



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 11:32 AM UTC

PDB ID : 6RJN / pdb_00006rjn
Title : Crystal structure of a Fungal Catalase at 2.3 Angstroms
Authors : Gomez, S.; Navas-Yuste, S.; Payne, A.M.; Rivera, W.; Lopez-Esteva, M.;
Brangbour, C.; Fulla, D.; Juanhuix, J.; Fernandez, F.J.; Vega, M.C.
Deposited on : 2019-04-28
Resolution : 2.29 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

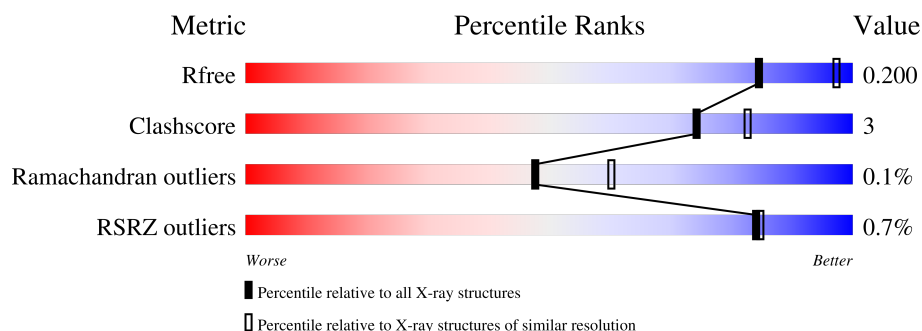
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	512	% 90% 8% .
1	B	512	91% 7% .
1	C	512	% 94% . .
1	D	512	% 91% 7% .

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	GOL	C	604	-	-	X	-
4	GOL	D	601	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 17948 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	502	Total	C	N	O	S	0	0	0
			4037	2566	705	754	12			
1	B	502	Total	C	N	O	S	0	0	0
			4037	2566	705	754	12			
1	C	502	Total	C	N	O	S	0	4	0
			4072	2586	716	758	12			
1	D	502	Total	C	N	O	S	0	5	0
			4068	2586	710	759	13			

There are 8 discrepancies between the modelled and reference sequences:

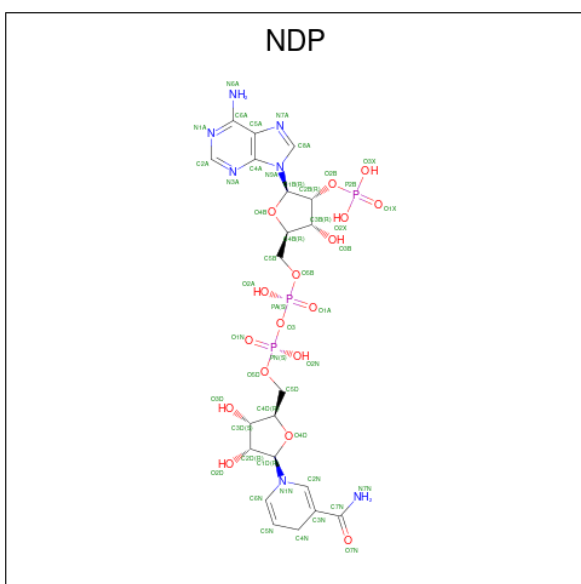
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP C1PHG1
A	0	ALA	-	expression tag	UNP C1PHG1
B	-1	GLY	-	expression tag	UNP C1PHG1
B	0	ALA	-	expression tag	UNP C1PHG1
C	-1	GLY	-	expression tag	UNP C1PHG1
C	0	ALA	-	expression tag	UNP C1PHG1
D	-1	GLY	-	expression tag	UNP C1PHG1
D	0	ALA	-	expression tag	UNP C1PHG1

