HIS:CE1	2.98	0.46
1:E:113:LYS:CE	1:L:142:ARG:HG2	2.46	0.46
1:J:27:TYR:CZ	1:J:62:CYS:HB3	2.51	0.46
1:C:27:TYR:CZ	1:C:62:CYS:HB3	2.50	0.46
1:E:45:VAL:HG11	1:E:59:HIS:CE1	2.51	0.46
1:F:188:ILE:HG21	1:F:207:ILE:HG13	1.98	0.46
1:H:59:HIS:O	1:H:63:ILE:HG13	2.16	0.46
1:B:84:ASN:HA	1:B:109:PRO:HB3	1.98	0.46
1:G:16:PRO:HD2	1:G:27:TYR:HD1	1.80	0.46
1:H:27:TYR:CZ	1:H:62:CYS:HB3	2.51	0.46
1:H:27:TYR:CE2	1:H:62:CYS:HB3	2.51	0.46
1:F:86:THR:HA	1:F:122:HIS:CE1	2.51	0.45
1:I:6:ILE:HG23	1:I:76:LYS:HE2	1.98	0.45
1:I:268:SER:OG	1:I:270:GLU:HG3	2.17	0.45
1:J:294:LYS:HA	1:J:294:LYS:HD3	1.75	0.45
1:E:95:PHE:O	1:E:95:PHE:CD1	2.69	0.45
1:H:60:ARG:CZ	1:H:99:HIS:HE1	2.30	0.45
1:J:111:TYR:O	1:K:142:ARG:NH1	2.49	0.45
1:A:124:LYS:HG2	1:A:128:GLN:HE21	1.82	0.45
1:B:47:THR:HG1	1:D:111:TYR:HH	1.55	0.45
1:C:23:ASP:CG	1:C:26:SER:HB3	2.42	0.45
1:E:247:LEU:HA	1:E:255:PRO:HB2	1.99	0.45
1:F:13:LEU:HD11	1:F:42:VAL:HB	1.98	0.45
1:I:287:GLU:HA	1:I:290:MET:CE	2.47	0.45
1:C:102:ASP:O	1:C:133:PRO:HD2	2.17	0.45
1:H:10:MET:HA	1:H:41:ALA:O	2.16	0.45
1:H:147:ILE:O	1:H:174:LYS:NZ	2.44	0.45
1:H:189:SER:HB3	1:H:206:VAL:HG12	1.98	0.45
1:A:294:LYS:HA	1:A:294:LYS:HD3	1.74	0.45
1:C:235:LYS:HD3	1:C:235:LYS:C	2.37	0.45
1:L:248:PHE:CD2	4:L:302:ACT:H2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:57:GLU:H	1:H:57:GLU:HG2	1.53	0.45
1:E:12:ALA:HA	1:E:43:VAL:HB	1.98	0.44
1:K:189:SER:HB3	1:K:206:VAL:HG12	1.98	0.44
1:J:10:MET:HA	1:J:41:ALA:O	2.17	0.44
1:A:86:THR:O	1:A:90:VAL:HG23	2.16	0.44
1:A:246:ILE:HD11	1:A:285:LYS:HB3	1.99	0.44
1:D:92:LEU:HA	1:D:95:PHE:HB3	1.99	0.44
1:G:245:LYS:HA	1:G:245:LYS:HD2	1.87	0.44
1:A:280:LYS:HD2	1:A:280:LYS:N	2.33	0.44
1:C:10:MET:HG2	1:C:41:ALA:HB3	1.99	0.44
1:F:243:ILE:HB	1:F:293:TYR:CE2	2.53	0.44
1:G:292:LYS:HE2	1:G:293:TYR:CZ	2.52	0.44
1:J:269:LEU:HD21	1:J:271:PHE:HE1	1.81	0.44
1:C:91:GLY:HA2	1:C:94:LYS:HE3	1.99	0.44
1:C:252:ASN:H	4:C:305:ACT:CH3	2.29	0.44
1:D:243:ILE:HB	1:D:293:TYR:CE2	2.53	0.44
1:F:229:ASN:HB2	6:F:434:HOH:O	2.17	0.44
1:G:13:LEU:HD11	1:G:42:VAL:HB	1.99	0.44
1:D:104:ILE:HG23	1:D:132:ILE:HD11	1.98	0.44
1:I:16:PRO:HD2	1:I:27:TYR:HD1	1.81	0.44
1:B:29:ARG:HG3	1:B:29:ARG:NH1	2.33	0.44
1:G:254:ILE:HA	1:G:271:PHE:CE2	2.52	0.44
1:L:63:ILE:O	1:L:67:VAL:HG23	2.18	0.44
1:A:154:LYS:NZ	2:A:301:PGE:H62	2.32	0.44
1:C:108:ALA:HB2	1:C:147:ILE:HD11	1.98	0.44
1:C:231:LYS:HE3	1:D:228:GLU:OE1	2.17	0.44
1:B:147:ILE:O	1:B:174:LYS:NZ	2.50	0.44
1:I:219:SER:O	1:I:223:HIS:ND1	2.51	0.44
1:K:64:GLU:O	1:K:68:GLU:HG2	2.18	0.43
1:C:142:ARG:HH21	4:C:305:ACT:H3	1.83	0.43
1:E:29:ARG:HA	1:E:32:LYS:CG	2.48	0.43
1:G:294:LYS:HA	1:G:294:LYS:HD3	1.72	0.43
1:H:149:THR:HG23	1:H:178:LEU:HD23	1.99	0.43
1:J:243:ILE:HA	1:J:246:ILE:HG22	2.00	0.43
1:K:256:ILE:O	1:K:260:MET:HG2	2.18	0.43
1:C:247:LEU:HD23	1:C:247:LEU:HA	1.87	0.43
1:G:166:KPI:HDA	1:G:207:ILE:HB	1.99	0.43
1:I:139:VAL:HG13	1:I:143:THR:HG23	2.00	0.43
1:C:135:LEU:HD13	1:C:163:TYR:CZ	2.54	0.43
1:D:31:ILE:HD13	1:D:66:ALA:HA	2.00	0.43
1:I:292:LYS:O	1:I:292:LYS:HG3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:261:TYR:O	1:J:261:TYR:CG	2.70	0.43
1:D:13:LEU:HD11	1:D:42:VAL:HB	2.01	0.43
1:D:252:ASN:OD1	1:D:253:PRO:HA	2.18	0.43
1:G:108:ALA:HB2	1:G:147:ILE:HD11	1.99	0.43
1:G:194:ILE:C	1:G:197:PRO:HD2	2.44	0.43
1:A:161:ASN:OD1	1:A:161:ASN:N	2.44	0.43
1:E:137:TYR:CZ	1:E:166:KPI:HE	2.53	0.43
1:I:204:LYS:HA	1:I:204:LYS:HD3	1.89	0.43
1:B:154:LYS:HD3	2:B:301:PGE:H32	2.01	0.43
1:E:16:PRO:HD2	1:E:27:TYR:HD1	1.84	0.43
1:B:245:LYS:HA	1:B:245:LYS:HD2	1.84	0.42
1:C:231:LYS:CE	1:D:228:GLU:OE1	2.67	0.42
1:E:102:ASP:O	1:E:133:PRO:HD2	2.19	0.42
1:F:94:LYS:O	1:F:98:GLU:HG3	2.20	0.42
1:H:246:ILE:HD12	1:H:246:ILE:HA	1.87	0.42
1:K:161:ASN:OD1	1:K:161:ASN:N	2.45	0.42
1:G:139:VAL:HG13	1:G:143:THR:HG23	2.02	0.42
1:I:243:ILE:HB	1:I:293:TYR:CE2	2.54	0.42
1:E:97:LYS:HG3	1:E:132:ILE:HG21	2.01	0.42
1:F:31:ILE:HD13	1:F:66:ALA:HA	2.00	0.42
1:J:269:LEU:HD23	1:J:269:LEU:O	2.19	0.42
1:J:274:PRO:HB3	1:K:122:HIS:HB2	2.02	0.42
1:B:43:VAL:HA	1:B:78:LEU:O	2.19	0.42
1:D:37:ASN:OD1	1:D:296:LYS:HD3	2.19	0.42
1:H:64:GLU:OE1	1:H:99:HIS:CE1	2.72	0.42
1:H:232:GLU:O	1:H:236:ILE:HG12	2.20	0.42
1:I:224:PHE:HB3	1:I:233:ALA:HB2	2.01	0.42
1:J:27:TYR:CE2	1:J:62:CYS:HB3	2.54	0.42
1:A:288:GLU:O	1:A:288:GLU:HG3	2.20	0.42
1:D:246:ILE:HD12	1:D:246:ILE:HA	1.90	0.42
1:I:58:GLU:OE2	6:I:401:HOH:O	2.21	0.42
1:L:124:LYS:O	1:L:128:GLN:HG3	2.20	0.42
1:A:249:CYS:SG	1:A:282:ASN:HB3	2.59	0.42
1:C:74:LYS:HB2	1:C:74:LYS:HE2	1.63	0.42
1:I:246:ILE:HD12	1:I:246:ILE:HA	1.79	0.42
1:J:112:ASN:HA	1:K:142:ARG:NH1	2.33	0.42
1:J:194:ILE:C	1:J:197:PRO:HD2	2.45	0.42
1:C:216:ASP:OD1	1:C:217:MET:N	2.52	0.42
1:I:3:LYS:N	6:I:413:HOH:O	2.52	0.42
1:B:231:LYS:H	1:B:231:LYS:CD	2.33	0.41
1:E:131:ASP:OD1	1:E:131:ASP:O	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:246:ILE:HD12	1:K:246:ILE:HA	1.91	0.41
1:F:29:ARG:HE	1:F:29:ARG:HB2	1.58	0.41
1:G:219:SER:O	1:G:223:HIS:ND1	2.53	0.41
1:L:31:ILE:HD13	1:L:66:ALA:HA	2.02	0.41
1:F:274:PRO:HB2	1:H:114:PRO:HB3	2.01	0.41
1:I:27:TYR:CE2	1:I:62:CYS:HB3	2.55	0.41
1:J:124:LYS:O	1:J:128:GLN:HG3	2.20	0.41
1:C:35:ILE:HD11	1:C:70:CYS:SG	2.59	0.41
1:G:113:LYS:HB2	1:G:113:LYS:HE3	1.90	0.41
1:B:246:ILE:C	1:B:248:PHE:H	2.27	0.41
1:G:104:ILE:HG23	1:G:132:ILE:HD11	2.01	0.41
1:J:45:VAL:HG12	1:J:45:VAL:O	2.20	0.41
1:L:254:ILE:HA	1:L:271:PHE:CE2	2.56	0.41
1:C:166:KPI:HE	1:C:207:ILE:HB	2.03	0.41
1:K:172:ILE:HG23	1:L:194:ILE:HG21	2.01	0.41
1:C:67:VAL:O	1:C:71:LYS:HG2	2.21	0.41
1:H:23:ASP:OD1	1:H:25:GLN:HG3	2.21	0.41
1:I:32:LYS:O	1:I:36:GLU:HB2	2.20	0.41
1:J:142:ARG:O	1:K:111:TYR:HA	2.21	0.41
1:L:45:VAL:HG11	1:L:59:HIS:CE1	2.55	0.41
1:L:213:LEU:HD11	1:L:295:ILE:HD13	2.01	0.41
1:E:256:ILE:O	1:E:260:MET:HG2	2.19	0.41
1:F:35:ILE:HG12	1:F:35:ILE:H	1.71	0.41
1:I:102:ASP:O	1:I:133:PRO:HD2	2.21	0.41
1:I:142:ARG:NH2	1:I:252:ASN:O	2.53	0.41
1:G:246:ILE:C	1:G:248:PHE:H	2.28	0.41
1:H:102:ASP:O	1:H:133:PRO:HD2	2.20	0.41
1:H:135:LEU:HD13	1:H:163:TYR:CZ	2.56	0.41
1:K:43:VAL:HA	1:K:78:LEU:O	2.21	0.41
1:D:216:ASP:OD1	1:D:216:ASP:N	2.54	0.41
1:J:116:GLN:HB3	2:J:301:PGE:H42	2.02	0.41
1:L:194:ILE:C	1:L:197:PRO:HD2	2.46	0.41
1:A:274:PRO:HB2	1:G:114:PRO:HB3	2.04	0.40
1:B:10:MET:HA	1:B:41:ALA:O	2.20	0.40
1:K:182:GLU:HB3	1:K:185:MET:HE3	2.03	0.40
1:A:21:LYS:HD3	1:A:21:LYS:HA	1.92	0.40
1:A:102:ASP:O	1:A:133:PRO:HD2	2.21	0.40
1:B:194:ILE:C	1:B:197:PRO:HD2	2.47	0.40
1:B:256:ILE:O	1:B:260:MET:HG2	2.21	0.40
1:D:280:LYS:HD3	1:D:280:LYS:HA	1.85	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	293/310 (94%)	284 (97%)	9 (3%)	0	100	100
1	B	293/310 (94%)	284 (97%)	9 (3%)	0	100	100
1	C	293/310 (94%)	284 (97%)	9 (3%)	0	100	100
1	D	293/310 (94%)	285 (97%)	8 (3%)	0	100	100
1	E	291/310 (94%)	283 (97%)	8 (3%)	0	100	100
1	F	293/310 (94%)	285 (97%)	8 (3%)	0	100	100
1	G	293/310 (94%)	283 (97%)	10 (3%)	0	100	100
1	H	293/310 (94%)	284 (97%)	9 (3%)	0	100	100
1	I	293/310 (94%)	285 (97%)	8 (3%)	0	100	100
1	J	293/310 (94%)	285 (97%)	8 (3%)	0	100	100
1	K	293/310 (94%)	283 (97%)	10 (3%)	0	100	100
1	L	293/310 (94%)	284 (97%)	9 (3%)	0	100	100
All	All	3514/3720 (94%)	3409 (97%)	105 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	247/260 (95%)	247 (100%)	0	100	100
1	B	246/260 (95%)	244 (99%)	2 (1%)	73	85

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	244/260 (94%)	242 (99%)	2 (1%)	73	85
1	D	242/260 (93%)	240 (99%)	2 (1%)	73	85
1	E	246/260 (95%)	244 (99%)	2 (1%)	73	85
1	F	246/260 (95%)	245 (100%)	1 (0%)	84	92
1	G	246/260 (95%)	243 (99%)	3 (1%)	63	79
1	H	248/260 (95%)	244 (98%)	4 (2%)	55	74
1	I	247/260 (95%)	245 (99%)	2 (1%)	73	85
1	J	248/260 (95%)	243 (98%)	5 (2%)	48	68
1	K	245/260 (94%)	238 (97%)	7 (3%)	37	57
1	L	246/260 (95%)	244 (99%)	2 (1%)	73	85
All	All	2951/3120 (95%)	2919 (99%)	32 (1%)	65	81

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	29	ARG
1	B	292	LYS
1	C	98	GLU
1	C	235	LYS
1	D	25	GLN
1	D	229	ASN
1	E	74	LYS
1	E	161	ASN
1	F	29	ARG
1	G	32	LYS
1	G	239	GLU
1	G	294	LYS
1	H	29	ARG
1	H	57	GLU
1	H	71	LYS
1	H	98	GLU
1	I	26	SER
1	I	291	LYS
1	J	25	GLN
1	J	59	HIS
1	J	97	LYS
1	J	98	GLU
1	J	262	LEU

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Mol	Chain	Res	Type
1	K	59	HIS
1	K	60	ARG
1	K	68	GLU
1	K	69	THR
1	K	98	GLU
1	K	131	ASP
1	K	235	LYS
1	L	18	LYS
1	L	60	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	128	GLN
1	B	242	ASN
1	C	4	ASN
1	C	117	GLN
1	C	252	ASN
1	D	25	GLN
1	D	56	HIS
1	F	99	HIS
1	G	56	HIS
1	H	19	ASN
1	H	99	HIS
1	J	19	ASN
1	J	59	HIS
1	J	201	ASN
1	K	59	HIS
1	L	4	ASN
1	L	99	HIS
1	L	252	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

12 non-standard protein/DNA/RNA residues are modelled in this entry.

