



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 07:59 AM UTC

PDB ID : 1YE9 / pdb_00001ye9
Title : Crystal structure of proteolytically truncated catalase HPII from E. coli
Authors : Loewen, P.C.; Chelikani, P.; Carpena, X.; Fita, I.; Perez-Luque, R.; Donald, L.J.; Switala, J.; Duckworth, H.W.
Deposited on : 2004-12-28
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

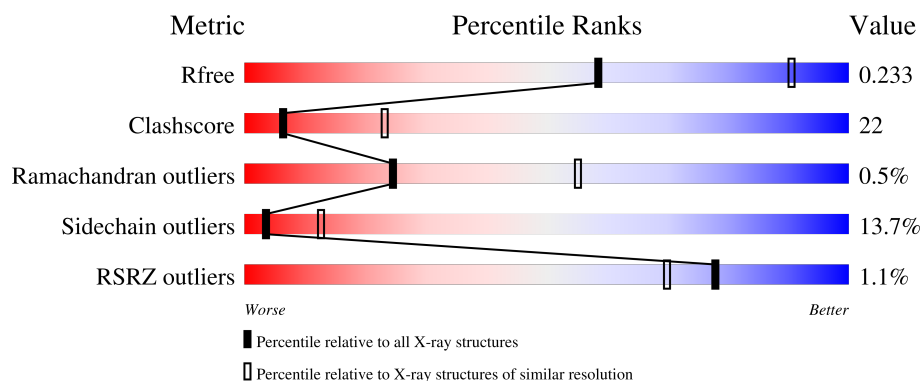
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	226	 61% 34% 5% . .
1	B	226	 55% 38% 5% 2% .
1	C	226	 66% 29% . .
1	D	226	 61% 33% . .
1	I	226	 56% 35% 8% .

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Mol	Chain	Length	Quality of chain
1	J	226	
1	K	226	
1	L	226	
2	E	259	
2	F	259	
2	G	259	
2	H	259	
2	M	259	
2	N	259	
2	O	259	
2	P	259	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	HDD	F	760	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 31715 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 catalase HPIL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	0	0
			1796	1148	319	326	3			
1	B	223	Total	C	N	O	S	0	1	0
			1803	1152	322	326	3			
1	C	223	Total	C	N	O	S	0	1	0
			1800	1151	320	326	3			
1	D	223	Total	C	N	O	S	0	1	0
			1802	1151	322	326	3			
1	I	223	Total	C	N	O	S	0	0	0
			1796	1148	319	326	3			
1	J	223	Total	C	N	O	S	0	0	0
			1796	1148	319	326	3			
1	K	223	Total	C	N	O	S	0	0	0
			1796	1148	319	326	3			
1	L	223	Total	C	N	O	S	0	1	0
			1801	1150	322	326	3			

- Molecule 2 is a protein called catalase HPIL.

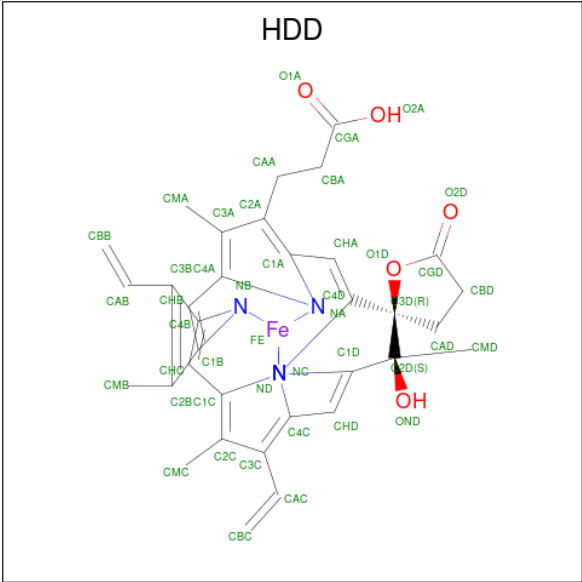
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	256	Total	C	N	O	S	0	3	0
			2107	1340	370	393	4			
2	F	256	Total	C	N	O	S	0	0	0
			2094	1333	367	390	4			
2	G	256	Total	C	N	O	S	0	2	0
			2106	1340	372	390	4			
2	H	256	Total	C	N	O	S	0	1	0
			2100	1337	369	390	4			
2	M	256	Total	C	N	O	S	0	1	0
			2098	1336	368	390	4			
2	N	256	Total	C	N	O	S	0	1	0
			2100	1337	369	390	4			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	O	256	Total	C	N	O	S	0	0	0
			2094	1333	367	390	4			
2	P	256	Total	C	N	O	S	0	0	0
			2094	1333	367	390	4			

- Molecule 3 is CIS-HEME D HYDROXYCHLORIN GAMMA-SPIROLACTONE (CCD ID: HDD) (formula: C₃₄H₃₂FeN₄O₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	E	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		
3	F	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		
3	G	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		
3	H	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		
3	M	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		
3	N	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		
3	O	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		
3	P	1	Total	C	Fe	N	O	0	0
			44	34	1	4	5		

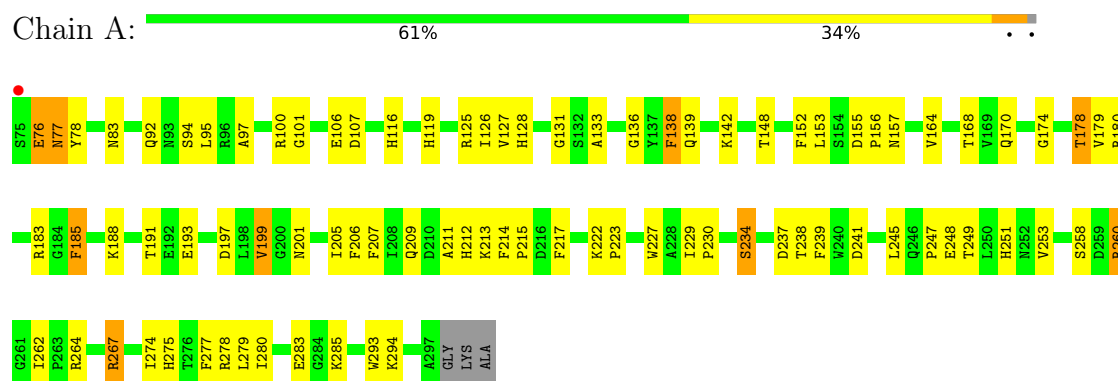
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	21	Total O 21 21	0	0
4	E	20	Total O 20 20	0	0
4	B	8	Total O 8 8	0	0
4	F	11	Total O 11 11	0	0
4	C	20	Total O 20 20	0	0
4	G	11	Total O 11 11	0	0
4	D	11	Total O 11 11	0	0
4	H	11	Total O 11 11	0	0
4	I	5	Total O 5 5	0	0
4	M	2	Total O 2 2	0	0
4	J	12	Total O 12 12	0	0
4	N	12	Total O 12 12	0	0
4	K	3	Total O 3 3	0	0
4	O	4	Total O 4 4	0	0
4	L	16	Total O 16 16	0	0
4	P	13	Total O 13 13	0	0

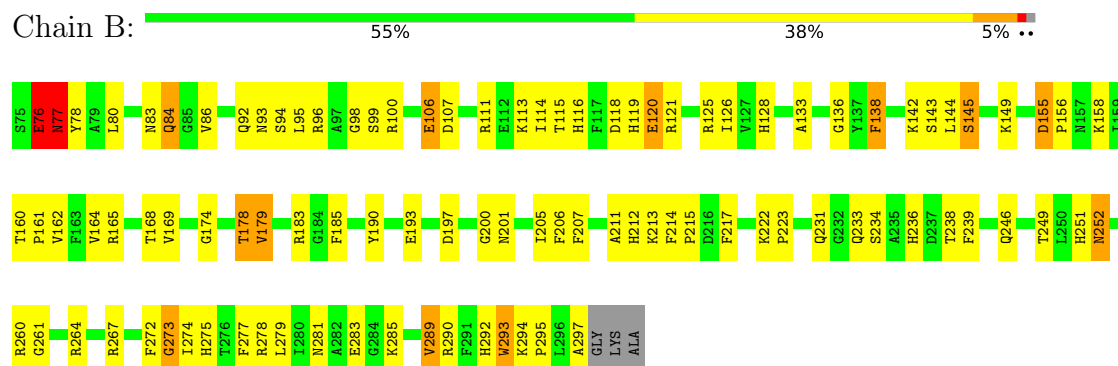
3 Residue-property plots

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

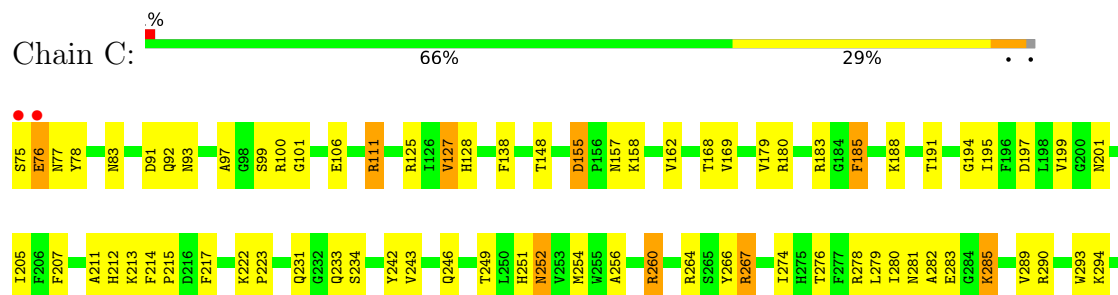
• Molecule 1: catalase HP11



• Molecule 1: catalase HP11



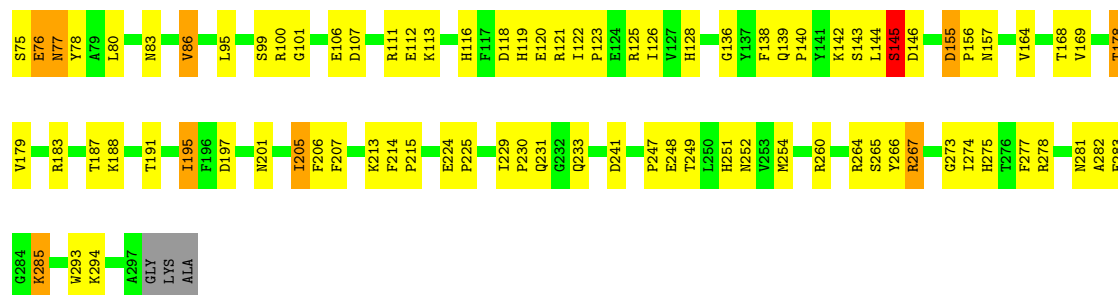
• Molecule 1: catalase HP11





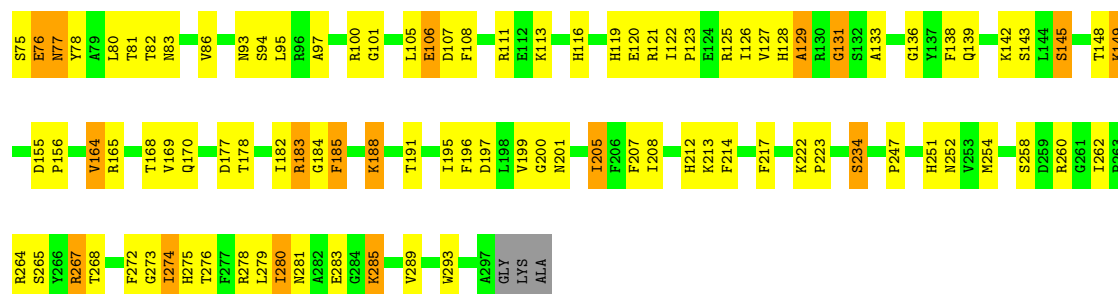
● Molecule 1: catalase HP1I

Chain D: 61% 33% . .