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



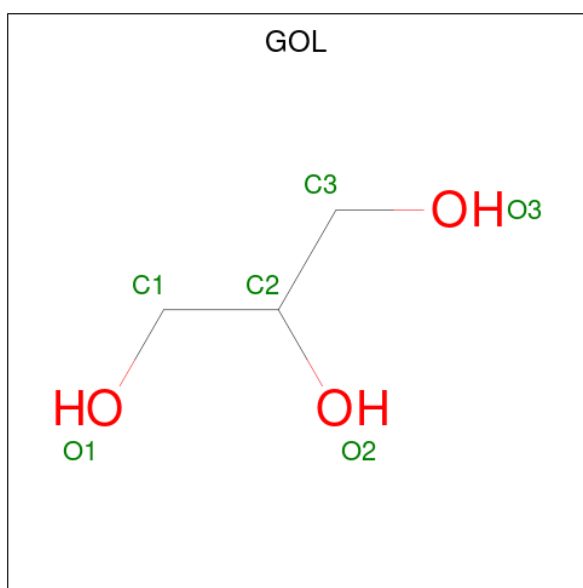
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $\text{C}_{21}\text{H}_{30}\text{N}_7\text{O}_{17}\text{P}_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



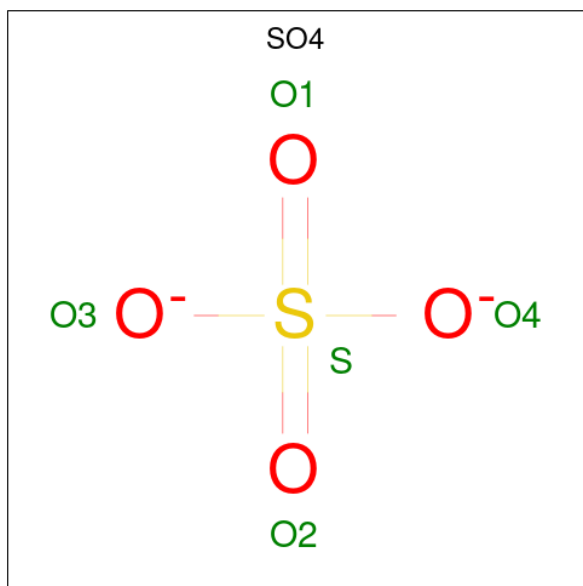
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total C O 6 3 3	0	0
4	C	1	Total C O 6 3 3	0	0
4	C	1	Total C O 6 3 3	0	0
4	C	1	Total C O 6 3 3	0	0
4	C	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0
4	D	1	Total C O 6 3 3	0	0

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



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

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	2	Total Cl 2 2	0	0
6	B	2	Total Cl 2 2	0	0
6	D	1	Total Cl 1 1	0	0

- Molecule 7 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total K 1 1	0	0
7	B	3	Total K 3 3	0	0
7	C	2	Total K 2 2	0	0
7	D	1	Total K 1 1	0	0

- Molecule 8 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total Na 1 1	0	0
8	B	1	Total Na 1 1	0	0
8	C	1	Total Na 1 1	0	0
8	D	1	Total Na 1 1	0	0

- Molecule 9 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			7	4	3		
9	C	1	Total	C	O	0	0
			7	4	3		

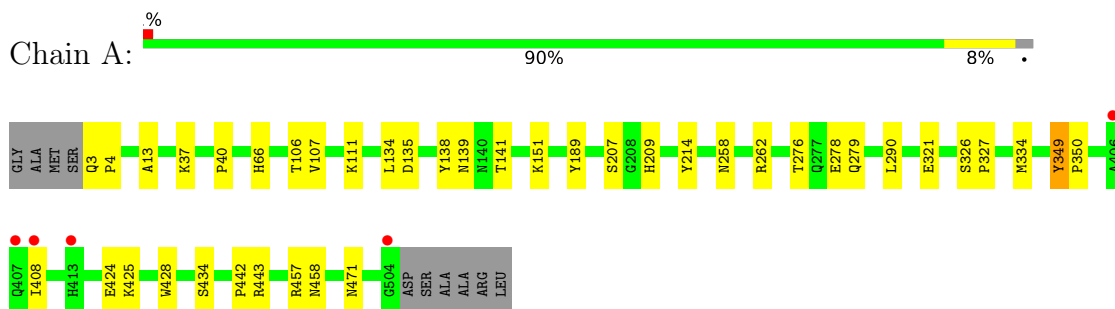
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	296	Total	O	0	0
			296	296		
10	B	321	Total	O	0	0
			321	321		
10	C	300	Total	O	0	0
			300	300		
10	D	311	Total	O	0	2
			311	311		