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$
1	KPI	F	166	1	11,13,14	1.37	2 (18%)	9,15,17	2.58	5 (55%)
1	KPI	L	166	1	11,13,14	1.91	2 (18%)	9,15,17	1.69	2 (22%)
1	KPI	I	166	1	11,13,14	1.93	2 (18%)	9,15,17	1.81	3 (33%)
1	KPI	G	166	1	11,13,14	2.32	3 (27%)	9,15,17	4.65	5 (55%)
1	KPI	J	166	1	11,13,14	1.96	2 (18%)	9,15,17	3.76	5 (55%)
1	KPI	K	166	1	11,13,14	1.59	2 (18%)	9,15,17	1.66	3 (33%)
1	KPI	C	166	1	11,13,14	2.31	2 (18%)	9,15,17	2.29	4 (44%)
1	KPI	D	166	1	11,13,14	1.55	2 (18%)	9,15,17	2.52	5 (55%)
1	KPI	H	166	1	11,13,14	2.33	3 (27%)	9,15,17	4.51	5 (55%)
1	KPI	E	166	1	11,13,14	1.94	2 (18%)	9,15,17	3.03	3 (33%)
1	KPI	A	166	1	11,13,14	2.25	2 (18%)	9,15,17	2.13	5 (55%)
1	KPI	B	166	1	11,13,14	2.22	4 (36%)	9,15,17	2.13	3 (33%)

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
1	KPI	F	166	1	-	0/13/14/16	-
1	KPI	L	166	1	-	0/13/14/16	-
1	KPI	I	166	1	-	0/13/14/16	-
1	KPI	G	166	1	-	6/13/14/16	-
1	KPI	J	166	1	-	7/13/14/16	-
1	KPI	K	166	1	-	3/13/14/16	-
1	KPI	C	166	1	-	0/13/14/16	-
1	KPI	D	166	1	-	0/13/14/16	-
1	KPI	H	166	1	-	6/13/14/16	-
1	KPI	E	166	1	-	7/13/14/16	-
1	KPI	A	166	1	-	4/13/14/16	-
1	KPI	B	166	1	-	0/13/14/16	-

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	166	KPI	CX2-CX1	-6.41	1.42	1.49
1	A	166	KPI	CX2-CX1	-5.89	1.43	1.49
1	J	166	KPI	O2-CX2	5.60	1.36	1.22
1	G	166	KPI	O2-CX2	5.58	1.36	1.22
1	H	166	KPI	O2-CX2	5.56	1.36	1.22
1	B	166	KPI	CX2-CX1	-5.54	1.43	1.49
1	L	166	KPI	CX2-CX1	-4.87	1.44	1.49
1	E	166	KPI	CX2-CX1	-4.77	1.44	1.49
1	I	166	KPI	CX2-CX1	-4.61	1.44	1.49
1	H	166	KPI	O-C	4.38	1.36	1.20
1	G	166	KPI	O-C	4.31	1.36	1.20
1	A	166	KPI	O1-CX2	-3.96	1.19	1.30
1	I	166	KPI	O1-CX2	-3.78	1.20	1.30
1	D	166	KPI	O1-CX2	-3.74	1.20	1.30
1	K	166	KPI	CX2-CX1	-3.72	1.45	1.49
1	C	166	KPI	O1-CX2	-3.50	1.21	1.30
1	E	166	KPI	O1-CX2	-3.42	1.21	1.30
1	L	166	KPI	O1-CX2	-3.27	1.21	1.30
1	B	166	KPI	O1-CX2	-3.25	1.21	1.30
1	K	166	KPI	O1-CX2	-3.22	1.21	1.30
1	F	166	KPI	O1-CX2	-2.95	1.22	1.30
1	D	166	KPI	CX2-CX1	-2.79	1.46	1.49
1	F	166	KPI	CX2-CX1	-2.62	1.46	1.49
1	H	166	KPI	O1-CX2	-2.59	1.23	1.30
1	J	166	KPI	O1-CX2	-2.55	1.23	1.30
1	G	166	KPI	O1-CX2	-2.54	1.23	1.30
1	B	166	KPI	CB-CA	-2.38	1.49	1.53
1	B	166	KPI	CA-N	-2.20	1.41	1.48

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	166	KPI	C1-CX1-CX2	-10.78	107.95	118.11
1	H	166	KPI	C1-CX1-CX2	-9.60	109.06	118.11
1	J	166	KPI	C1-CX1-CX2	-7.19	111.33	118.11
1	E	166	KPI	CX2-CX1-NZ	5.87	128.70	114.88
1	H	166	KPI	O1-CX2-CX1	5.60	128.42	116.50
1	H	166	KPI	O2-CX2-CX1	-5.59	114.52	121.35
1	J	166	KPI	O2-CX2-CX1	-5.47	114.67	121.35
1	E	166	KPI	CE-NZ-CX1	5.43	136.49	121.58
1	J	166	KPI	O1-CX2-CX1	5.16	127.50	116.50
1	G	166	KPI	O1-CX2-CX1	4.98	127.10	116.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	166	KPI	C1-CX1-CX2	4.90	122.74	118.11
1	G	166	KPI	O2-CX2-CX1	-4.87	115.40	121.35
1	G	166	KPI	CD-CE-NZ	4.50	118.63	110.67
1	D	166	KPI	CE-NZ-CX1	4.25	133.26	121.58
1	H	166	KPI	CD-CE-NZ	3.98	117.73	110.67
1	C	166	KPI	CE-NZ-CX1	3.65	131.61	121.58
1	C	166	KPI	CD-CE-NZ	-3.64	104.23	110.67
1	B	166	KPI	O2-CX2-CX1	-3.51	117.07	121.35
1	A	166	KPI	CE-NZ-CX1	3.47	131.12	121.58
1	D	166	KPI	CX2-CX1-NZ	3.40	122.90	114.88
1	C	166	KPI	O2-CX2-CX1	-3.18	117.46	121.35
1	F	166	KPI	CE-NZ-CX1	3.12	130.15	121.58
1	B	166	KPI	CX2-CX1-NZ	3.12	122.23	114.88
1	I	166	KPI	C1-CX1-CX2	3.12	121.05	118.11
1	D	166	KPI	O1-CX2-O2	-3.08	116.57	123.90
1	F	166	KPI	O1-CX2-CX1	3.06	123.01	116.50
1	D	166	KPI	O2-CX2-CX1	3.00	125.02	121.35
1	K	166	KPI	CE-NZ-CX1	2.98	129.76	121.58
1	H	166	KPI	O1-CX2-O2	-2.87	117.06	123.90
1	L	166	KPI	CX2-CX1-NZ	2.84	121.57	114.88
1	E	166	KPI	O1-CX2-O2	-2.79	117.26	123.90
1	L	166	KPI	O2-CX2-CX1	-2.76	117.98	121.35
1	K	166	KPI	CX2-CX1-NZ	2.75	121.34	114.88
1	F	166	KPI	O1-CX2-O2	-2.73	117.40	123.90
1	A	166	KPI	O1-CX2-O2	-2.72	117.42	123.90
1	G	166	KPI	O1-CX2-O2	-2.69	117.50	123.90
1	I	166	KPI	CE-NZ-CX1	2.68	128.95	121.58
1	A	166	KPI	O1-CX2-CX1	2.68	122.21	116.50
1	B	166	KPI	O1-CX2-CX1	2.57	121.97	116.50
1	J	166	KPI	O1-CX2-O2	-2.55	117.83	123.90
1	D	166	KPI	CD-CE-NZ	-2.54	106.18	110.67
1	A	166	KPI	CD-CE-NZ	-2.53	106.20	110.67
1	J	166	KPI	CD-CE-NZ	2.47	115.05	110.67
1	C	166	KPI	O1-CX2-CX1	2.40	121.62	116.50
1	I	166	KPI	O1-CX2-O2	-2.32	118.38	123.90
1	K	166	KPI	O1-CX2-O2	-2.13	118.83	123.90
1	F	166	KPI	CD-CE-NZ	-2.09	106.97	110.67
1	A	166	KPI	CX2-CX1-NZ	2.09	119.81	114.88

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	166	KPI	NZ-CX1-CX2-O1
1	A	166	KPI	NZ-CX1-CX2-O2
1	A	166	KPI	C1-CX1-CX2-O1
1	A	166	KPI	C1-CX1-CX2-O2
1	E	166	KPI	N-CA-CB-CG
1	E	166	KPI	C-CA-CB-CG
1	G	166	KPI	C1-CX1-NZ-CE
1	G	166	KPI	CX2-CX1-NZ-CE
1	G	166	KPI	NZ-CX1-CX2-O1
1	G	166	KPI	NZ-CX1-CX2-O2
1	H	166	KPI	C1-CX1-NZ-CE
1	H	166	KPI	CX2-CX1-NZ-CE
1	H	166	KPI	NZ-CX1-CX2-O1
1	H	166	KPI	NZ-CX1-CX2-O2
1	J	166	KPI	C1-CX1-NZ-CE
1	J	166	KPI	CX2-CX1-NZ-CE
1	J	166	KPI	NZ-CX1-CX2-O1
1	J	166	KPI	NZ-CX1-CX2-O2
1	K	166	KPI	NZ-CX1-CX2-O1
1	K	166	KPI	NZ-CX1-CX2-O2
1	E	166	KPI	CA-CB-CG-CD
1	E	166	KPI	C1-CX1-NZ-CE
1	E	166	KPI	CE-CD-CG-CB
1	E	166	KPI	CX2-CX1-NZ-CE
1	G	166	KPI	C-CA-CB-CG
1	H	166	KPI	C-CA-CB-CG
1	E	166	KPI	CG-CD-CE-NZ
1	J	166	KPI	C1-CX1-CX2-O1
1	K	166	KPI	C1-CX1-CX2-O1
1	J	166	KPI	CE-CD-CG-CB
1	H	166	KPI	CE-CD-CG-CB
1	G	166	KPI	CE-CD-CG-CB
1	J	166	KPI	C1-CX1-CX2-O2

There are no ring outliers.

5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	G	166	KPI	3	0
1	J	166	KPI	1	0
1	C	166	KPI	1	0
1	H	166	KPI	3	0
1	E	166	KPI	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 7 are monoatomic - leaving 18 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	$\# Z > 2$
2	PGE	J	301	-	9,9,9	0.31	0	8,8,8	0.34	0
2	PGE	B	301	-	9,9,9	0.32	0	8,8,8	0.26	0
4	ACT	F	303	-	3,3,3	1.42	1 (33%)	3,3,3	1.37	0
2	PGE	C	301	-	9,9,9	0.32	0	8,8,8	0.24	0
2	PGE	L	301	-	9,9,9	0.33	0	8,8,8	0.26	0
4	ACT	C	305	-	3,3,3	1.23	0	3,3,3	1.50	0
4	ACT	L	302	-	3,3,3	1.04	0	3,3,3	1.50	1 (33%)
2	PGE	H	301	-	9,9,9	0.31	0	8,8,8	0.27	0
5	EDO	J	303	-	3,3,3	0.45	0	2,2,2	0.29	0
2	PGE	E	302	-	9,9,9	0.38	0	8,8,8	0.28	0
2	PGE	E	301	-	9,9,9	0.32	0	8,8,8	0.27	0
2	PGE	D	301	-	9,9,9	0.33	0	8,8,8	0.25	0
2	PGE	F	301	-	9,9,9	0.31	0	8,8,8	0.25	0
2	PGE	I	301	-	9,9,9	0.31	0	8,8,8	0.27	0
2	PGE	K	301	-	9,9,9	0.32	0	8,8,8	0.30	0
2	PGE	C	302	-	9,9,9	0.31	0	8,8,8	0.29	0
2	PGE	G	301	-	9,9,9	0.33	0	8,8,8	0.26	0
2	PGE	A	301	-	9,9,9	0.32	0	8,8,8	0.29	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
2	PGE	J	301	-	-	5/7/7/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PGE	B	301	-	-	2/7/7/7	-
2	PGE	C	301	-	-	2/7/7/7	-
2	PGE	L	301	-	-	2/7/7/7	-
5	EDO	J	303	-	-	1/1/1/1	-
2	PGE	H	301	-	-	4/7/7/7	-
2	PGE	E	302	-	-	4/7/7/7	-
2	PGE	E	301	-	-	0/7/7/7	-
2	PGE	D	301	-	-	2/7/7/7	-
2	PGE	F	301	-	-	3/7/7/7	-
2	PGE	I	301	-	-	5/7/7/7	-
2	PGE	K	301	-	-	3/7/7/7	-
2	PGE	C	302	-	-	1/7/7/7	-
2	PGE	G	301	-	-	2/7/7/7	-
2	PGE	A	301	-	-	5/7/7/7	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	303	ACT	CH3-C	2.08	1.57	1.49

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	302	ACT	OXT-C-O	2.07	129.72	122.03

There are no chirality outliers.