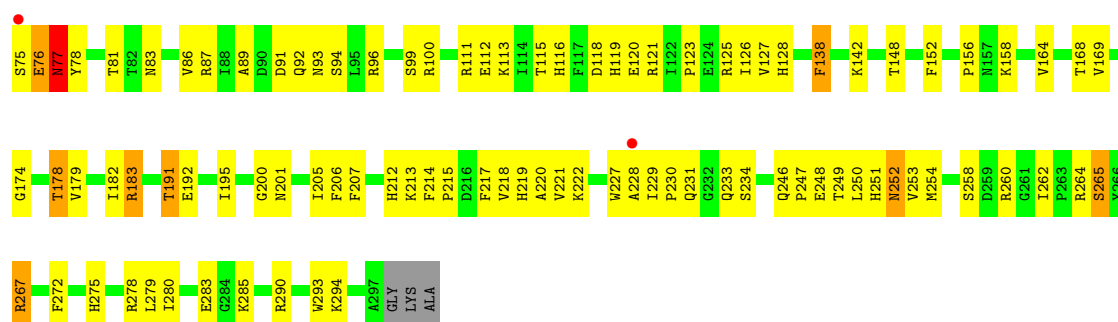
● Molecule 1: catalase HP1I

Chain I: 56% 35% 8% .



● Molecule 1: catalase HP1I

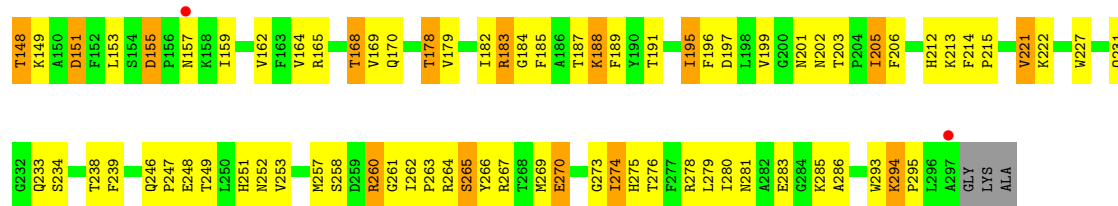
Chain J: 57% 38% . .



● Molecule 1: catalase HP1I

Chain K: 47% 43% 8% .

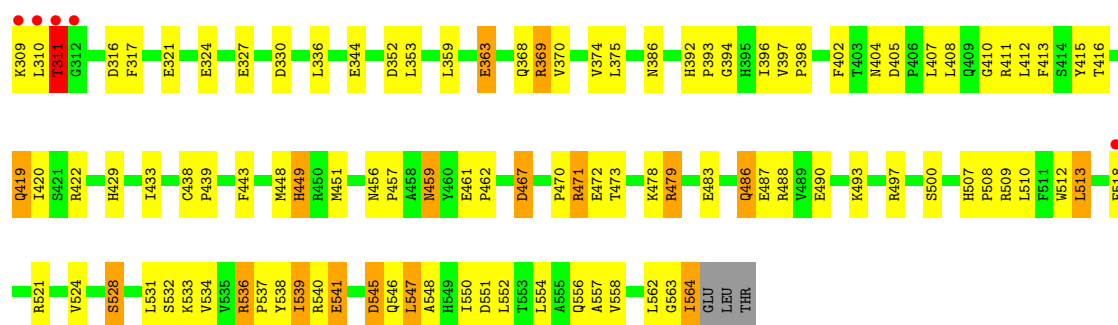




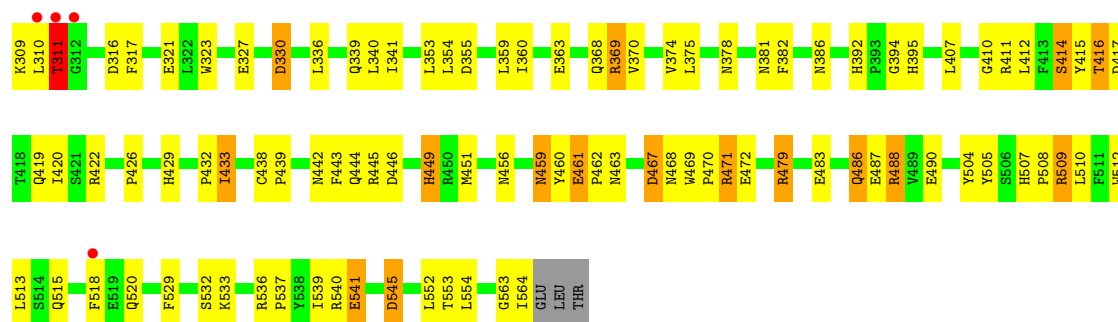
● Molecule 1: catalase HPII



● Molecule 2: catalase HPII

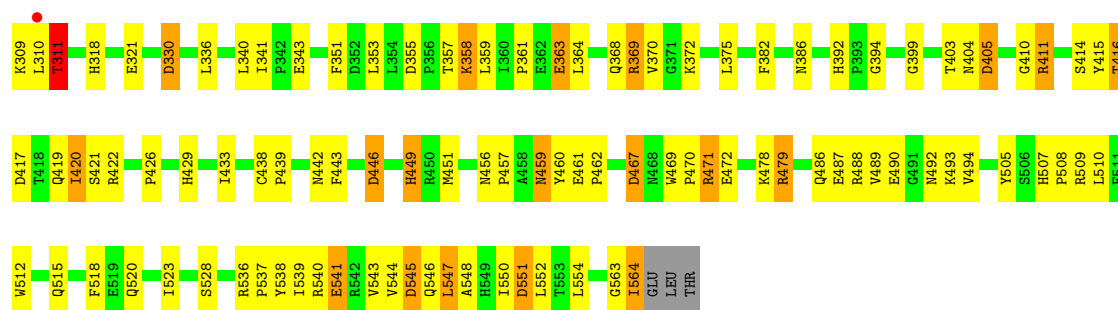


● Molecule 2: catalase HPII



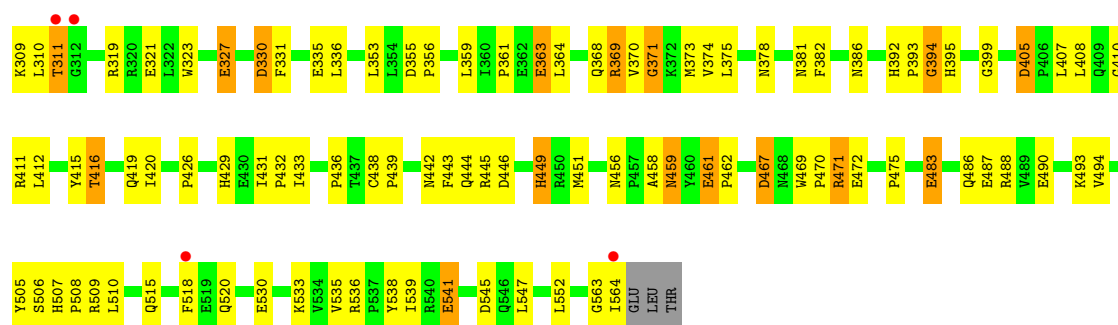
● Molecule 2: catalase HPII

Chain G:  59% 32% 7% .



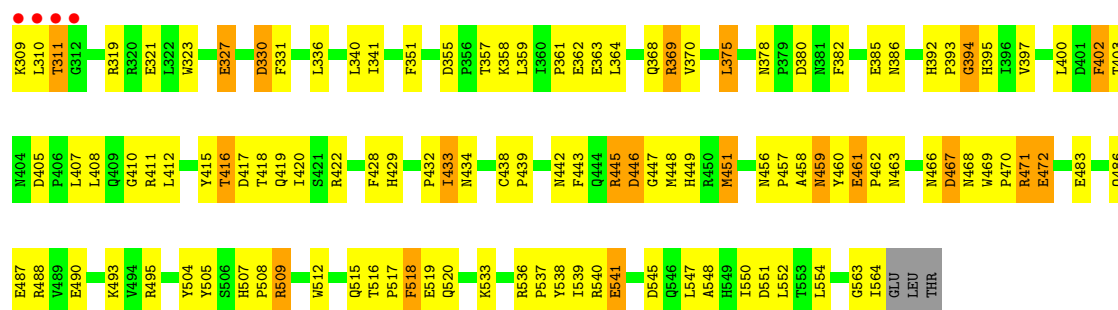
• Molecule 2: catalase HPII

Chain H:  61% 32% 6% .



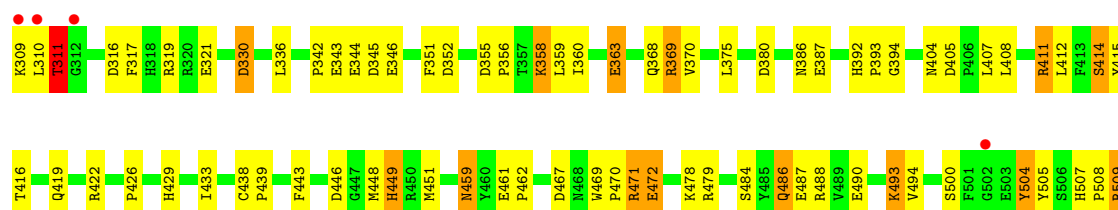
• Molecule 2: catalase HPII

Chain M:  54% 37% 8% .