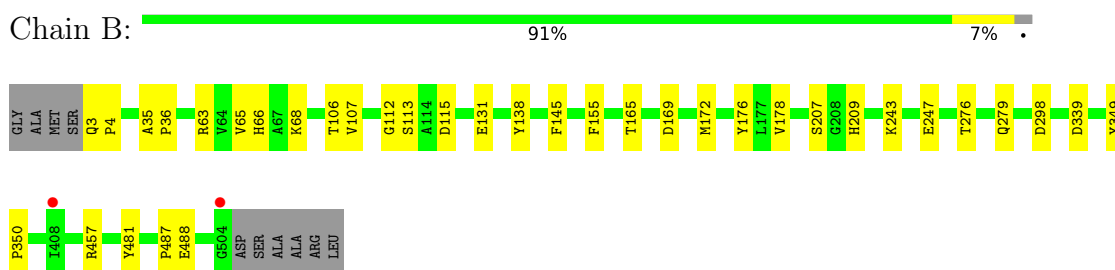
3 Residue-property plots [i](#)

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

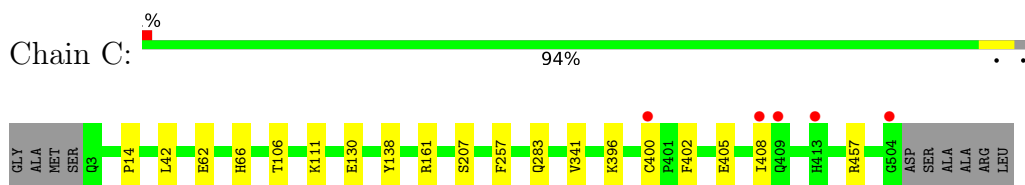
- Molecule 1: Catalase



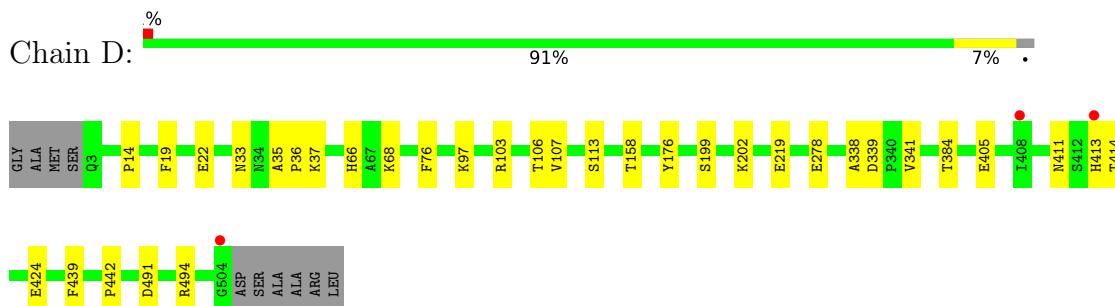
- Molecule 1: Catalase



- Molecule 1: Catalase



- Molecule 1: Catalase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	165.94Å 173.69Å 96.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.66 – 2.29 54.66 – 2.29	Depositor EDS
% Data completeness (in resolution range)	99.3 (54.66-2.29) 99.1 (54.66-2.29)	Depositor EDS
R_{merge}	0.25	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.142 , 0.196 0.153 , 0.200	Depositor DCC
R_{free} test set	6078 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.596	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.019 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17948	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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: NDP, PEG, GOL, CL, SO4, HEM, K, NA