All (41) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	PGE	O2-C3-C4-O3
2	J	301	PGE	O2-C3-C4-O3
2	K	301	PGE	O2-C3-C4-O3
2	I	301	PGE	O1-C1-C2-O2
2	D	301	PGE	O2-C3-C4-O3
2	E	302	PGE	O2-C3-C4-O3
2	A	301	PGE	O1-C1-C2-O2
2	K	301	PGE	C3-C4-O3-C5
2	E	302	PGE	O1-C1-C2-O2
2	I	301	PGE	O2-C3-C4-O3

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Mol	Chain	Res	Type	Atoms
2	H	301	PGE	O3-C5-C6-O4
2	E	302	PGE	O3-C5-C6-O4
2	H	301	PGE	O2-C3-C4-O3
5	J	303	EDO	O1-C1-C2-O2
2	A	301	PGE	O3-C5-C6-O4
2	B	301	PGE	O1-C1-C2-O2
2	G	301	PGE	C1-C2-O2-C3
2	G	301	PGE	C6-C5-O3-C4
2	F	301	PGE	C1-C2-O2-C3
2	H	301	PGE	C3-C4-O3-C5
2	J	301	PGE	C6-C5-O3-C4
2	F	301	PGE	C6-C5-O3-C4
2	F	301	PGE	O3-C5-C6-O4
2	J	301	PGE	O1-C1-C2-O2
2	A	301	PGE	C1-C2-O2-C3
2	I	301	PGE	C6-C5-O3-C4
2	K	301	PGE	C1-C2-O2-C3
2	L	301	PGE	O3-C5-C6-O4
2	C	302	PGE	O2-C3-C4-O3
2	B	301	PGE	C6-C5-O3-C4
2	J	301	PGE	C1-C2-O2-C3
2	D	301	PGE	C6-C5-O3-C4
2	A	301	PGE	C6-C5-O3-C4
2	I	301	PGE	C4-C3-O2-C2
2	E	302	PGE	C1-C2-O2-C3
2	I	301	PGE	O3-C5-C6-O4
2	H	301	PGE	C1-C2-O2-C3
2	J	301	PGE	O3-C5-C6-O4
2	C	301	PGE	O2-C3-C4-O3
2	C	301	PGE	C1-C2-O2-C3
2	L	301	PGE	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 16 short contacts:

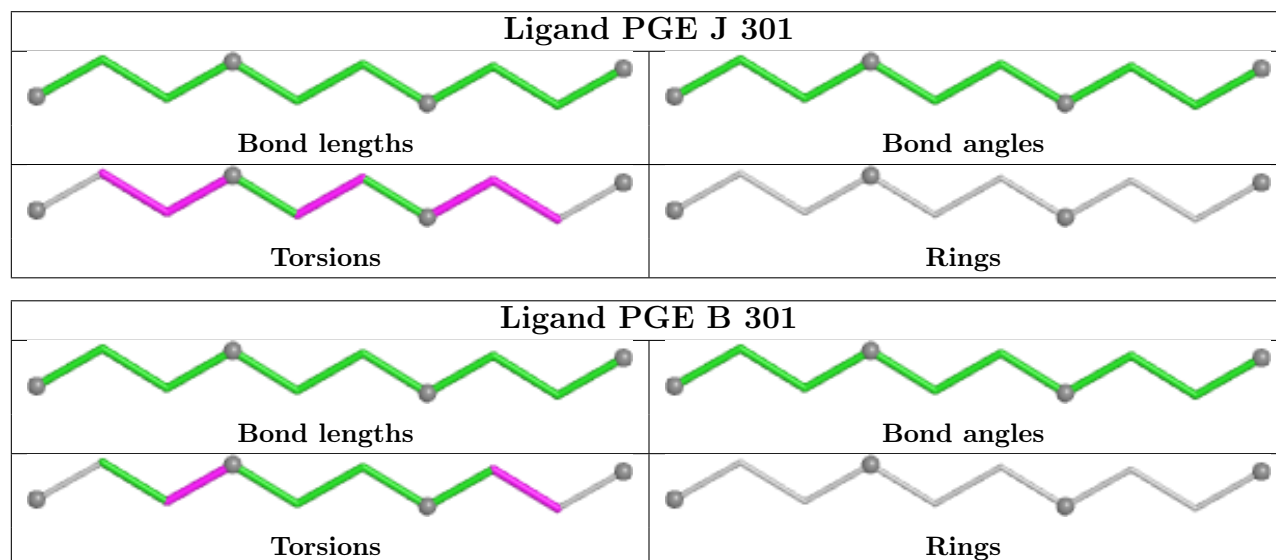
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	J	301	PGE	1	0
2	B	301	PGE	1	0
4	C	305	ACT	5	0
4	L	302	ACT	4	0
2	D	301	PGE	1	0
2	G	301	PGE	1	0

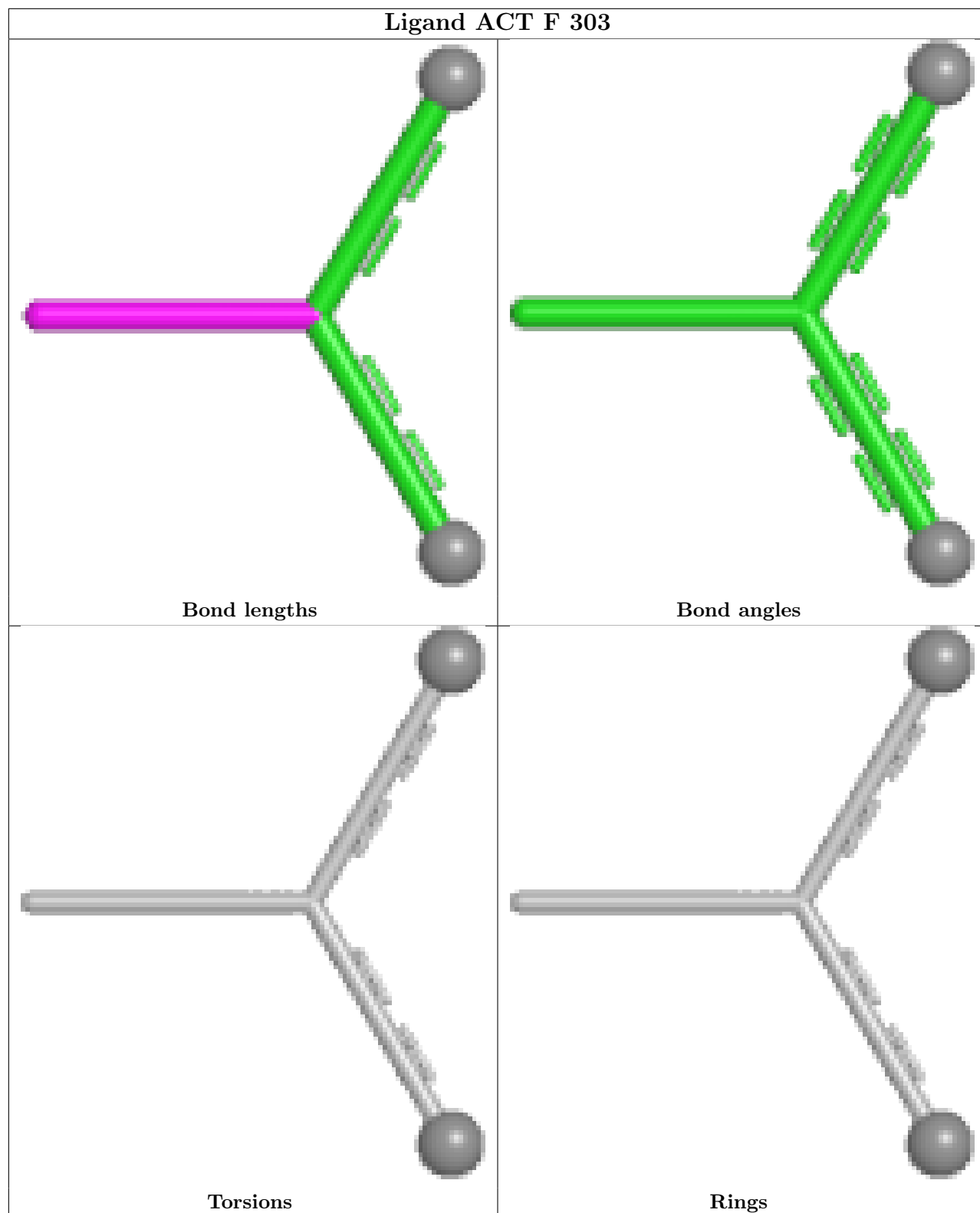
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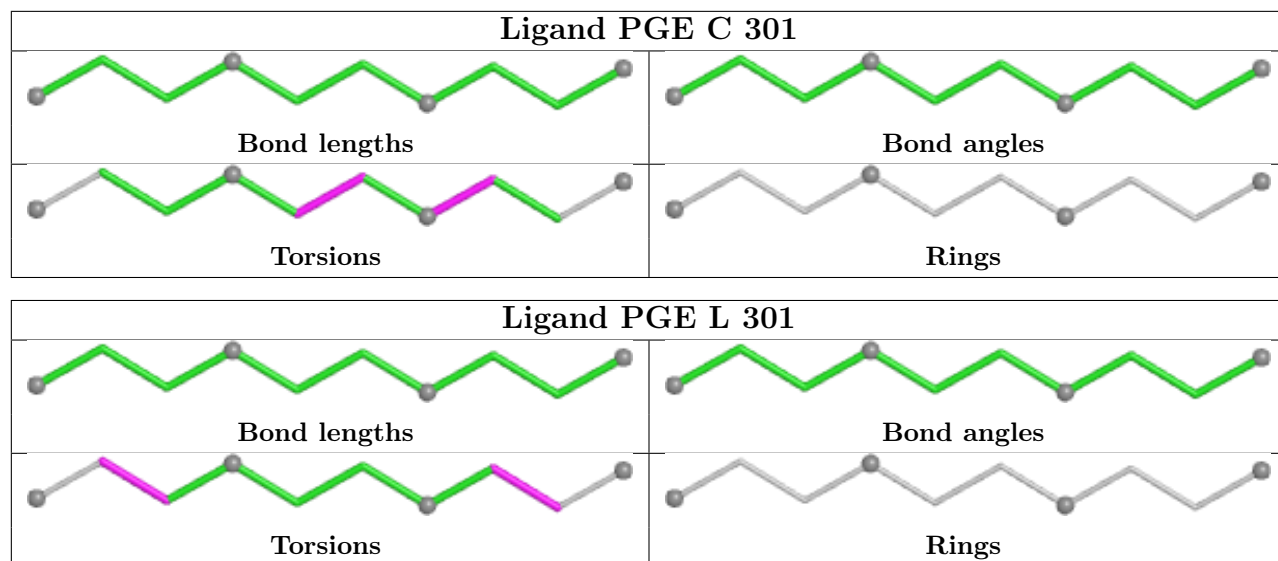
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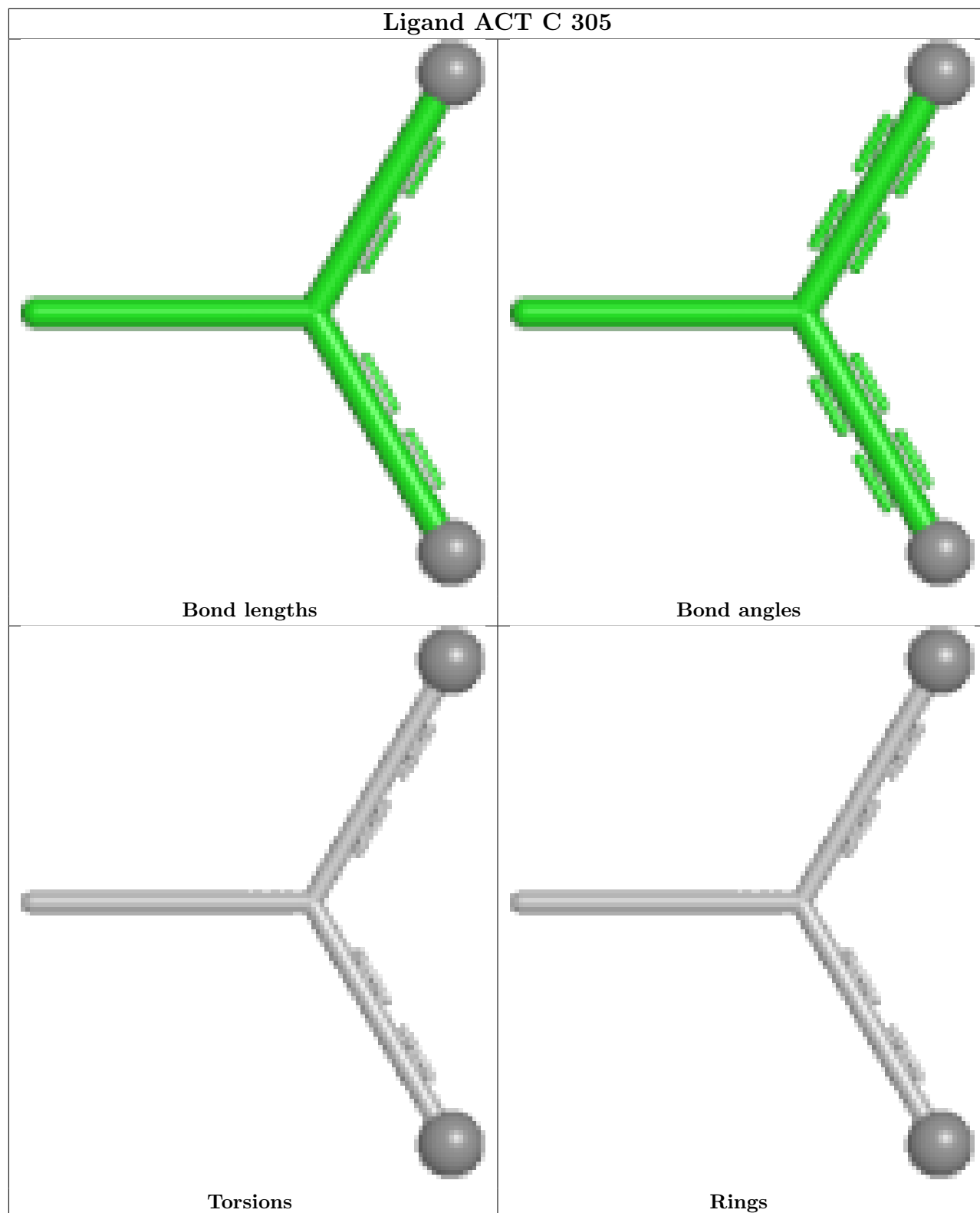
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	PGE	3	0