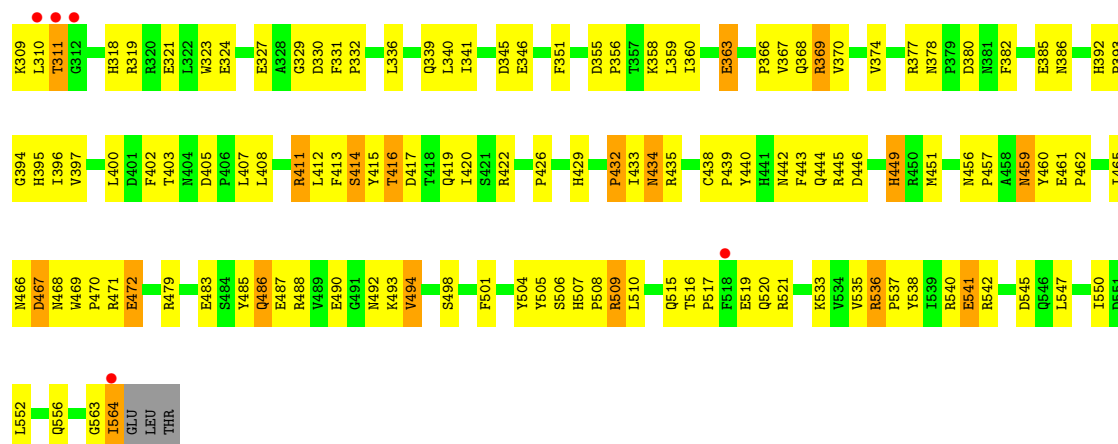
• Molecule 2: catalase HPII

Chain N:  60% 31% 8% .

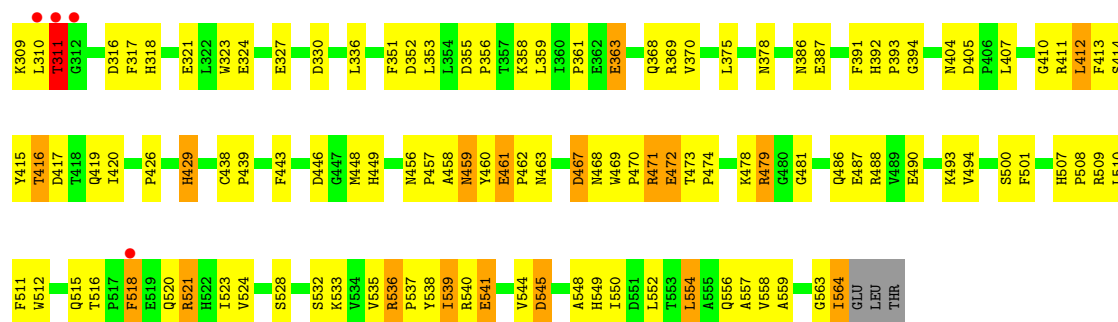




● Molecule 2: catalase HP11



● Molecule 2: catalase HP11



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	111.01Å 152.89Å 135.29Å 90.00° 97.54° 90.00°	Depositor
Resolution (Å)	30.00 – 2.80 30.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.00-2.80) 95.3 (30.00-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.217 , 0.269 (Not available) , 0.233	Depositor DCC
R_{free} test set	5289 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	51.6	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	31715	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: HDD

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	0.94	0/1851	1.09	4/2513 (0.2%)
1	B	0.91	1/1862 (0.1%)	1.15	8/2527 (0.3%)
1	C	0.92	0/1860	1.11	3/2525 (0.1%)
1	D	0.91	1/1862 (0.1%)	1.15	2/2527 (0.1%)
1	I	0.95	1/1851 (0.1%)	1.17	9/2513 (0.4%)
1	J	0.89	2/1851 (0.1%)	1.15	3/2513 (0.1%)
1	K	0.98	0/1851	1.16	6/2513 (0.2%)
1	L	0.88	0/1862	1.09	3/2527 (0.1%)
2	E	0.90	1/2188 (0.0%)	1.07	0/2975
2	F	0.88	1/2159 (0.0%)	1.10	4/2936 (0.1%)
2	G	0.92	2/2181 (0.1%)	1.08	5/2965 (0.2%)
2	H	0.89	0/2170	1.11	4/2951 (0.1%)
2	M	0.88	0/2168	1.12	5/2947 (0.2%)
2	N	0.89	2/2170 (0.1%)	1.08	4/2951 (0.1%)
2	O	0.91	1/2159 (0.0%)	1.13	5/2936 (0.2%)
2	P	0.88	1/2159 (0.0%)	1.09	4/2936 (0.1%)
All	All	0.91	13/32204 (0.0%)	1.11	69/43755 (0.2%)

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	311	THR	CA-CB	6.70	1.65	1.53
1	B	178	THR	CA-C	-6.60	1.46	1.53
2	O	451	MET	SD-CE	6.08	1.94	1.79
1	J	164	VAL	CA-CB	-5.94	1.47	1.54
2	G	442	ASN	C-O	-5.69	1.18	1.24

The worst 5 of 69 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	76	GLU	N-CA-C	7.40	122.37	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	414	SER	N-CA-C	6.79	118.48	111.14
2	H	371	GLY	N-CA-C	6.67	121.62	110.56
2	M	428	PHE	N-CA-C	-6.64	105.14	113.18
2	M	394	GLY	N-CA-C	-6.62	107.38	114.67

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1796	0	1720	77	0
1	B	1803	0	1729	106	0
1	C	1800	0	1729	68	0
1	D	1802	0	1729	99	1
1	I	1796	0	1720	102	0
1	J	1796	0	1720	80	0
1	K	1796	0	1720	117	0
1	L	1801	0	1725	85	0
2	E	2107	0	1980	84	0
2	F	2094	0	1975	109	0
2	G	2106	0	1987	97	0
2	H	2100	0	1978	92	0
2	M	2098	0	1984	128	0
2	N	2100	0	1978	99	0
2	O	2094	0	1975	138	3
2	P	2094	0	1975	123	2
3	E	44	0	28	12	0
3	F	44	0	28	25	0
3	G	44	0	28	8	0
3	H	44	0	28	14	0
3	M	44	0	28	15	0
3	N	44	0	28	8	0
3	O	44	0	28	20	0
3	P	44	0	28	9	0
4	A	21	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	8	0	0	3	0
4	C	20	0	0	5	0
4	D	11	0	0	5	0
4	E	20	0	0	6	0
4	F	11	0	0	9	0
4	G	11	0	0	1	0
4	H	11	0	0	2	0
4	I	5	0	0	4	0
4	J	12	0	0	4	0
4	K	3	0	0	1	0
4	L	16	0	0	4	0
4	M	2	0	0	2	0
4	N	12	0	0	3	0
4	O	4	0	0	2	0
4	P	13	0	0	4	0
All	All	31715	0	29848	1352	3

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 22.

The worst 5 of 1352 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:L:267:ARG:HH11	1:L:267:ARG:CB	1.11	1.61
2:O:392:HIS:ND1	2:O:415:TYR:CB	1.71	1.53
2:P:392:HIS:ND1	2:P:415:TYR:CB	1.72	1.50
2:G:392:HIS:ND1	2:G:415:TYR:CB	1.74	1.48
2:N:392:HIS:ND1	2:N:415:TYR:CB	1.74	1.48

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:509:ARG:NH1	2:P:521:ARG:NE[2_756]	1.70	0.50
1:D:157:ASN:OD1	2:O:369:ARG:NH2[2_645]	2.16	0.04
2:O:509:ARG:NH1	2:P:521:ARG:CD[2_756]	2.19	0.01

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	221/226 (98%)	208 (94%)	12 (5%)	1 (0%)	24	55
1	B	222/226 (98%)	207 (93%)	13 (6%)	2 (1%)	14	41
1	C	222/226 (98%)	214 (96%)	8 (4%)	0	100	100
1	D	222/226 (98%)	209 (94%)	11 (5%)	2 (1%)	14	41
1	I	221/226 (98%)	206 (93%)	13 (6%)	2 (1%)	14	41
1	J	221/226 (98%)	207 (94%)	14 (6%)	0	100	100
1	K	221/226 (98%)	202 (91%)	17 (8%)	2 (1%)	14	41
1	L	222/226 (98%)	206 (93%)	15 (7%)	1 (0%)	24	55
2	E	257/259 (99%)	240 (93%)	15 (6%)	2 (1%)	16	44
2	F	254/259 (98%)	238 (94%)	15 (6%)	1 (0%)	30	60
2	G	256/259 (99%)	237 (93%)	19 (7%)	0	100	100
2	H	255/259 (98%)	237 (93%)	18 (7%)	0	100	100
2	M	255/259 (98%)	230 (90%)	23 (9%)	2 (1%)	16	44
2	N	255/259 (98%)	237 (93%)	17 (7%)	1 (0%)	30	60
2	O	254/259 (98%)	228 (90%)	23 (9%)	3 (1%)	10	34
2	P	254/259 (98%)	233 (92%)	19 (8%)	2 (1%)	16	44
All	All	3812/3880 (98%)	3539 (93%)	252 (7%)	21 (1%)	24	51

5 of 21 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	SER
1	I	274	ILE
1	B	77	ASN
1	B	274	ILE
1	D	145	SER

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	189/190 (100%)	172 (91%)	17 (9%)	9	28
1	B	190/190 (100%)	165 (87%)	25 (13%)	4	14
1	C	190/190 (100%)	169 (89%)	21 (11%)	6	20
1	D	190/190 (100%)	170 (90%)	20 (10%)	6	22
1	I	189/190 (100%)	162 (86%)	27 (14%)	3	11
1	J	189/190 (100%)	165 (87%)	24 (13%)	4	15
1	K	189/190 (100%)	155 (82%)	34 (18%)	2	6
1	L	190/190 (100%)	169 (89%)	21 (11%)	6	20
2	E	229/231 (99%)	190 (83%)	39 (17%)	2	7
2	F	226/231 (98%)	195 (86%)	31 (14%)	3	13
2	G	228/231 (99%)	194 (85%)	34 (15%)	3	10
2	H	227/231 (98%)	196 (86%)	31 (14%)	3	13
2	M	227/231 (98%)	196 (86%)	31 (14%)	3	13
2	N	227/231 (98%)	190 (84%)	37 (16%)	2	8
2	O	226/231 (98%)	191 (84%)	35 (16%)	2	9
2	P	226/231 (98%)	193 (85%)	33 (15%)	3	11
All	All	3332/3368 (99%)	2872 (86%)	460 (14%)	3	12

5 of 460 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	188	LYS
2	P	479	ARG
1	J	191	THR
2	P	459	ASN
2	O	533	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 61 such sidechains are listed below:

Mol	Chain	Res	Type
2	H	466	ASN
2	O	459	ASN
2	M	449	HIS
2	O	449	HIS
2	P	419	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 ligands 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
3	HDD	P	760	2	46,52,52	2.93	24 (52%)	62,89,89	4.79	39 (62%)
3	HDD	O	760	2	46,52,52	2.89	24 (52%)	62,89,89	4.62	33 (53%)
3	HDD	H	760	2	46,52,52	2.90	23 (50%)	62,89,89	4.33	32 (51%)
3	HDD	E	760	2	46,52,52	3.11	24 (52%)	62,89,89	4.95	36 (58%)
3	HDD	M	760	2	46,52,52	2.76	23 (50%)	62,89,89	4.54	32 (51%)
3	HDD	F	760	2	46,52,52	2.64	21 (45%)	62,89,89	4.57	34 (54%)
3	HDD	N	760	2	46,52,52	3.03	22 (47%)	62,89,89	4.49	34 (54%)
3	HDD	G	760	2	46,52,52	2.80	25 (54%)	62,89,89	4.16	29 (46%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HDD	P	760	2	-	4/9/89/89	0/1/9/9
3	HDD	O	760	2	-	8/9/89/89	0/1/9/9
3	HDD	H	760	2	-	4/9/89/89	0/1/9/9
3	HDD	E	760	2	-	2/9/89/89	0/1/9/9
3	HDD	M	760	2	-	2/9/89/89	0/1/9/9
3	HDD	F	760	2	-	5/9/89/89	0/1/9/9
3	HDD	N	760	2	-	2/9/89/89	0/1/9/9
3	HDD	G	760	2	-	5/9/89/89	0/1/9/9

The worst 5 of 186 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	760	HDD	O1D-CGD	7.50	1.47	1.35
3	E	760	HDD	C4C-NC	-7.03	1.26	1.39
3	O	760	HDD	O1D-CGD	6.67	1.46	1.35
3	H	760	HDD	O1D-CGD	6.63	1.46	1.35
3	F	760	HDD	O1D-CGD	6.62	1.46	1.35

The worst 5 of 269 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	O	760	HDD	C1A-CHA-C4D	16.46	157.77	125.81
3	E	760	HDD	C4C-CHD-C1D	16.37	157.61	125.81
3	E	760	HDD	C1A-CHA-C4D	15.91	156.71	125.81
3	M	760	HDD	C1A-CHA-C4D	15.83	156.55	125.81
3	N	760	HDD	C4C-CHD-C1D	15.71	156.33	125.81

There are no chirality outliers.

5 of 32 torsion outliers are listed below:

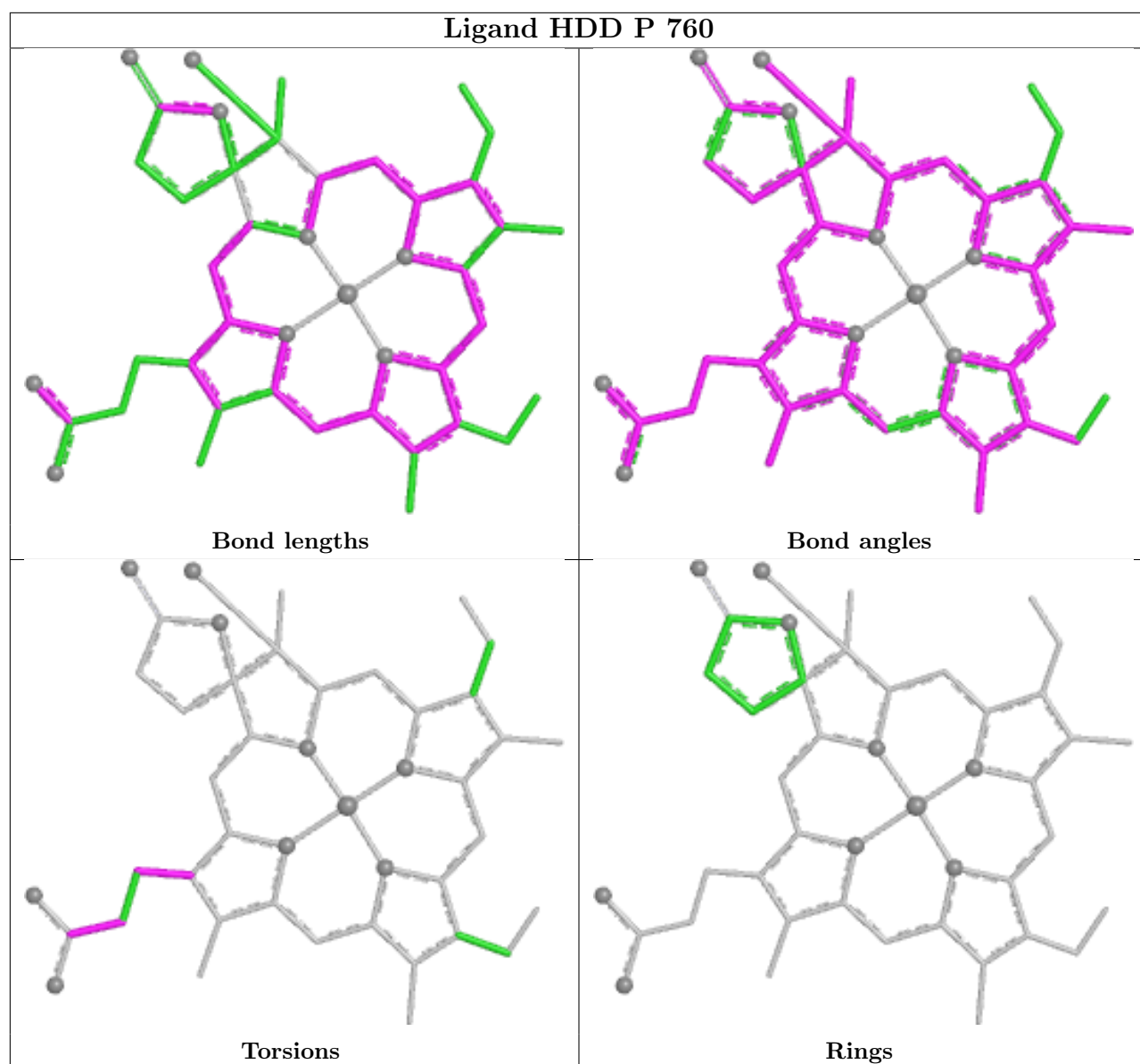
Mol	Chain	Res	Type	Atoms
3	G	760	HDD	C2C-C3C-CAC-CBC
3	G	760	HDD	C4C-C3C-CAC-CBC
3	H	760	HDD	C2B-C3B-CAB-CBB
3	O	760	HDD	C2B-C3B-CAB-CBB
3	O	760	HDD	C4B-C3B-CAB-CBB

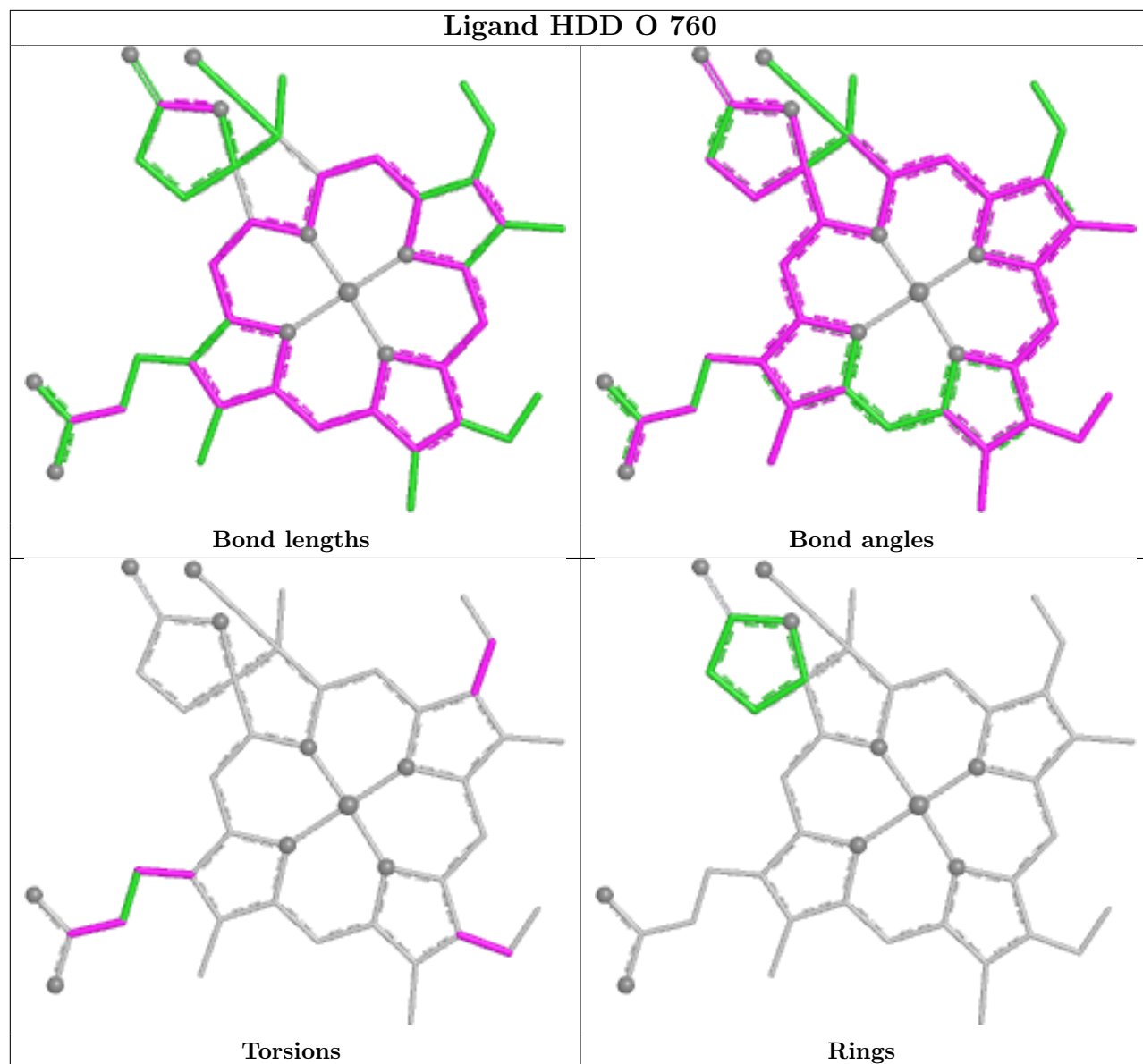
There are no ring outliers.