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	0.60	3/4160 (0.1%)	0.69	0/5652
1	B	0.60	2/4160 (0.0%)	0.70	0/5652
1	C	0.59	0/4195	0.63	1/5696 (0.0%)
1	D	0.53	0/4197	0.62	1/5700 (0.0%)
All	All	0.58	5/16712 (0.0%)	0.66	2/22700 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	112	GLY	C-O	-5.94	1.20	1.24
1	A	442	PRO	C-O	-5.81	1.16	1.24
1	B	487	PRO	C-O	-5.67	1.17	1.24
1	A	349	TYR	C-O	-5.41	1.20	1.24
1	A	327	PRO	C-O	-5.02	1.18	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	161	ARG	N-CA-C	5.76	119.44	110.17
1	D	278	GLU	CB-CG-CD	5.26	121.55	112.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4037	0	3835	35	0
1	B	4037	0	3835	21	0
1	C	4072	0	3874	22	0
1	D	4068	0	3874	25	0
2	A	43	0	30	4	0
2	B	43	0	30	1	0
2	C	43	0	30	2	0
2	D	43	0	30	3	0
3	A	48	0	26	0	0
3	B	48	0	26	0	0
3	C	48	0	26	0	0
3	D	48	0	26	0	0
4	A	24	0	32	1	0
4	B	24	0	32	0	0
4	C	30	0	40	9	0
4	D	24	0	32	7	0
5	A	10	0	0	0	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
6	D	1	0	0	0	0
7	A	1	0	0	0	0
7	B	3	0	0	0	0
7	C	2	0	0	0	0
7	D	1	0	0	0	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
8	C	1	0	0	0	0
8	D	1	0	0	0	0
9	B	7	0	10	1	0
9	C	7	0	10	3	0
10	A	296	0	0	1	0
10	B	321	0	0	2	0
10	C	300	0	0	2	0
10	D	311	0	0	2	0
All	All	17948	0	15798	101	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (101) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:400:CYS:HB3	4:C:604:GOL:O2	1.39	1.22
1:C:400:CYS:HB3	4:C:604:GOL:C2	2.08	0.83
1:A:428:TRP:HE1	4:D:601:GOL:H32	1.52	0.73
1:C:457[B]:ARG:NE	4:C:605:GOL:O3	2.27	0.68
1:C:400:CYS:HB3	4:C:604:GOL:HO2	1.55	0.66
1:A:434:SER:O	1:A:443:ARG:NH2	2.27	0.62
1:A:334:MET:HE3	10:A:911:HOH:O	1.99	0.62
1:D:68[A]:LYS:NZ	1:D:113:SER:O	2.30	0.62
1:C:283:GLN:HE22	9:C:608:PEG:H22	1.64	0.61
1:C:400:CYS:SG	4:C:604:GOL:H2	2.40	0.60
1:C:457[B]:ARG:CD	4:C:605:GOL:O3	2.50	0.59
1:A:134:LEU:HD21	1:A:334:MET:HE1	1.86	0.57
1:B:66:HIS:HA	1:B:106:THR:O	2.06	0.56
1:C:138:TYR:CD1	1:C:207:SER:HA	2.40	0.56
2:D:604:HEM:HMB1	2:D:604:HEM:HBB2	1.88	0.56
1:D:36:PRO:HG3	4:D:601:GOL:H31	1.88	0.55
1:B:488:GLU:HG2	10:B:935:HOH:O	2.07	0.54
1:A:278:GLU:OE1	1:A:278:GLU:N	2.40	0.54
1:A:111:LYS:HE3	1:A:321:GLU:CD	2.33	0.54
1:D:33:ASN:OD1	4:D:601:GOL:H12	2.09	0.53