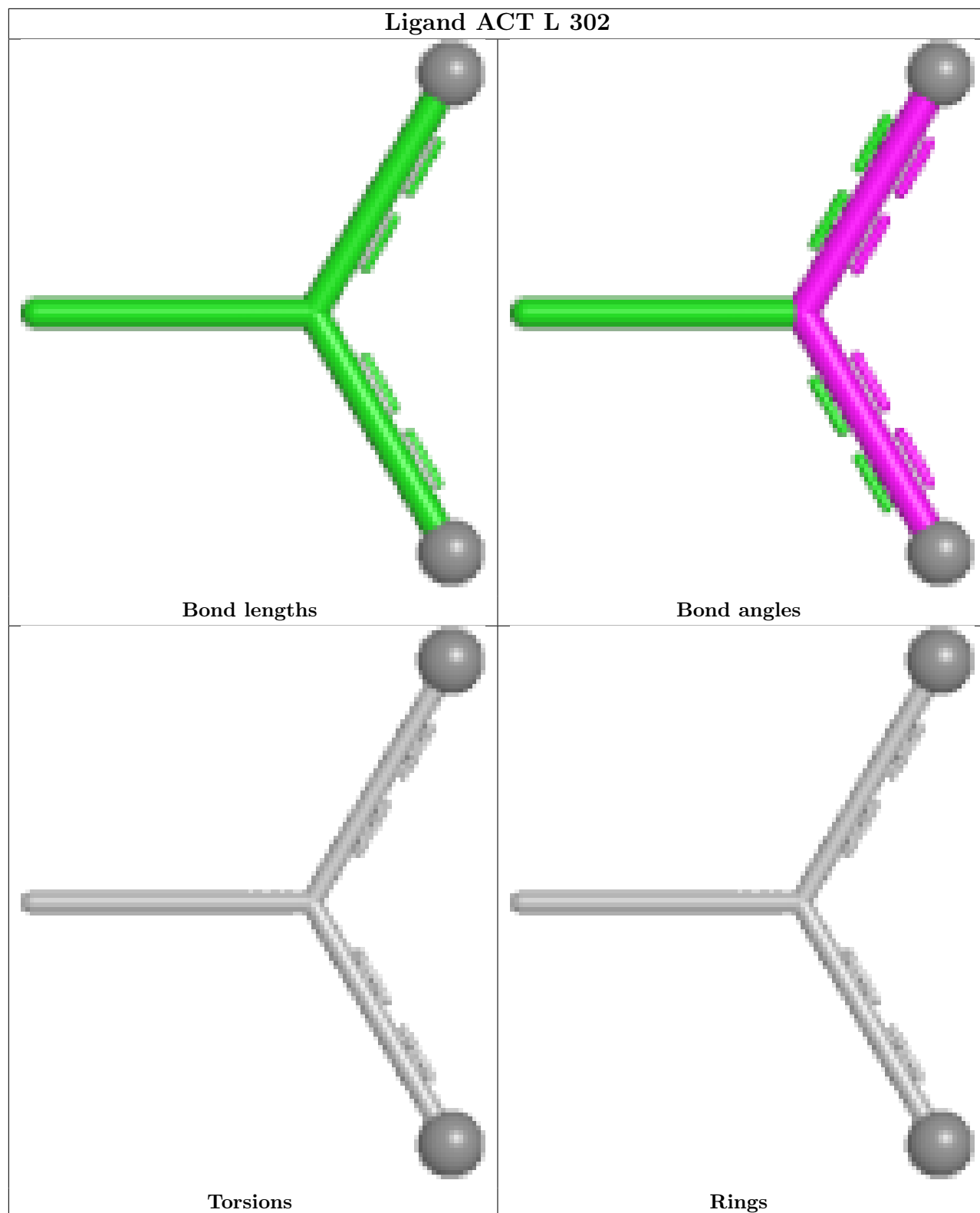
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

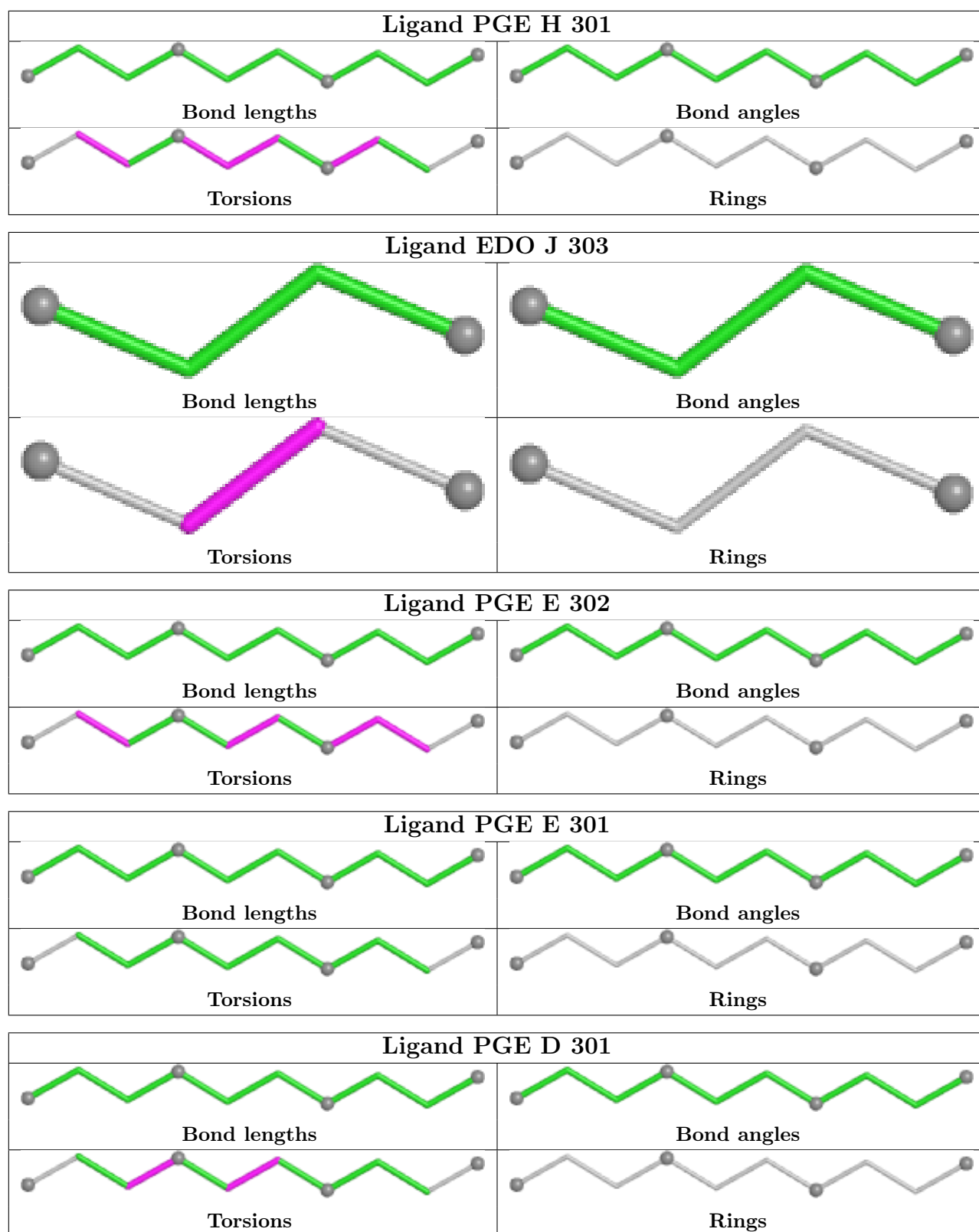


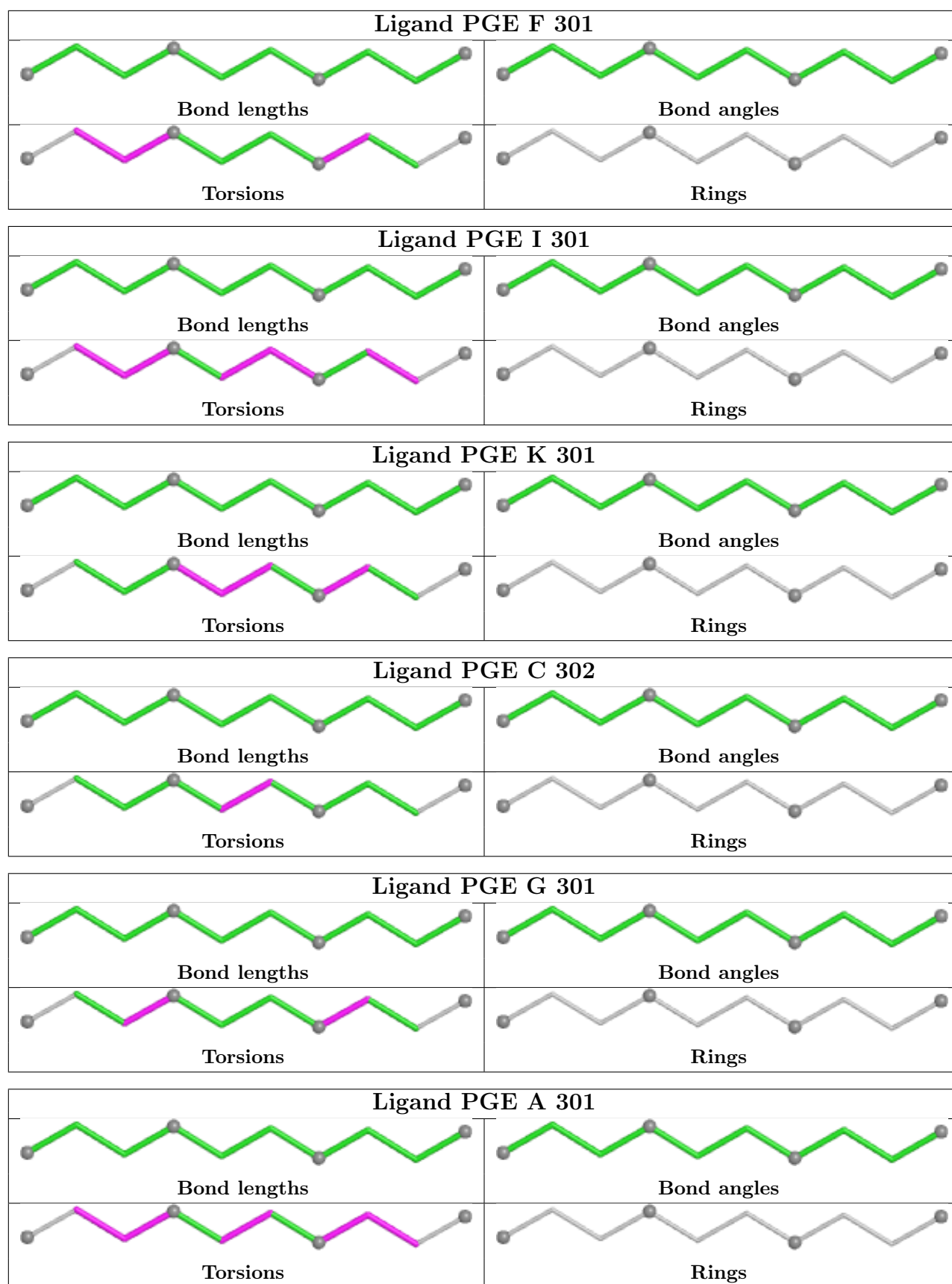












5.7 Other polymers [i](#)

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	295/310 (95%)	-1.12	0 100 100	34, 49, 72, 91	0
1	B	295/310 (95%)	-1.11	0 100 100	37, 50, 70, 98	0
1	C	295/310 (95%)	-1.11	0 100 100	34, 50, 71, 87	0
1	D	295/310 (95%)	-1.14	0 100 100	36, 49, 73, 89	0
1	E	293/310 (94%)	-1.13	1 (0%) 90 88	30, 54, 75, 97	0
1	F	295/310 (95%)	-1.12	0 100 100	36, 51, 73, 105	0
1	G	295/310 (95%)	-1.14	0 100 100	36, 51, 70, 89	0
1	H	295/310 (95%)	-1.08	0 100 100	38, 53, 73, 103	0
1	I	295/310 (95%)	-1.12	0 100 100	37, 53, 73, 90	0
1	J	295/310 (95%)	-1.11	0 100 100	40, 56, 83, 96	0
1	K	295/310 (95%)	-1.14	0 100 100	41, 55, 77, 108	0
1	L	295/310 (95%)	-1.14	0 100 100	24, 57, 81, 97	0
All	All	3538/3720 (95%)	-1.12	1 (0%) 100 100	24, 53, 75, 108	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	164	GLY	2.7

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

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
1	KPI	A	166	14/15	0.98	0.06	33,43,52,54	0
1	KPI	I	166	14/15	0.98	0.05	55,58,61,64	0
1	KPI	K	166	14/15	0.98	0.05	45,52,59,61	0
1	KPI	D	166	14/15	0.99	0.06	45,49,57,57	0
1	KPI	E	166	14/15	0.99	0.05	42,48,59,64	0
1	KPI	F	166	14/15	0.99	0.05	52,54,64,66	0
1	KPI	G	166	14/15	0.99	0.07	46,51,57,59	0
1	KPI	H	166	14/15	0.99	0.04	33,41,50,50	0
1	KPI	B	166	14/15	0.99	0.05	36,42,59,61	0
1	KPI	J	166	14/15	0.99	0.06	44,51,62,64	0
1	KPI	C	166	14/15	0.99	0.05	40,46,54,58	0
1	KPI	L	166	14/15	0.99	0.05	45,50,51,52	0

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
2	PGE	J	301	10/10	0.97	0.06	56,61,64,66	0
4	ACT	C	305	4/4	0.97	0.06	62,68,70,71	0
2	PGE	E	302	10/10	0.98	0.05	57,68,70,70	0
2	PGE	F	301	10/10	0.98	0.05	52,59,62,62	0
2	PGE	C	302	10/10	0.98	0.05	63,66,69,70	0
3	MG	D	302	1/1	0.98	0.09	59,59,59,59	0
3	MG	E	303	1/1	0.98	0.09	59,59,59,59	0
3	MG	F	302	1/1	0.98	0.10	54,54,54,54	0
3	MG	J	302	1/1	0.98	0.10	61,61,61,61	0
2	PGE	E	301	10/10	0.98	0.05	72,72,74,76	0
4	ACT	L	302	4/4	0.98	0.05	73,79,80,82	0
5	EDO	J	303	4/4	0.98	0.07	63,64,65,67	0
2	PGE	K	301	10/10	0.99	0.05	66,72,80,81	0
2	PGE	L	301	10/10	0.99	0.05	65,66,70,72	0
3	MG	A	302	1/1	0.99	0.09	55,55,55,55	0
3	MG	C	303	1/1	0.99	0.04	47,47,47,47	0
2	PGE	B	301	10/10	0.99	0.05	62,63,66,66	0

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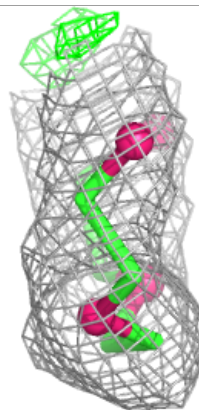
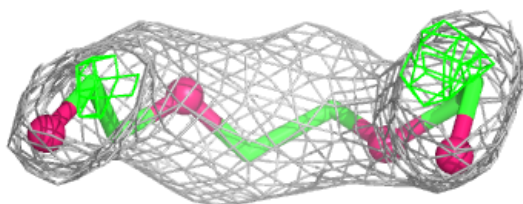
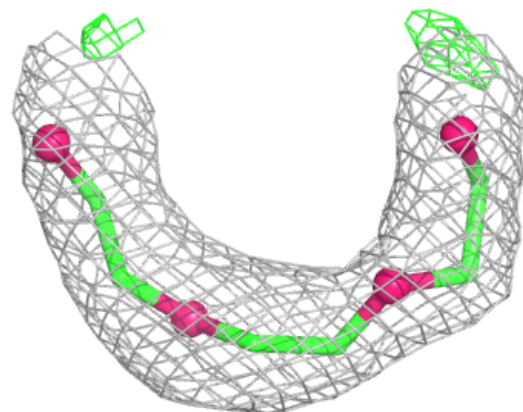
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PGE	C	301	10/10	0.99	0.04	52,58,61,63	0
2	PGE	A	301	10/10	0.99	0.04	59,62,68,68	0
2	PGE	G	301	10/10	0.99	0.05	55,58,60,62	0
2	PGE	H	301	10/10	0.99	0.04	60,61,65,67	0
4	ACT	F	303	4/4	0.99	0.04	53,54,55,56	0
2	PGE	I	301	10/10	0.99	0.04	60,66,71,71	0
2	PGE	D	301	10/10	0.99	0.04	59,61,65,67	0
3	MG	C	304	1/1	1.00	0.09	53,53,53,53	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.