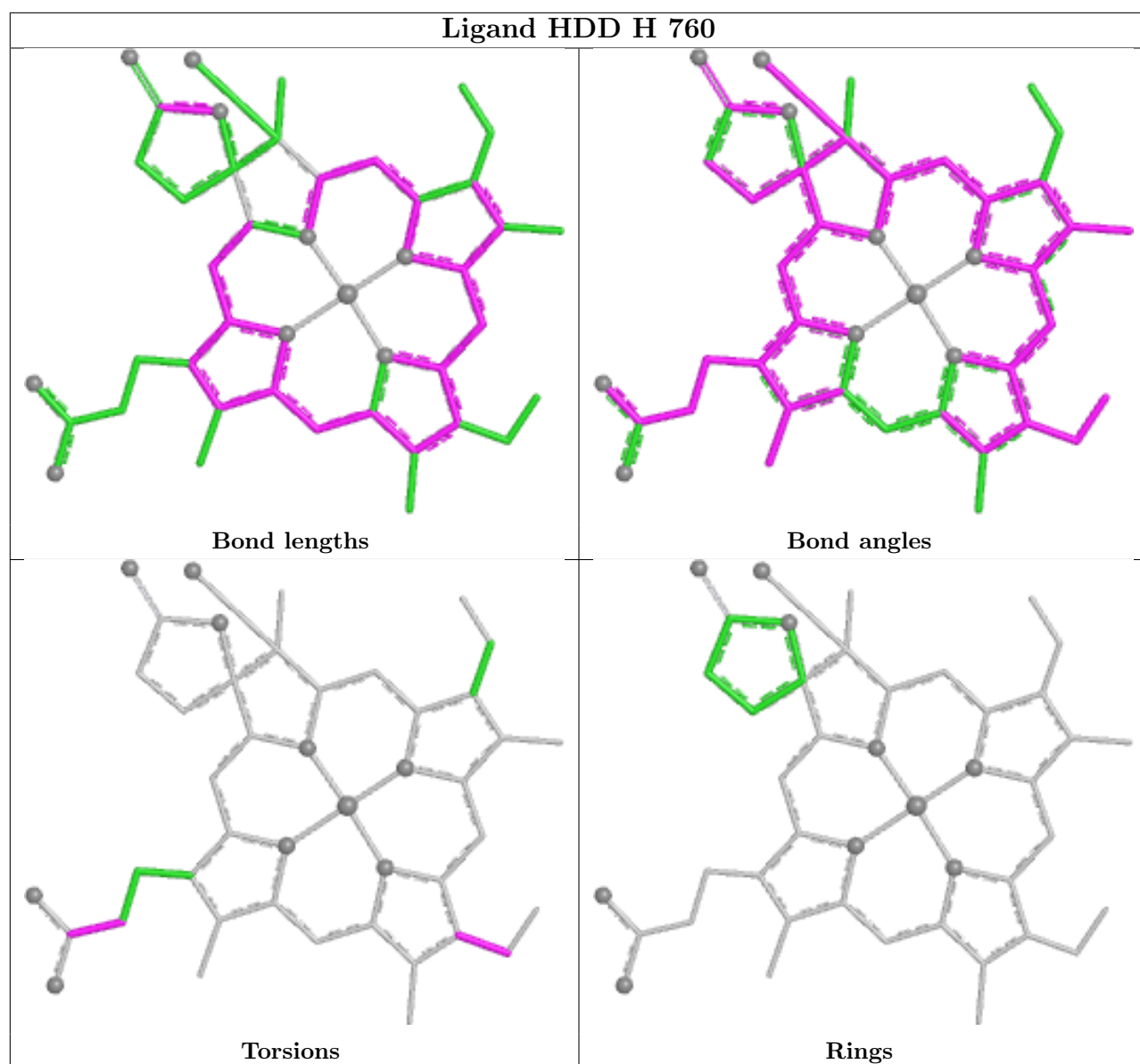
8 monomers are involved in 111 short contacts:

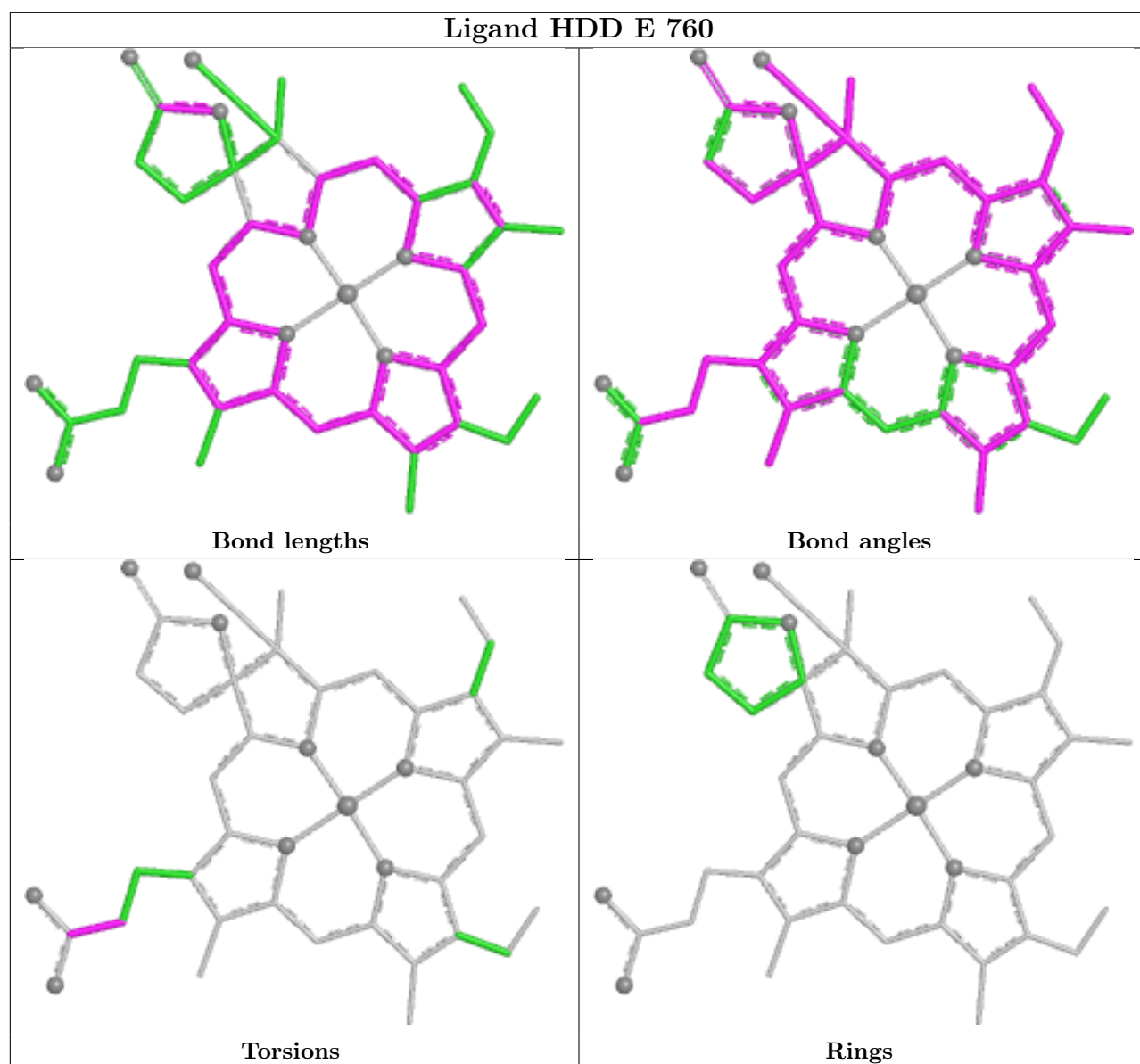
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	760	HDD	9	0
3	O	760	HDD	20	0
3	H	760	HDD	14	0
3	E	760	HDD	12	0
3	M	760	HDD	15	0
3	F	760	HDD	25	0
3	N	760	HDD	8	0
3	G	760	HDD	8	0

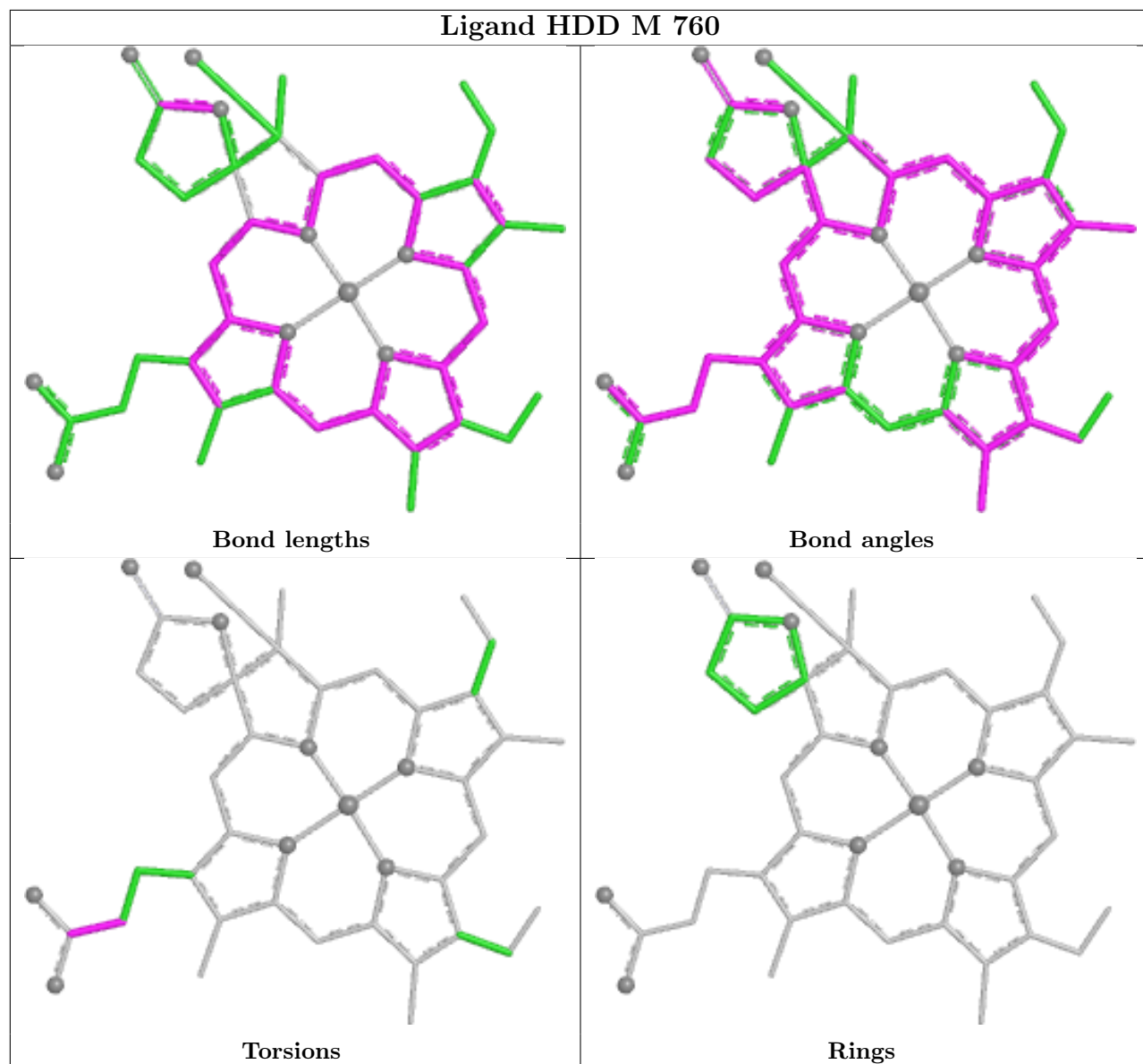
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