1:A:66:HIS:HA	1:A:106:THR:O	2.09	0.52
2:C:601:HEM:HMB2	2:C:601:HEM:HBB2	1.90	0.52
1:C:62:GLU:OE1	1:C:111[B]:LYS:HG2	2.09	0.52
1:D:66:HIS:HA	1:D:106:THR:O	2.10	0.51
1:A:66:HIS:CE1	1:A:107:VAL:HG22	2.45	0.51
1:A:428:TRP:NE1	4:D:601:GOL:H32	2.23	0.51
1:D:36:PRO:HG3	4:D:601:GOL:C3	2.41	0.51
1:B:138:TYR:CD1	1:B:207:SER:HA	2.46	0.50
2:A:601:HEM:HBB2	2:A:601:HEM:HMB2	1.94	0.50
1:A:40:PRO:HB2	1:C:42:LEU:HD22	1.95	0.49
1:C:405:GLU:HG3	1:D:14:PRO:HB2	1.94	0.49
1:A:471:ASN:OD1	1:A:471:ASN:C	2.55	0.49
1:A:138:TYR:CD1	1:A:207:SER:HA	2.47	0.49
1:A:457:ARG:NH1	1:A:458:ASN:OD1	2.46	0.49
1:B:66:HIS:CE1	1:B:107:VAL:HG22	2.48	0.49
1:A:425:LYS:HG2	1:D:22:GLU:HG2	1.95	0.48
1:A:214:TYR:OH	4:A:606:GOL:O2	2.23	0.48
1:D:199:SER:HB3	1:D:202:LYS:HG3	1.97	0.47
1:D:341:VAL:HG13	2:D:604:HEM:HAB	1.97	0.47
1:B:243:LYS:NZ	1:B:247:GLU:OE2	2.47	0.47
1:A:424:GLU:OE1	1:D:37:LYS:HD3	2.14	0.47
1:B:349:TYR:HB2	1:B:350:PRO:HD3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:601:HEM:HBB2	2:A:601:HEM:CMB	2.46	0.46
1:C:66:HIS:HA	1:C:106:THR:O	2.15	0.46
1:D:76:PHE:O	1:D:97:LYS:HA	2.16	0.46
1:A:3:GLN:N	1:A:4:PRO:CD	2.79	0.45
1:C:138:TYR:CE1	1:C:207:SER:HA	2.52	0.45
1:A:209:HIS:CE1	1:A:290:LEU:HD22	2.52	0.45
2:B:602:HEM:HMB2	2:B:602:HEM:HBB2	1.99	0.45
1:A:151:LYS:HE2	1:D:19:PHE:CD2	2.52	0.45
1:A:428:TRP:HE1	4:D:601:GOL:C3	2.25	0.45
1:B:178:VAL:HA	1:B:481:TYR:CD1	2.52	0.45
1:D:491:ASP:OD1	1:D:494:ARG:NH2	2.43	0.45
1:A:276:THR:OG1	1:A:279:GLN:HG3	2.17	0.44
1:A:139:ASN:CG	2:A:601:HEM:HAC	2.43	0.44
1:A:13:ALA:HA	1:B:131:GLU:O	2.18	0.44
1:A:138:TYR:CE1	1:A:207:SER:HA	2.53	0.43
1:C:283:GLN:NE2	9:C:608:PEG:H22	2.32	0.43
1:C:396:LYS:HE2	1:C:402:PHE:HZ	1.83	0.43
1:B:145:PHE:CD1	1:B:155:PHE:CZ	3.06	0.43
1:C:457[B]:ARG:HD2	4:C:605:GOL:O3	2.18	0.43
1:A:37:LYS:HD3	1:D:424:GLU:OE1	2.18	0.43
1:C:400:CYS:HB3	4:C:604:GOL:H2	1.96	0.43
1:C:14:PRO:HB2	1:D:405:GLU:HG3	2.01	0.42
1:A:408:ILE:O	1:A:408:ILE:HG22	2.19	0.42
1:D:411:ASN:OD1	1:D:413:HIS:HB3	2.18	0.42
1:D:219:GLU:HA	1:D:414:THR:OG1	2.20	0.42
1:D:103:ARG:HD3	2:D:604:HEM:O1D	2.19	0.42
1:B:165:THR:HG22	1:C:257:PHE:CZ	2.54	0.42
1:A:3:GLN:N	1:A:4:PRO:HD2	2.34	0.42
1:D:384:THR:HG23	10:D:816:HOH:O	2.20	0.42
1:B:35:ALA:N	1:B:36:PRO:CD	2.82	0.42