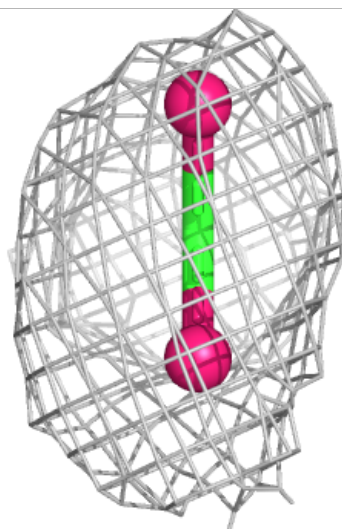
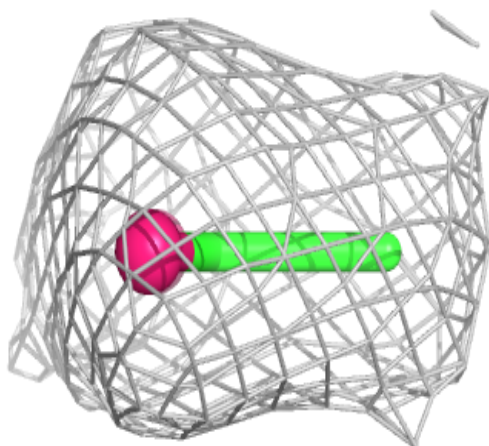
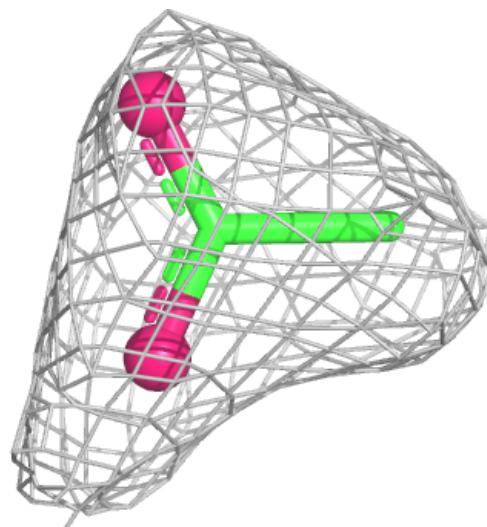
Electron density around PGE J 301:

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



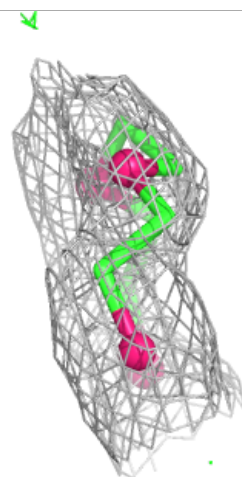
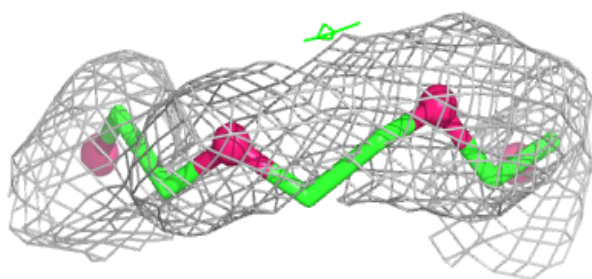
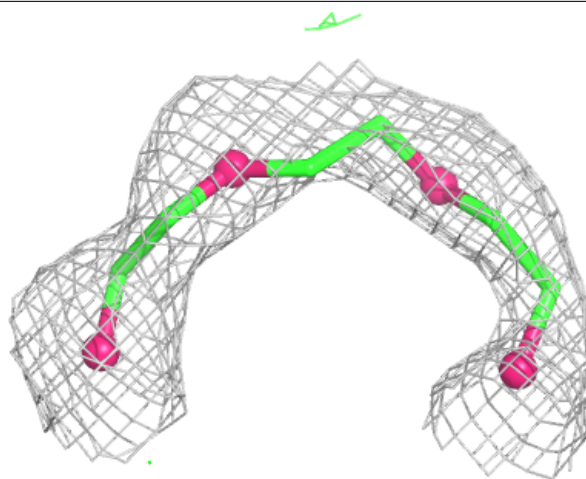
Electron density around ACT C 305:

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and green (positive)



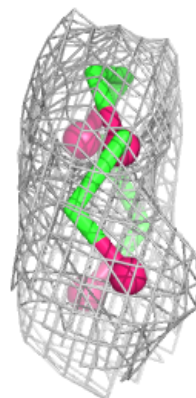
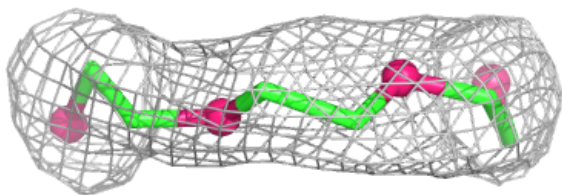
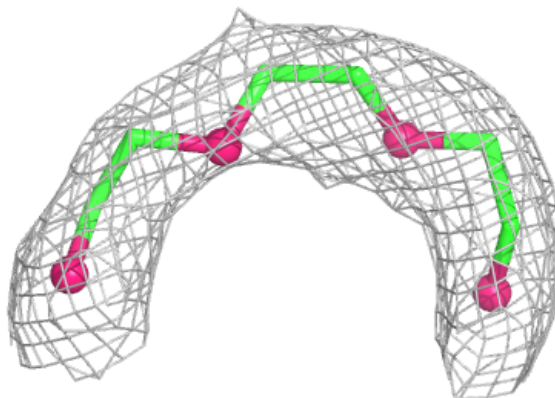
Electron density around PGE E 302:

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and green (positive)

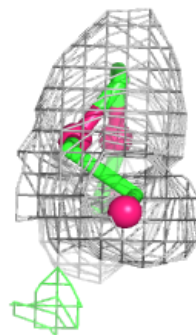
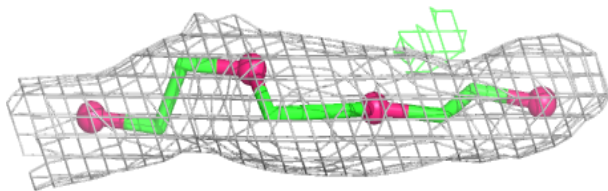
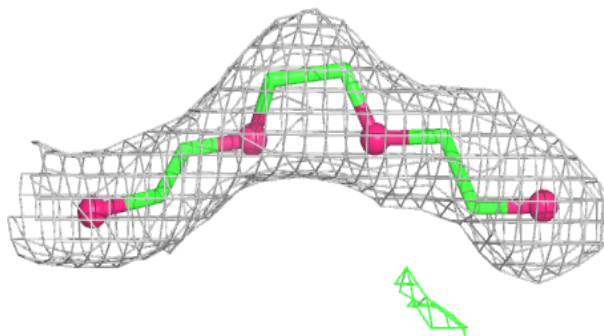


Electron density around PGE F 301:

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and green (positive)

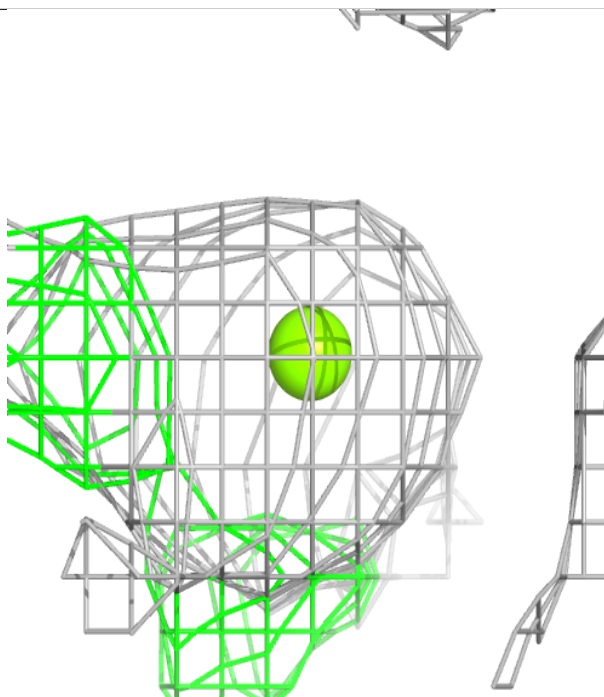
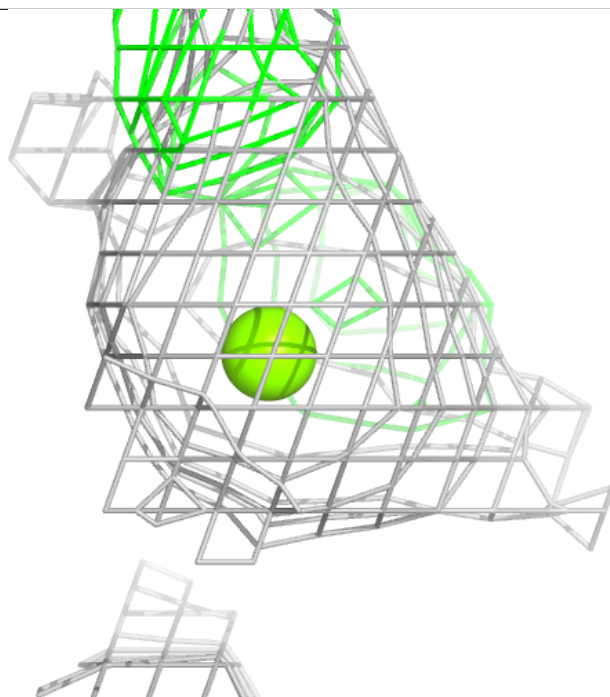
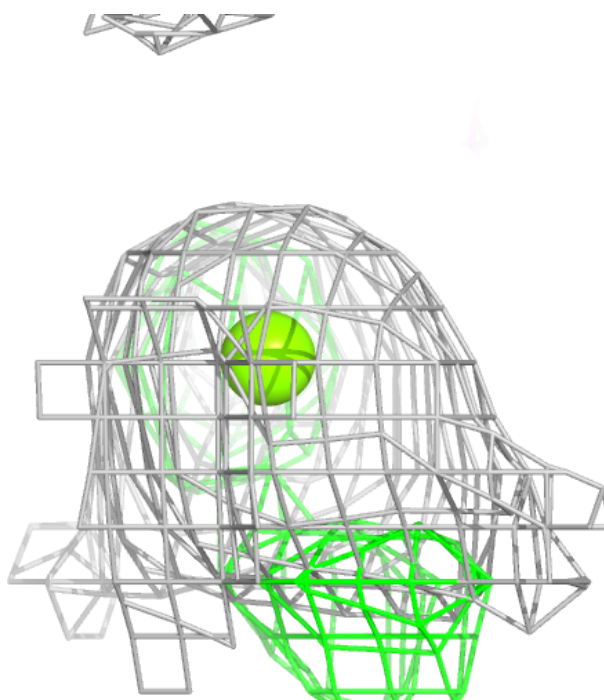
**Electron density around PGE C 302:**

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



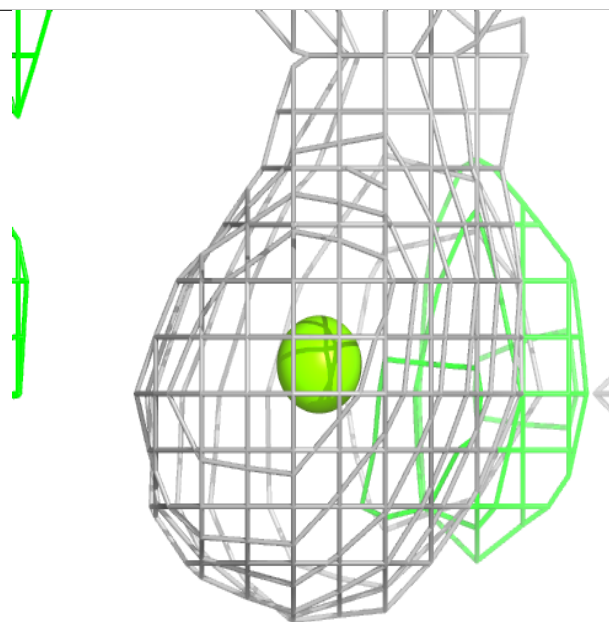
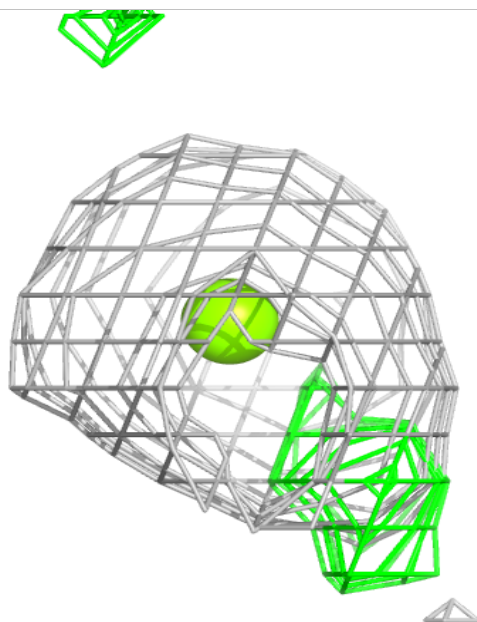
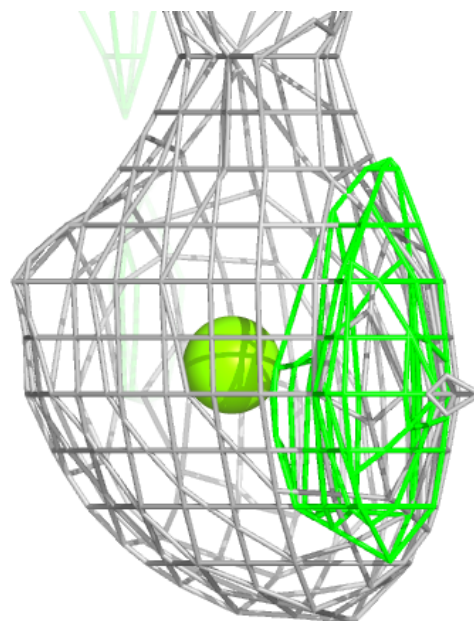
Electron density around MG D 302:

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and green (positive)



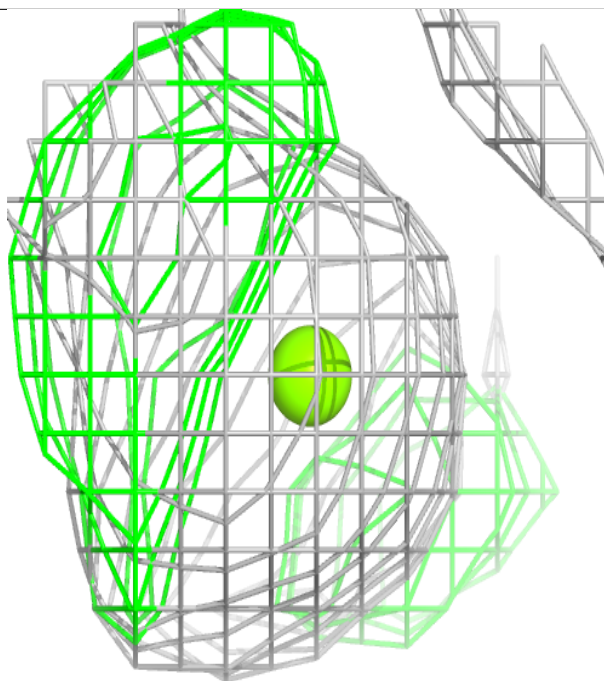
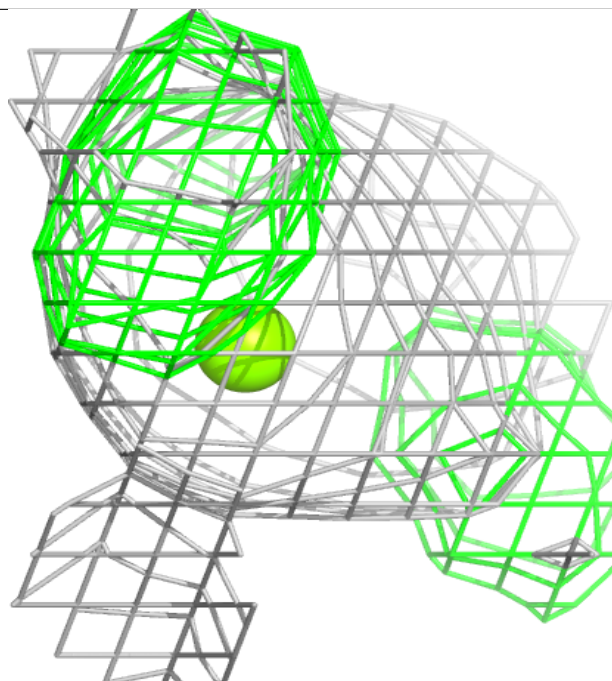
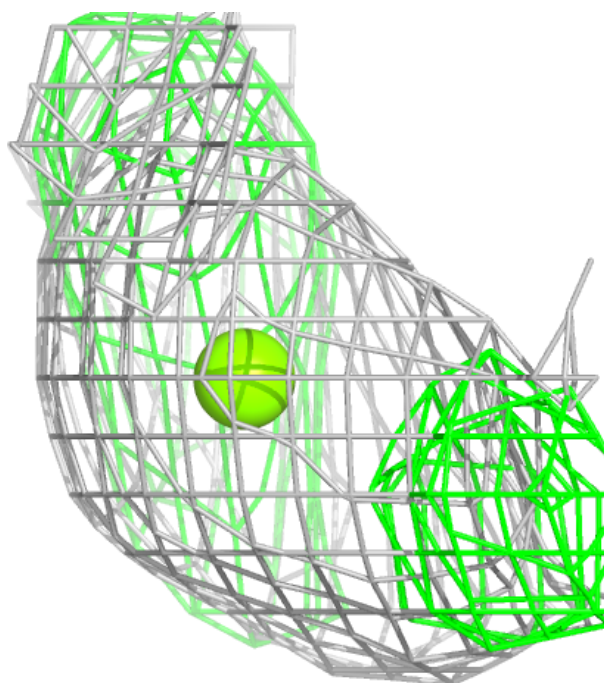
Electron density around MG E 303:

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



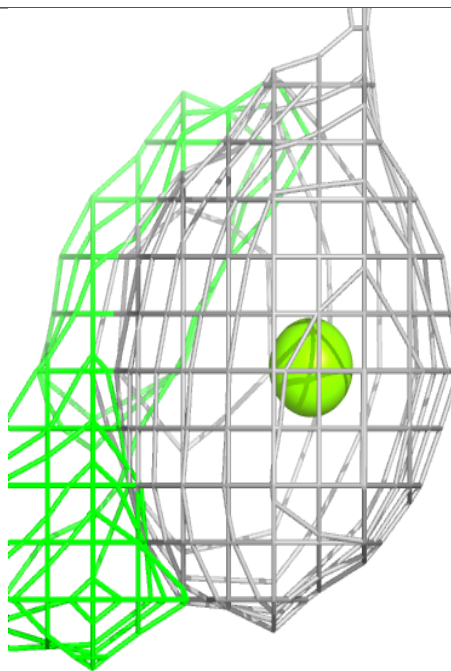
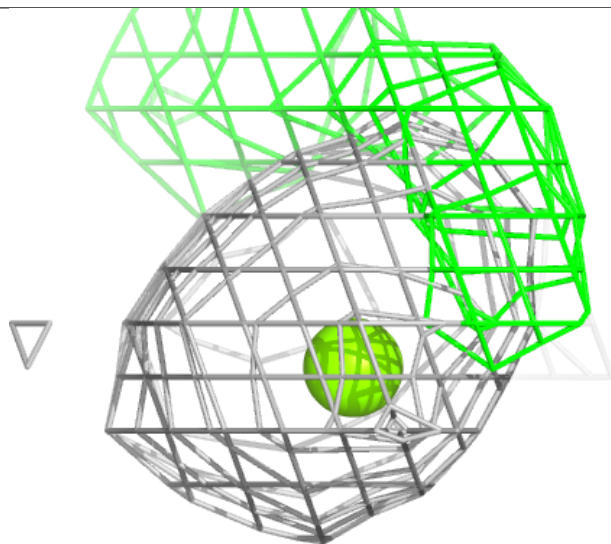
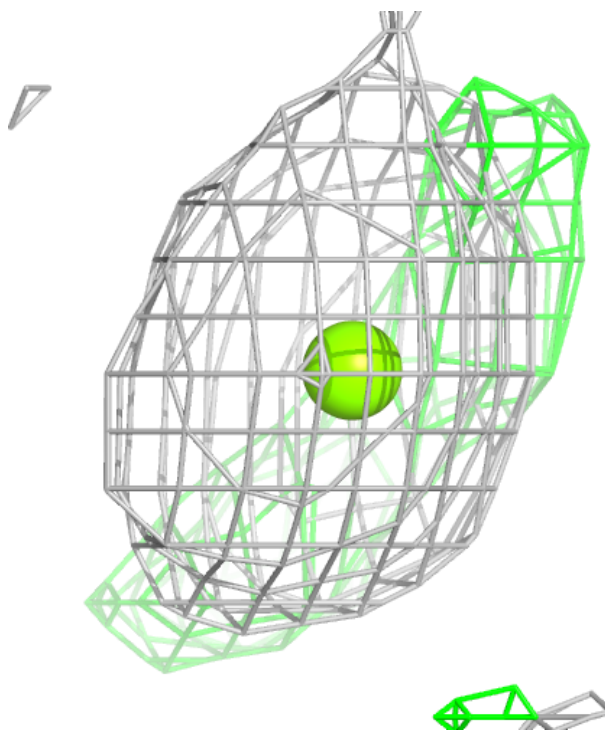
Electron density around MG F 302:

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and green (positive)



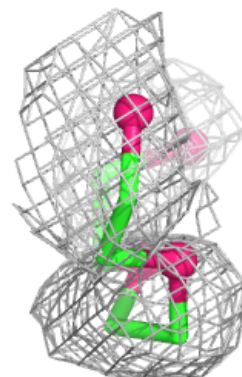
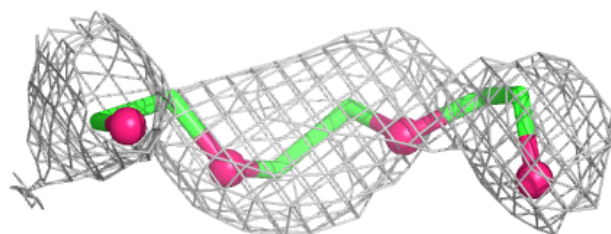
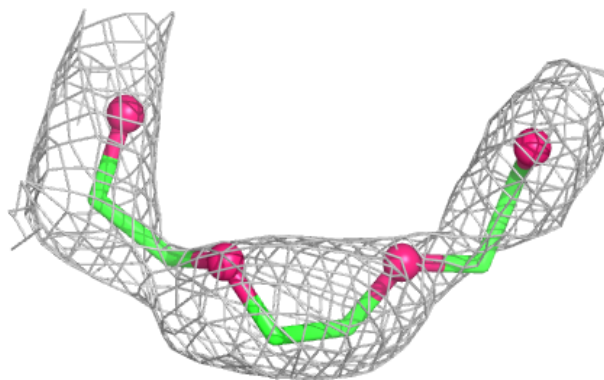
Electron density around MG J 302:

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and green (positive)



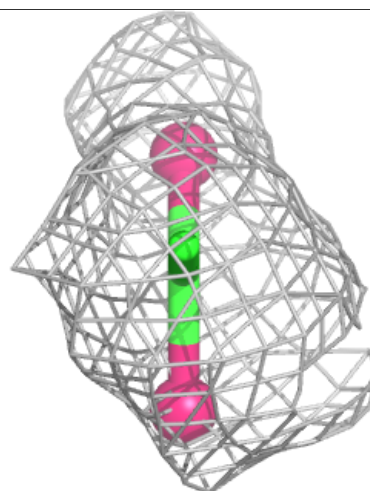
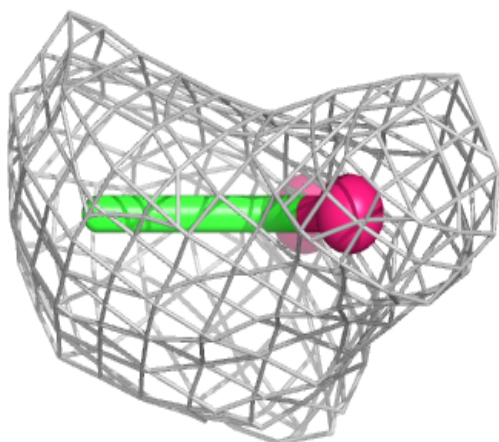
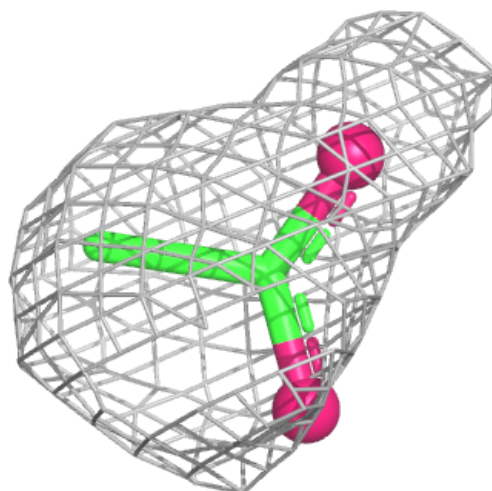
Electron density around PGE E 301:

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and green (positive)



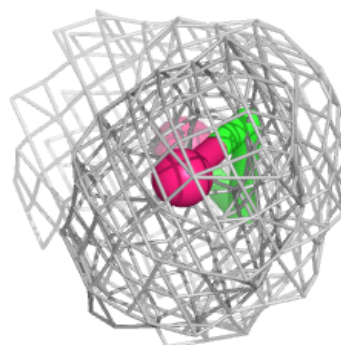
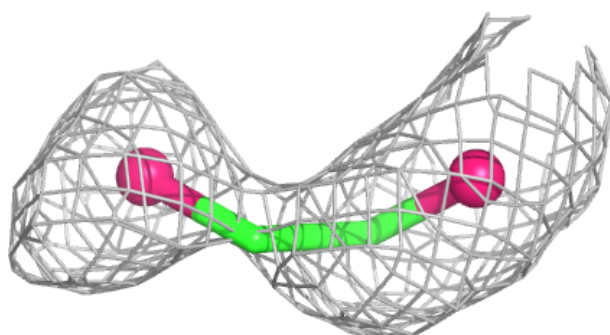
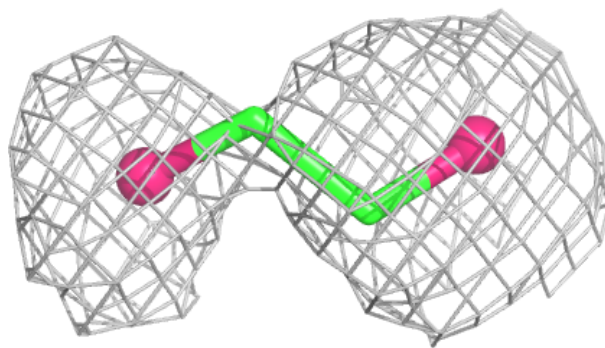
Electron density around ACT L 302:

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and green (positive)



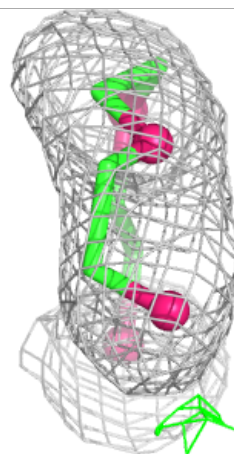
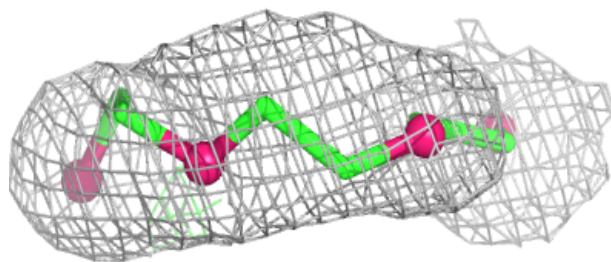
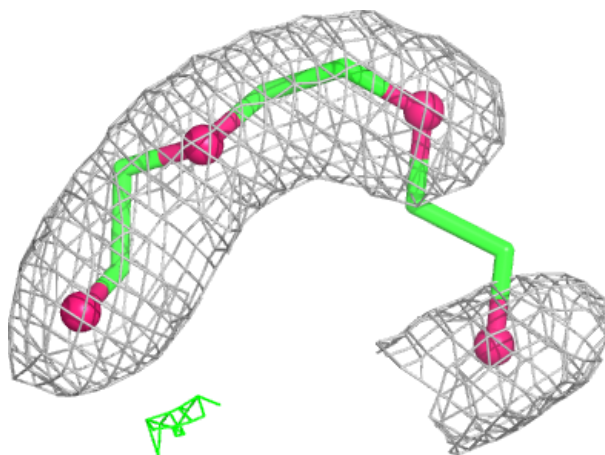
Electron density around EDO J 303:

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and green (positive)



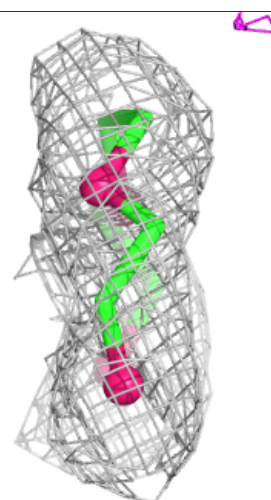
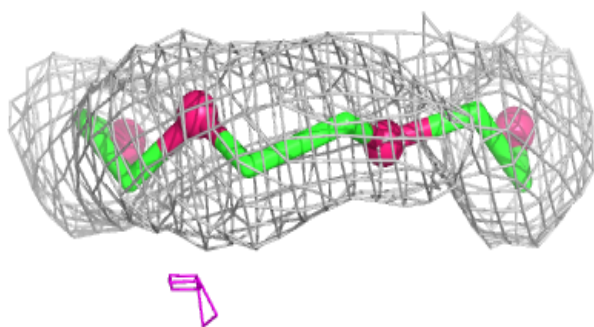
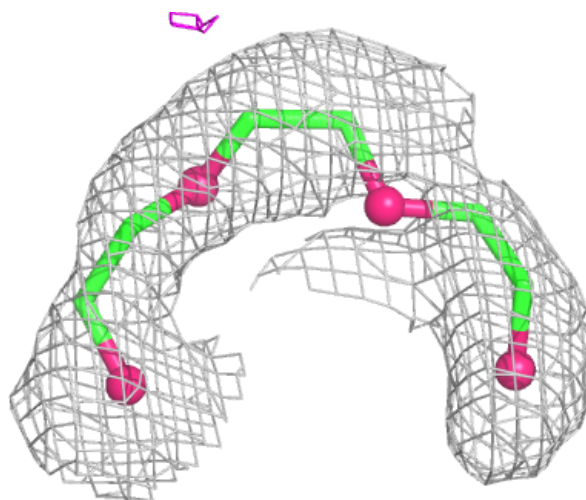
Electron density around PGE K 301:

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and green (positive)



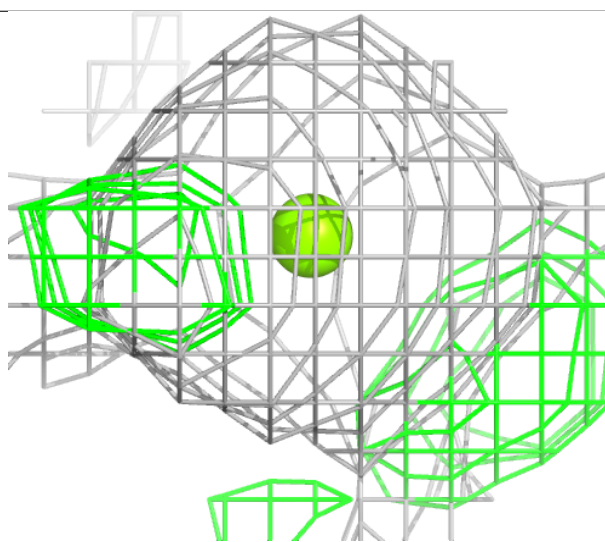
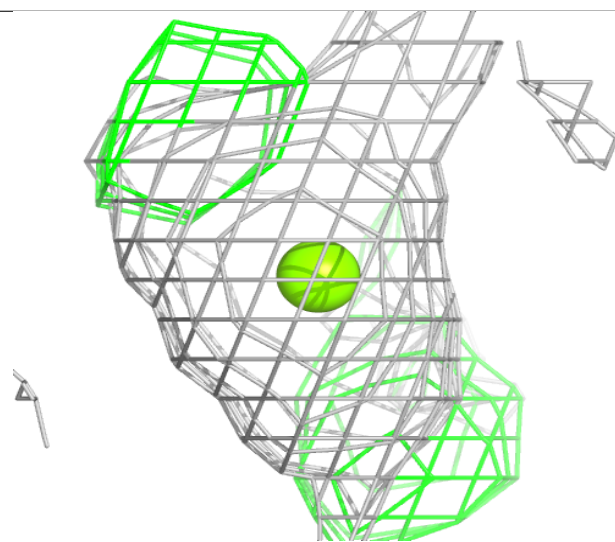
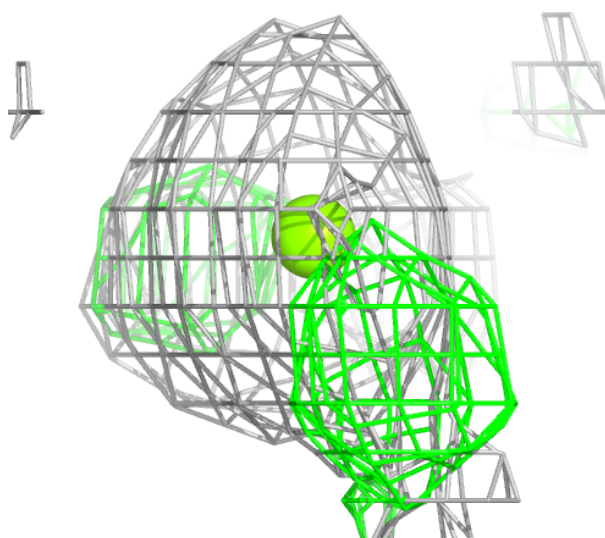
Electron density around PGE L 301:

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and green (positive)



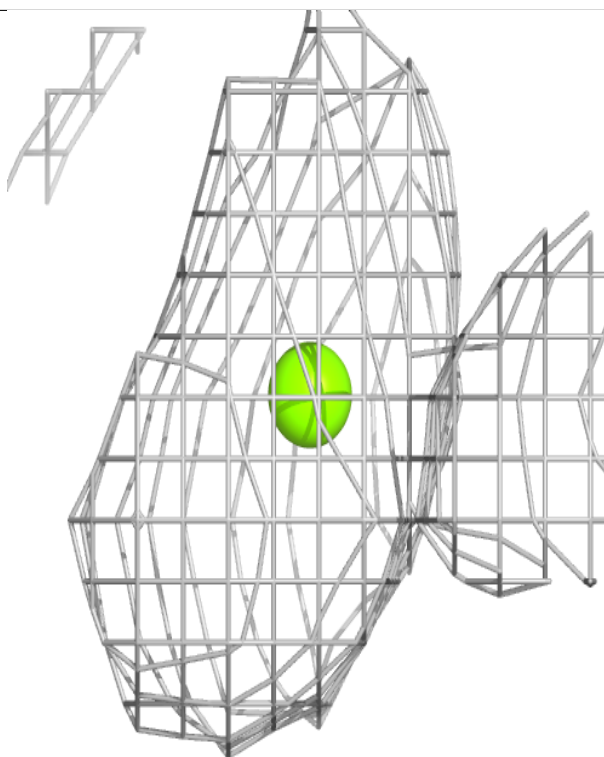
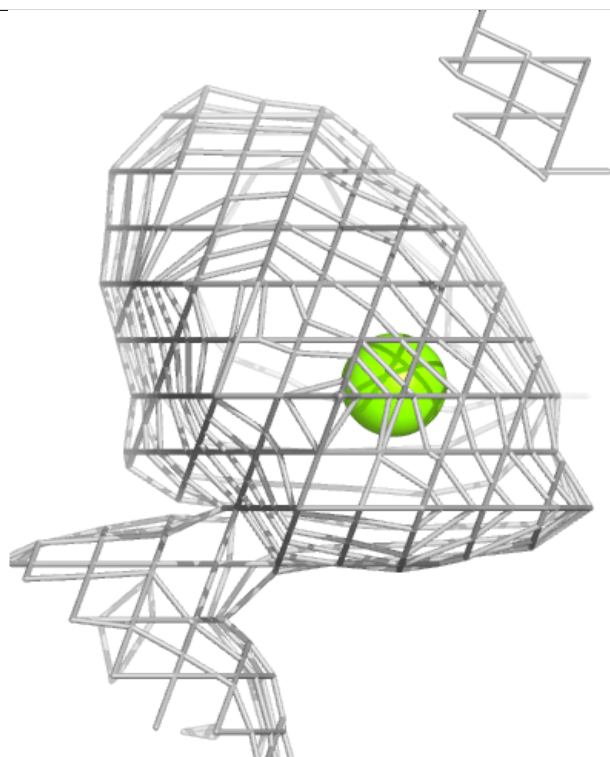
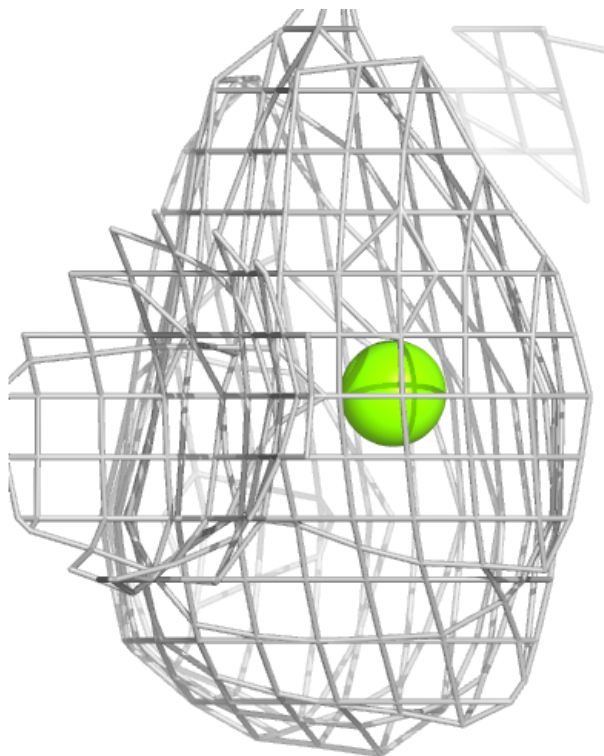
Electron density around MG A 302:

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and green (positive)



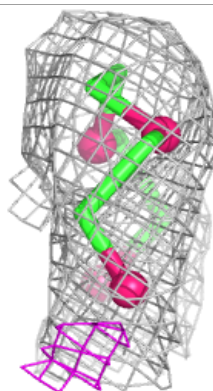
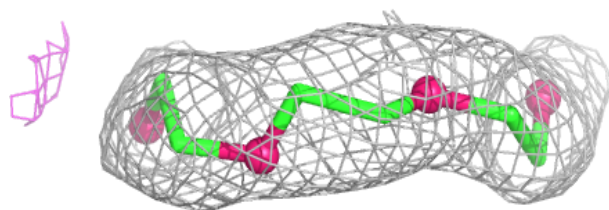
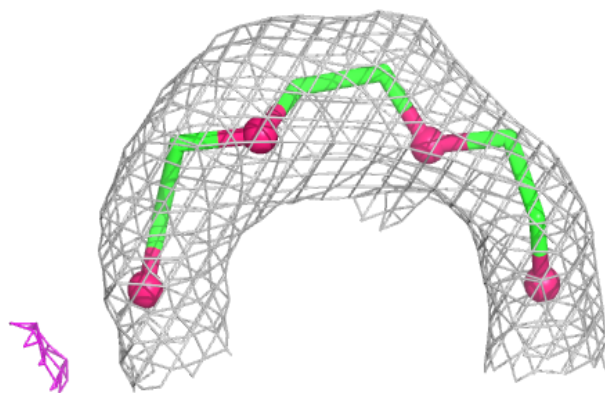
Electron density around MG C 303:

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



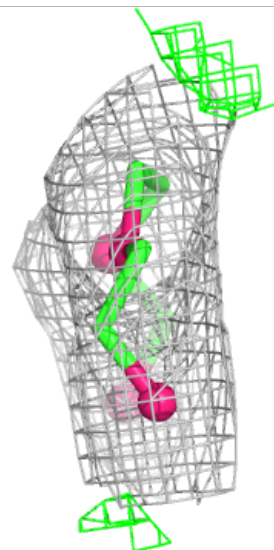
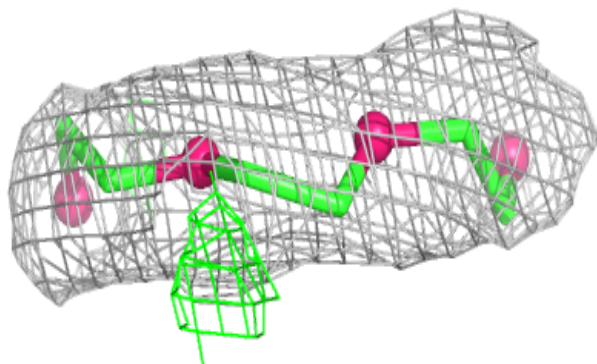
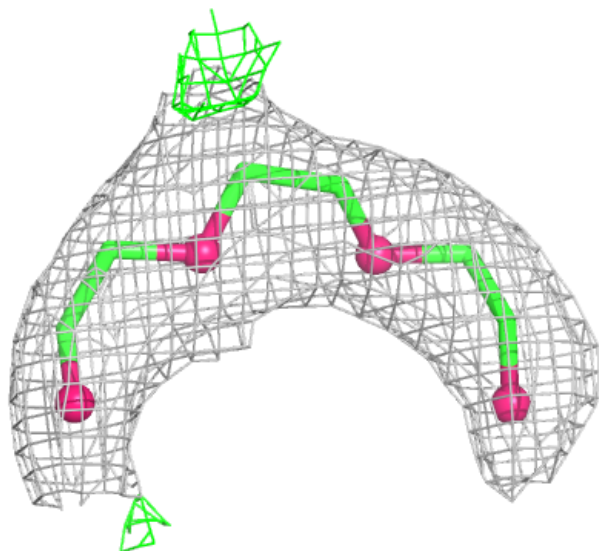
Electron density around PGE B 301:

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



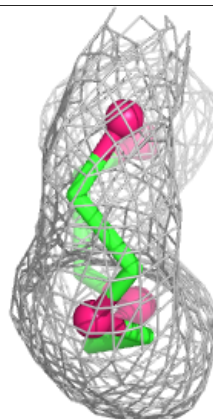
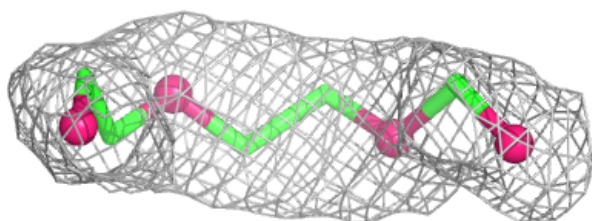
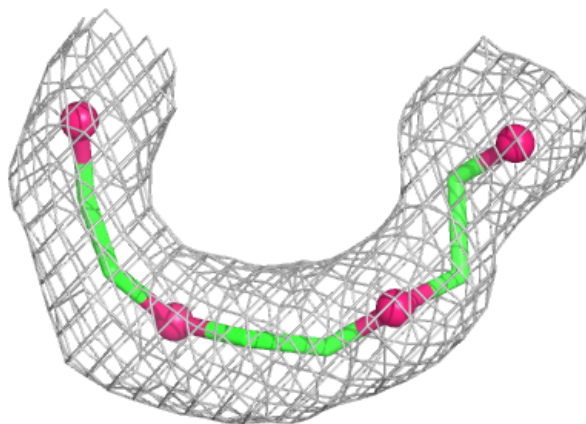
Electron density around PGE C 301:

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

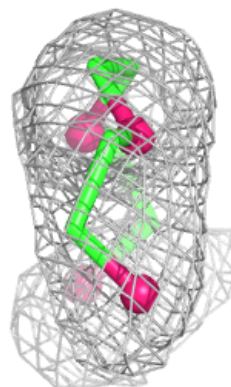
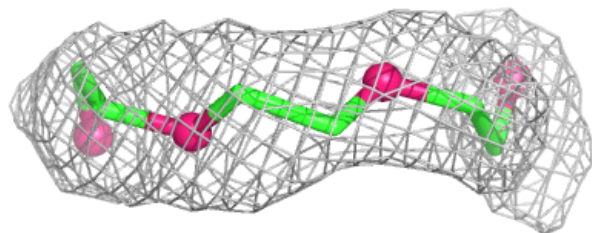
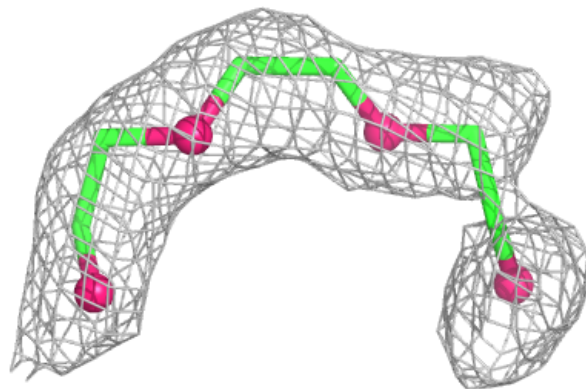


Electron density around PGE A 301:

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

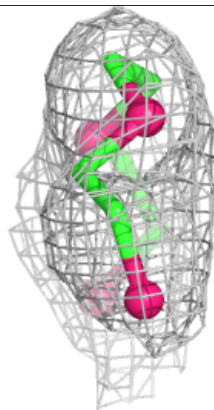
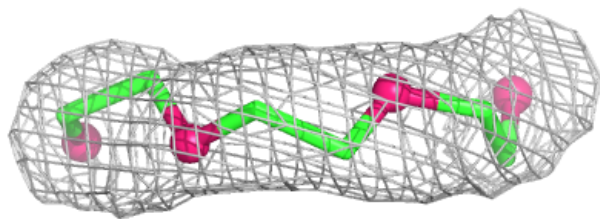
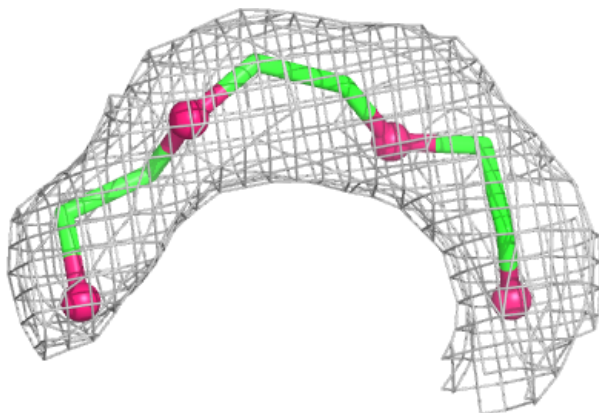
**Electron density around PGE G 301:**

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



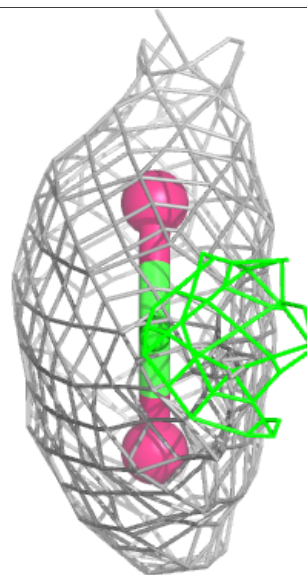
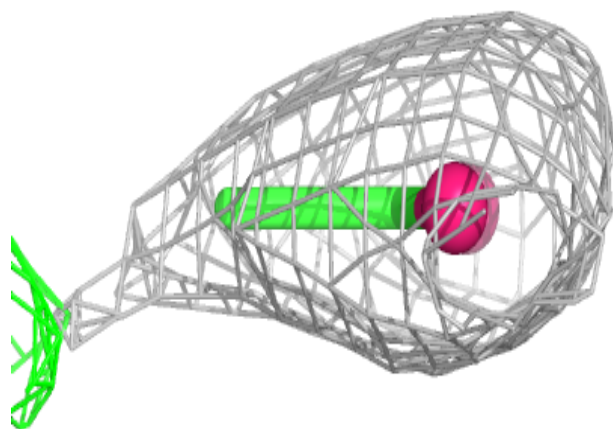
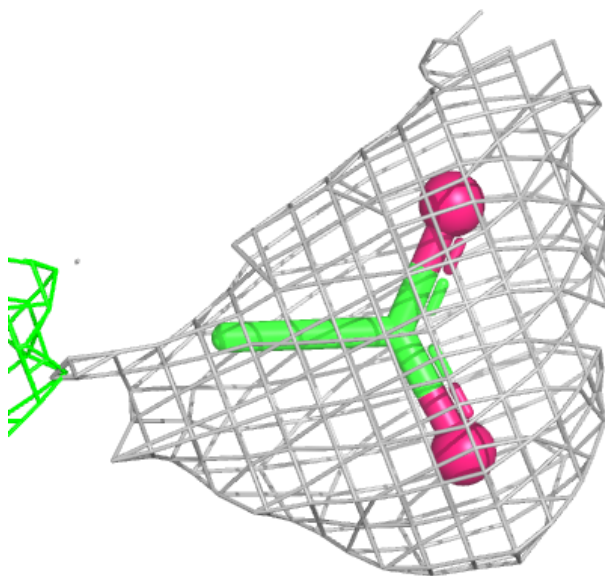
Electron density around PGE H 301:

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



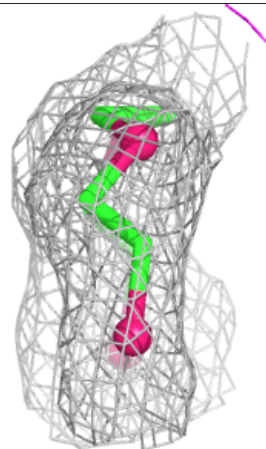
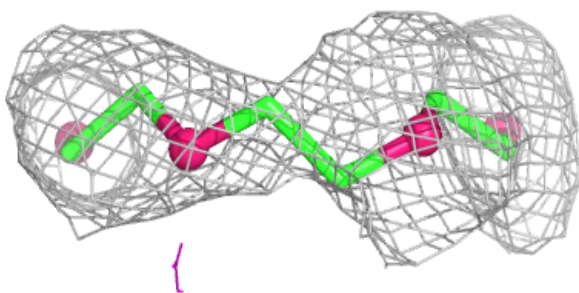
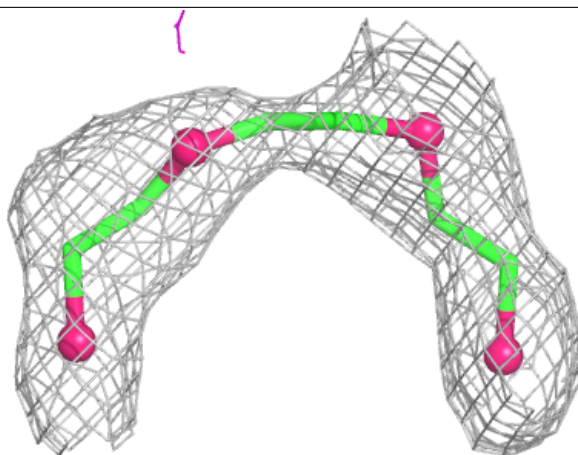
Electron density around ACT F 303:

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



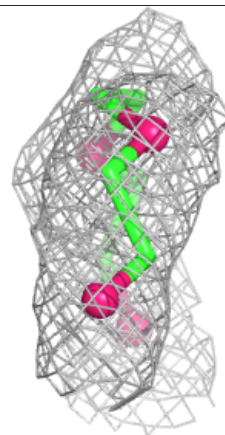
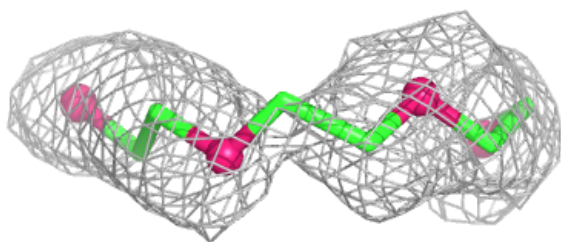
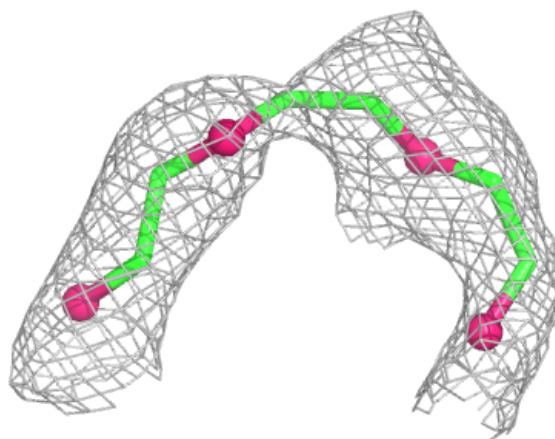
Electron density around PGE I 301:

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



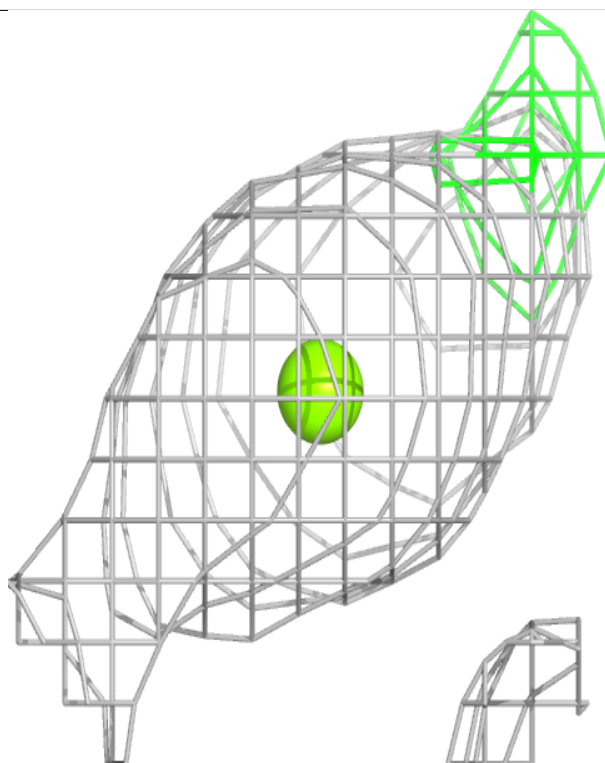
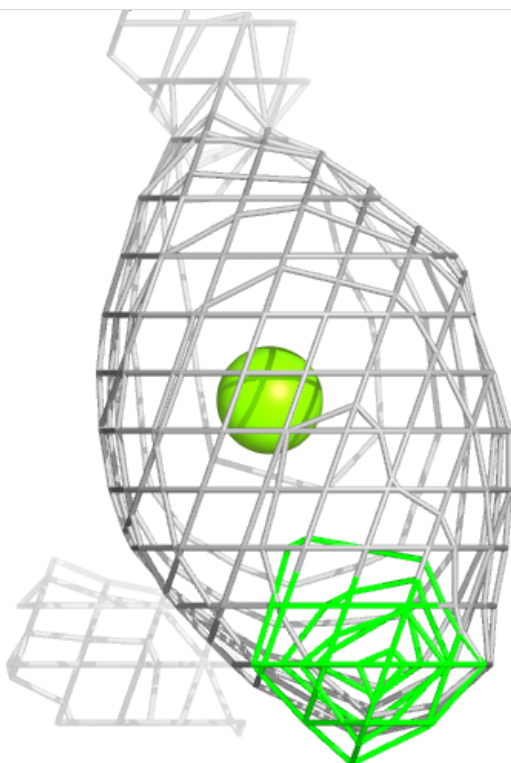
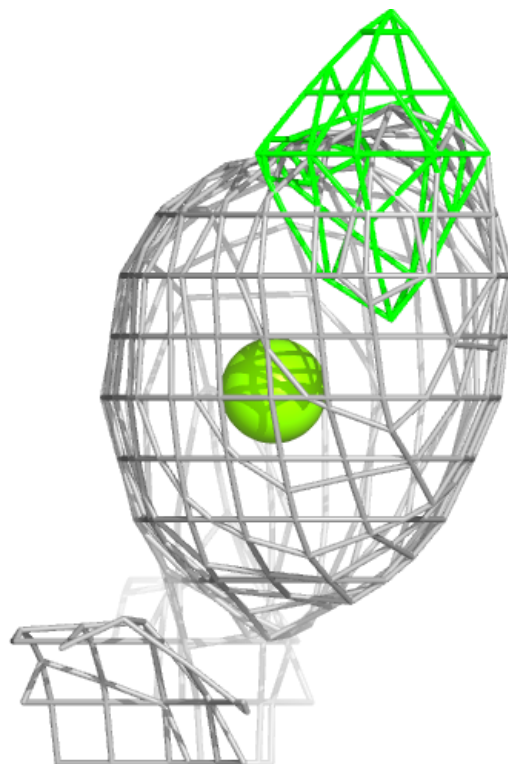
Electron density around PGE D 301:

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



Electron density around MG C 304:

$2mF_o-DF_c$ (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.