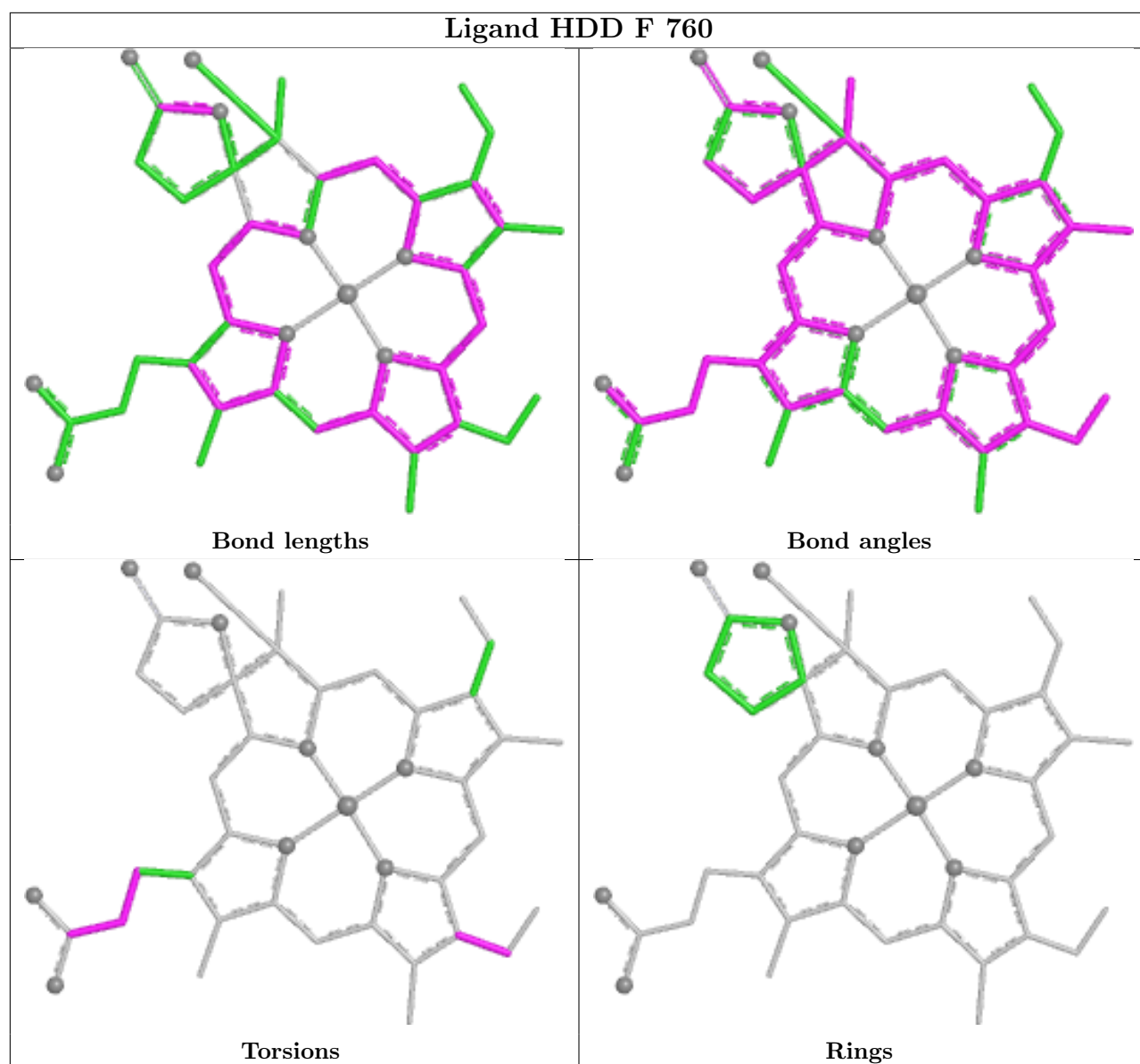


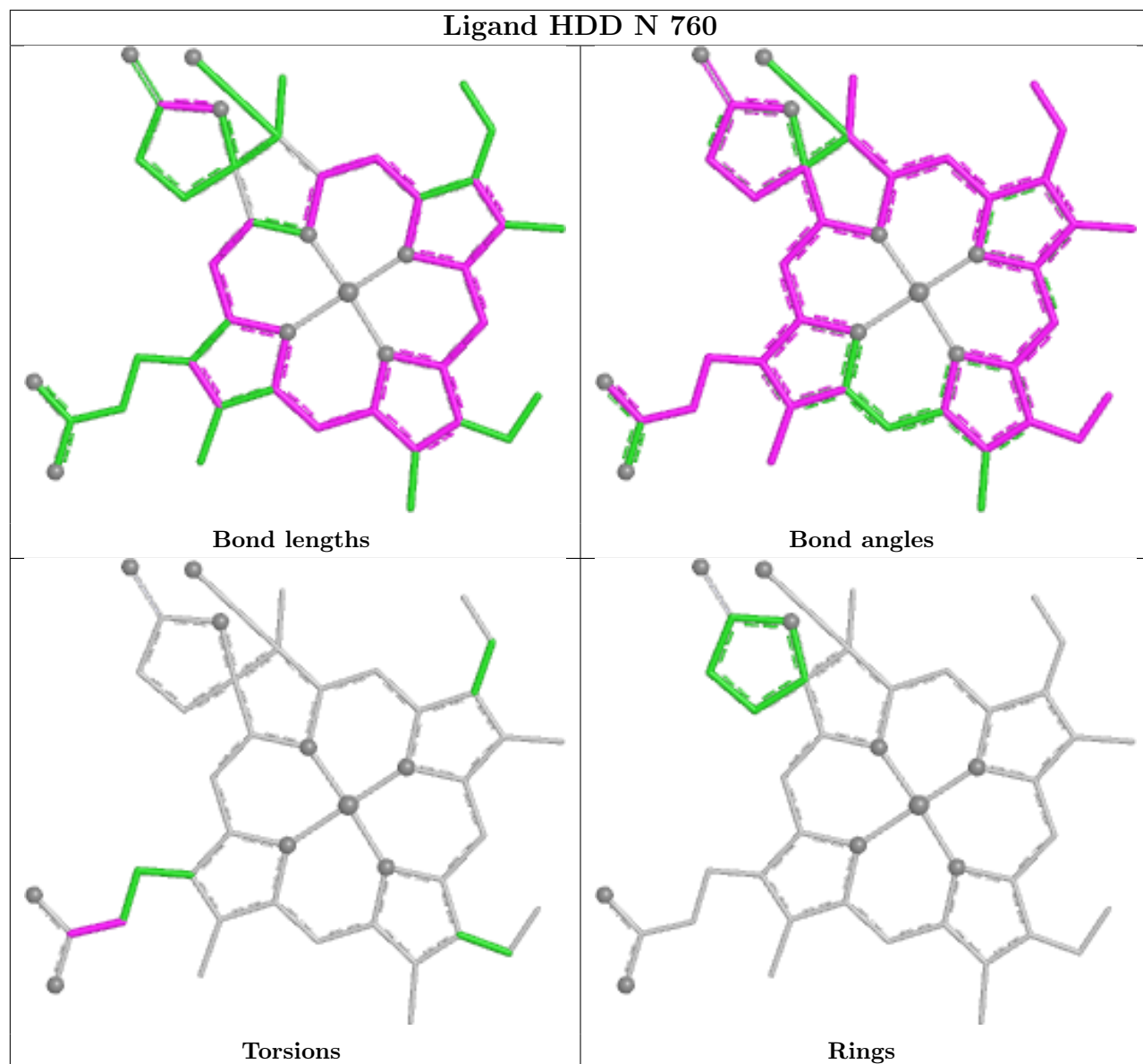


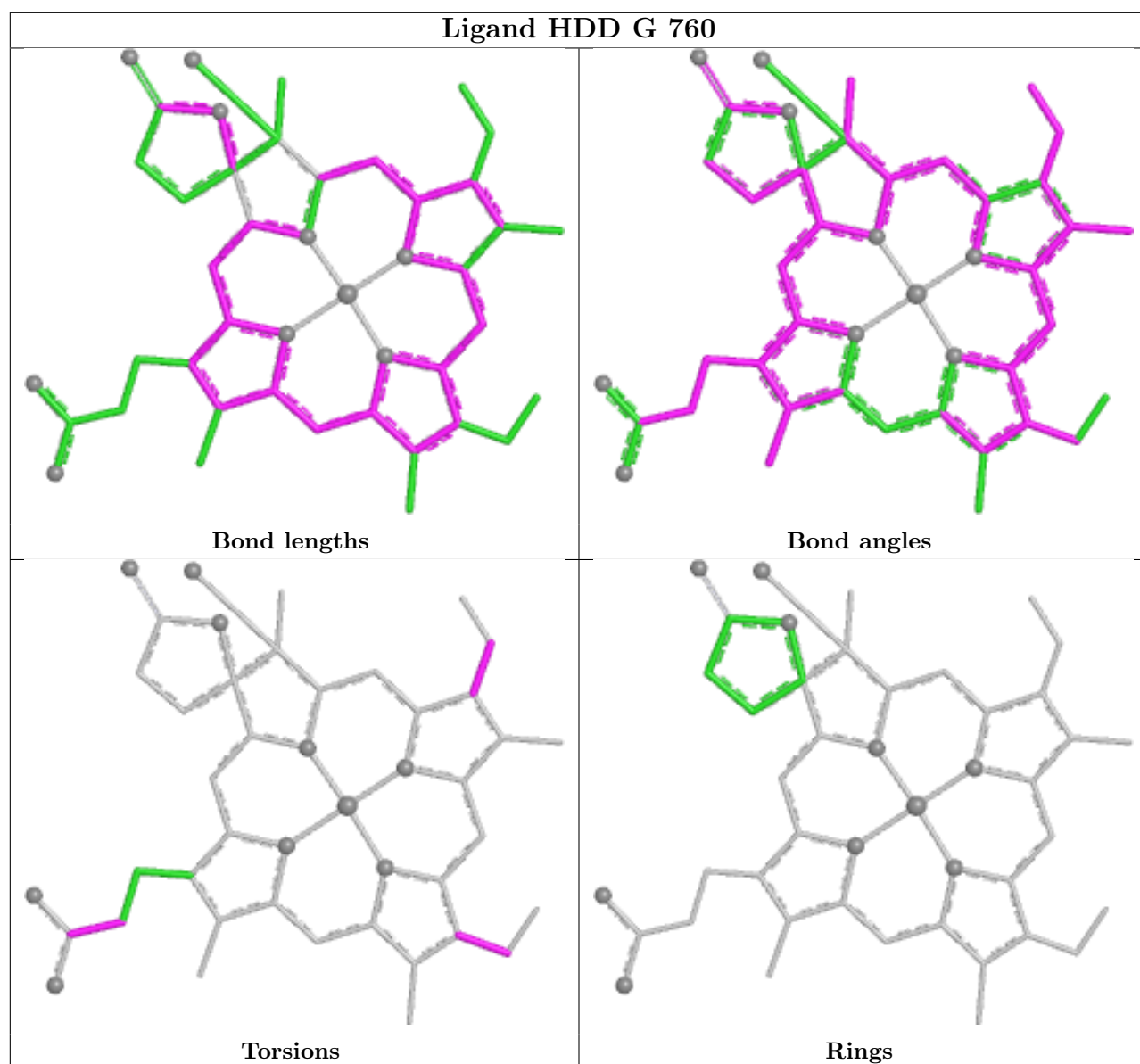












5.7 Other polymers [i](#)

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	223/226 (98%)	-0.14	1 (0%) 88 84	28, 40, 51, 80	0
1	B	223/226 (98%)	-0.27	0 100 100	26, 40, 51, 80	1 (0%)
1	C	223/226 (98%)	-0.21	3 (1%) 75 66	28, 40, 51, 80	1 (0%)
1	D	223/226 (98%)	-0.21	0 100 100	28, 40, 51, 80	1 (0%)
1	I	223/226 (98%)	-0.04	0 100 100	28, 41, 53, 80	0
1	J	223/226 (98%)	-0.18	2 (0%) 81 74	28, 41, 51, 80	0
1	K	223/226 (98%)	0.07	3 (1%) 75 66	28, 41, 53, 80	0
1	L	223/226 (98%)	-0.17	2 (0%) 81 74	25, 40, 52, 80	1 (0%)
2	E	256/259 (98%)	-0.03	5 (1%) 65 56	14, 44, 61, 81	3 (1%)
2	F	256/259 (98%)	-0.08	4 (1%) 70 61	29, 45, 61, 81	0
2	G	256/259 (98%)	-0.05	1 (0%) 88 84	19, 45, 61, 81	2 (0%)
2	H	256/259 (98%)	-0.06	4 (1%) 70 61	19, 45, 61, 81	1 (0%)
2	M	256/259 (98%)	0.07	4 (1%) 70 61	22, 45, 61, 81	1 (0%)
2	N	256/259 (98%)	-0.08	4 (1%) 70 61	20, 45, 61, 81	1 (0%)
2	O	256/259 (98%)	0.15	5 (1%) 65 56	30, 46, 61, 81	0
2	P	256/259 (98%)	-0.00	4 (1%) 70 61	29, 45, 61, 81	0
All	All	3832/3880 (98%)	-0.07	42 (1%) 78 70	14, 42, 59, 81	12 (0%)

The worst 5 of 42 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	O	312	GLY	4.2
1	A	75	SER	4.1
2	E	311	THR	4.1
2	F	518	PHE	4.0
1	L	75	SER	3.8

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

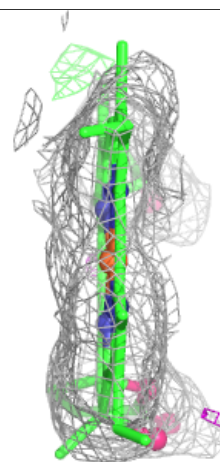
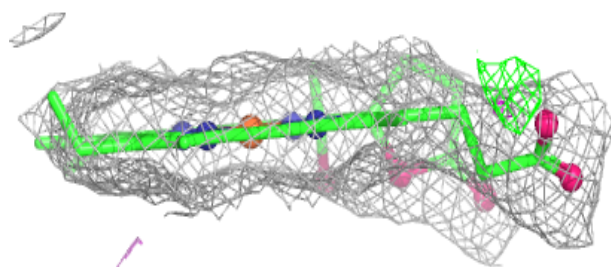
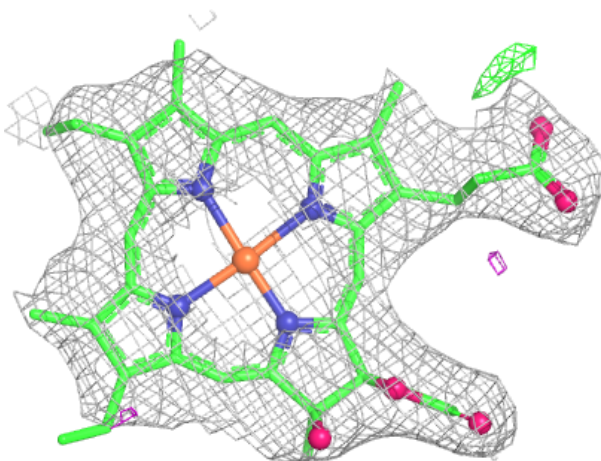
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	HDD	H	760	44/44	0.96	0.09	31,43,49,58	0
3	HDD	M	760	44/44	0.96	0.10	24,52,64,66	0
3	HDD	N	760	44/44	0.96	0.08	25,38,47,57	0
3	HDD	O	760	44/44	0.96	0.09	32,53,62,63	0
3	HDD	F	760	44/44	0.97	0.07	10,24,39,42	0
3	HDD	G	760	44/44	0.97	0.08	24,35,48,55	0
3	HDD	P	760	44/44	0.97	0.08	12,28,43,45	0
3	HDD	E	760	44/44	0.98	0.07	10,23,34,42	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.

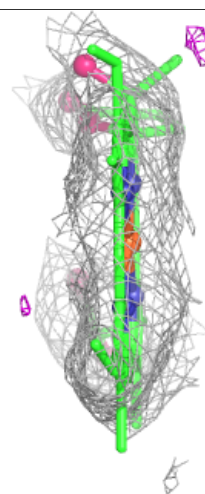
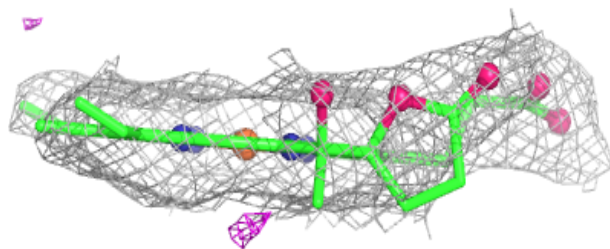
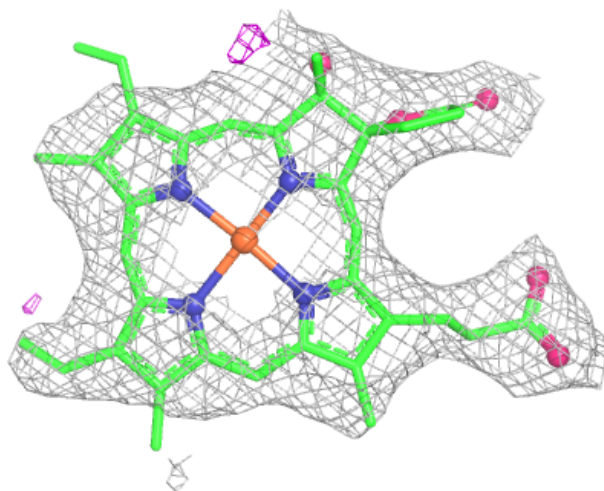
Electron density around HDD H 760:

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



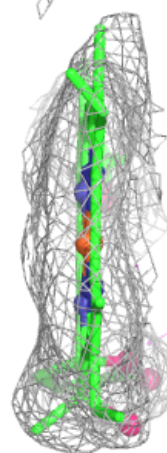
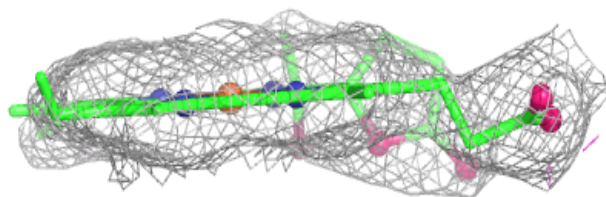
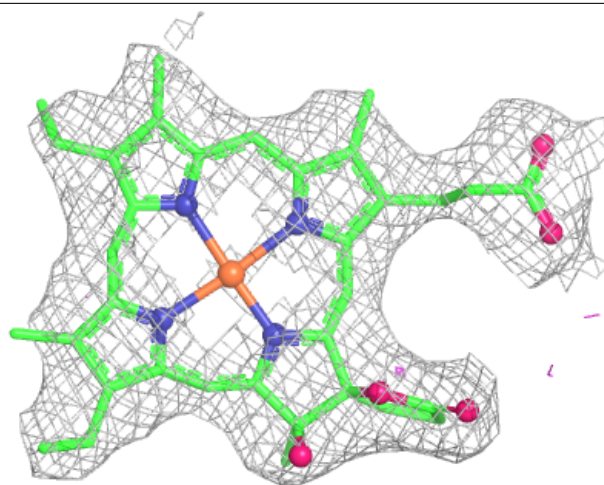
Electron density around HDD M 760:

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



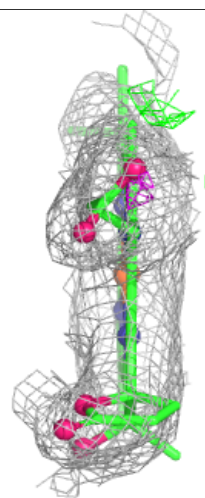
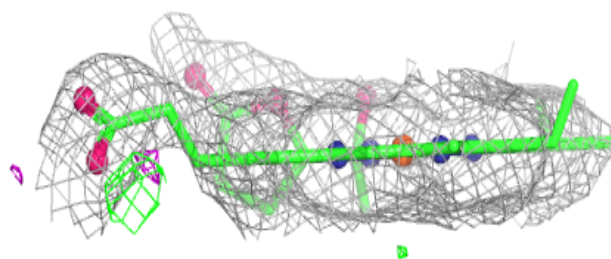
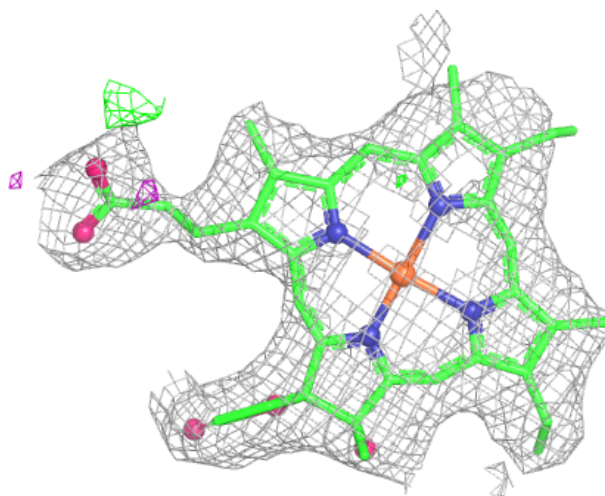
Electron density around HDD N 760:

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



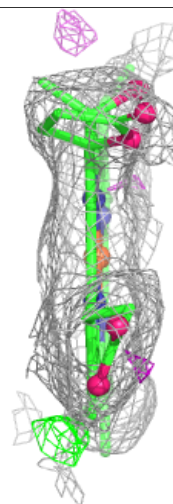
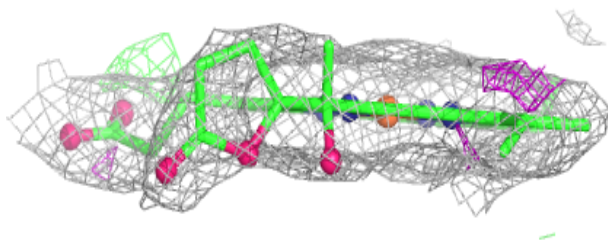
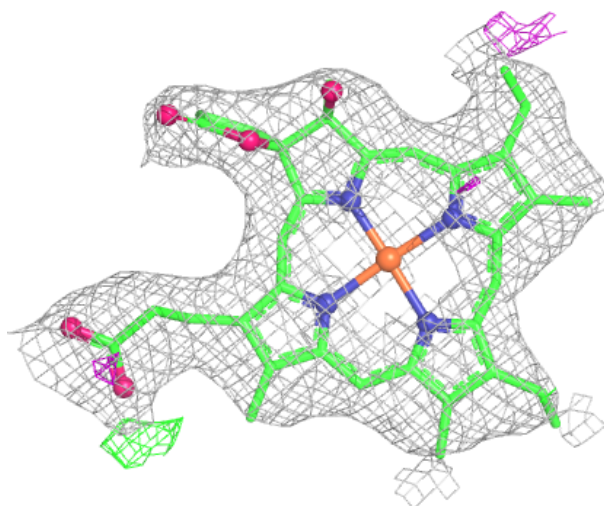
Electron density around HDD O 760:

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



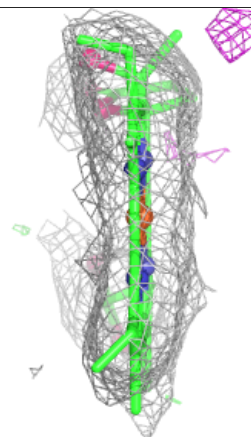
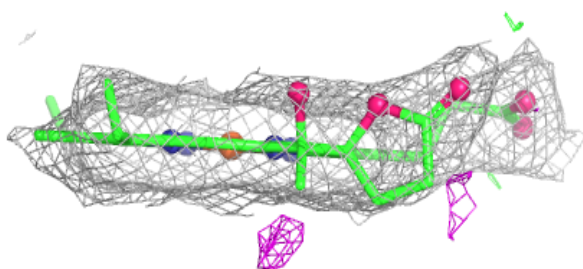
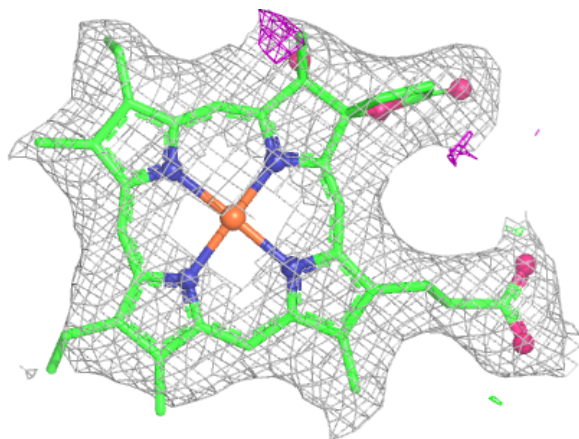
Electron density around HDD F 760:

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



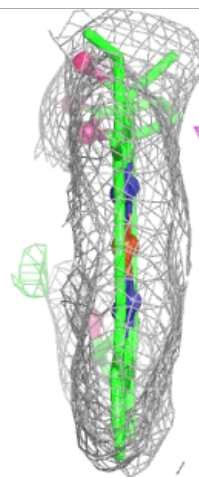
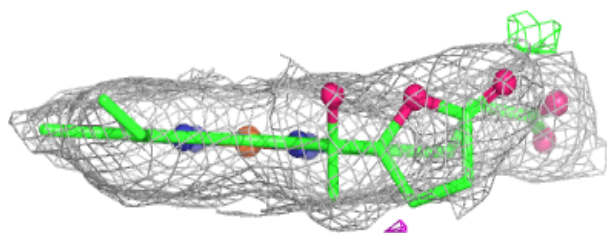
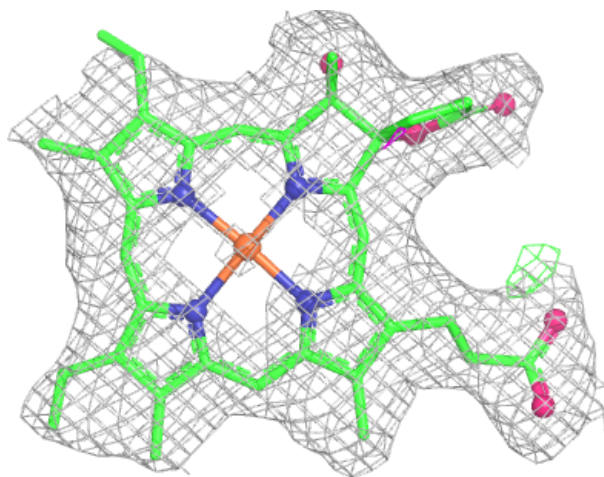
Electron density around HDD G 760:

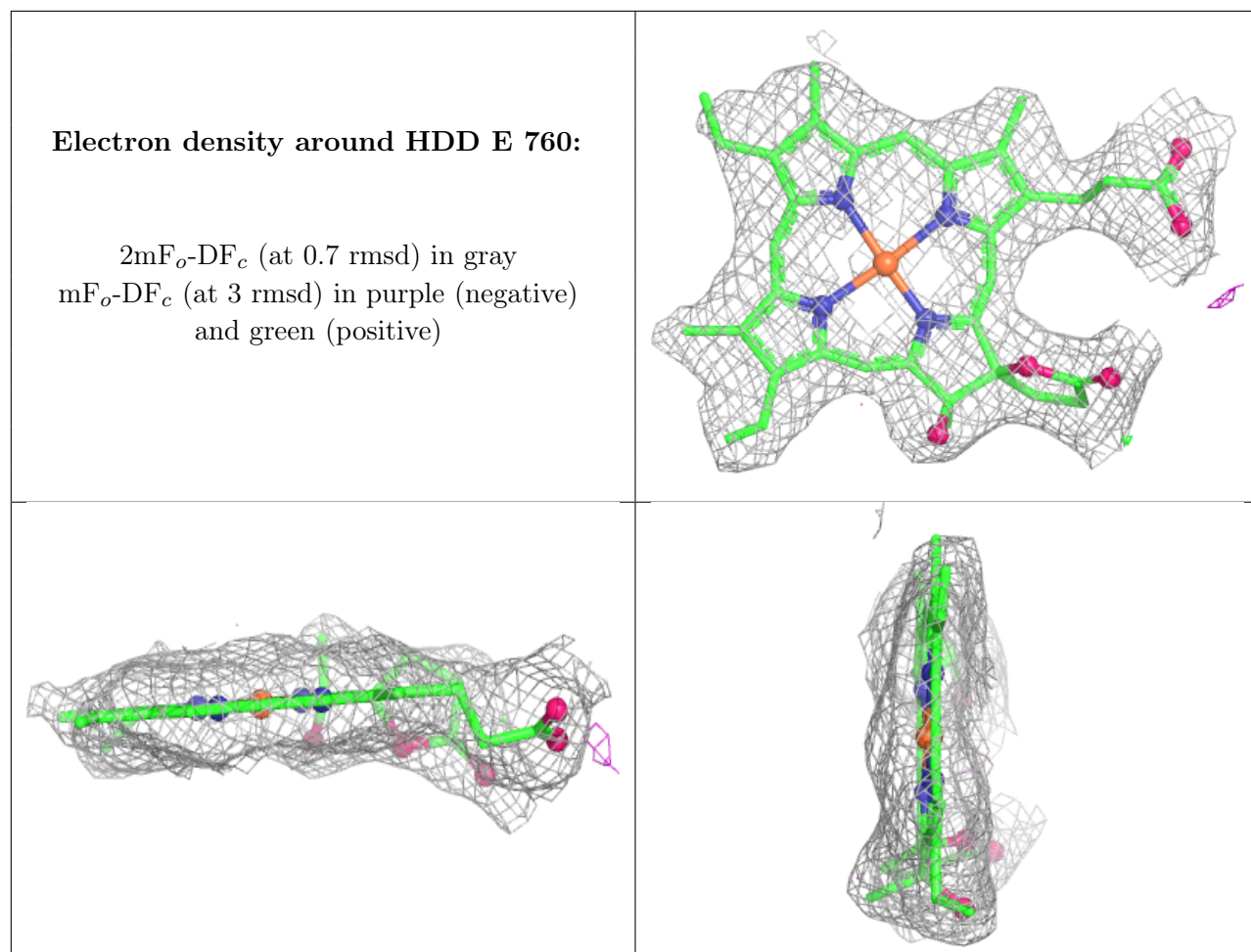
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HDD P 760:

$2mF_o-DF_c$ (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.