1:C:283:GLN:HE22	9:C:608:PEG:H31	1.84	0.42
1:B:68:LYS:HE3	1:B:115:ASP:OD1	2.19	0.42
1:A:134:LEU:CD2	1:A:334:MET:HE1	2.50	0.41
1:A:139:ASN:ND2	2:A:601:HEM:HAC	2.35	0.41
1:A:141:THR:HB	1:A:189:TYR:CE1	2.55	0.41
1:B:209:HIS:HD1	1:B:339:ASP:CG	2.28	0.41
1:D:176:TYR:CD2	1:D:176:TYR:C	2.98	0.41
1:B:63:ARG:HB3	1:B:65:VAL:O	2.20	0.41
1:D:158:THR:HG21	1:D:176:TYR:HB2	2.02	0.41
4:D:603:GOL:O1	10:D:701:HOH:O	2.22	0.41
1:B:68:LYS:HE2	1:B:113:SER:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:603:GOL:H31	10:C:882:HOH:O	2.20	0.41
1:A:135:ASP:HB2	1:A:326:SER:O	2.21	0.41
1:A:278:GLU:H	1:A:278:GLU:CD	2.28	0.41
1:A:349:TYR:HB2	1:A:350:PRO:HD3	2.02	0.41
1:B:298:ASP:HB3	9:B:608:PEG:H21	2.03	0.41
1:B:457:ARG:NH1	10:B:714:HOH:O	2.50	0.41
1:B:276:THR:OG1	1:B:279:GLN:HG3	2.21	0.41
1:D:439:PHE:C	1:D:442:PRO:HD2	2.46	0.41
1:D:66:HIS:CE1	1:D:107:VAL:HG22	2.56	0.40
1:B:169:ASP:HB3	1:B:172:MET:HG2	2.03	0.40
1:B:176:TYR:CD2	1:B:176:TYR:C	2.99	0.40
1:C:130:GLU:HG2	10:C:876:HOH:O	2.20	0.40
1:D:35:ALA:N	1:D:36:PRO:CD	2.84	0.40
1:A:258:ASN:O	1:A:262:ARG:HG3	2.22	0.40
1:B:3:GLN:N	1:B:4:PRO:CD	2.85	0.40
1:D:338:ALA:O	1:D:339:ASP:C	2.63	0.40
1:A:111:LYS:HE3	1:A:321:GLU:OE1	2.21	0.40
1:C:341:VAL:HG13	2:C:601:HEM:HAB	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	500/512 (98%)	487 (97%)	13 (3%)	0	100	100
1	B	500/512 (98%)	486 (97%)	14 (3%)	0	100	100
1	C	504/512 (98%)	490 (97%)	13 (3%)	1 (0%)	43	55
1	D	505/512 (99%)	491 (97%)	14 (3%)	0	100	100
All	All	2009/2048 (98%)	1954 (97%)	54 (3%)	1 (0%)	48	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	408	ILE

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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](#)

Of 45 ligands modelled in this entry, 16 are monoatomic - leaving 29 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	HEM	A	601	1	50,50,50	1.94	14 (28%)	67,82,82	1.47	11 (16%)
2	HEM	D	604	1	50,50,50	1.70	9 (18%)	67,82,82	1.58	14 (20%)
3	NDP	C	602	-	51,52,52	2.69	14 (27%)	71,80,80	1.51	13 (18%)
4	GOL	C	605	-	5,5,5	0.51	0	5,5,5	0.64	0
4	GOL	D	606	-	5,5,5	0.12	0	5,5,5	0.42	0
4	GOL	B	606	-	5,5,5	0.41	0	5,5,5	0.35	0
2	HEM	C	601	1	50,50,50	1.61	6 (12%)	67,82,82	1.57	16 (23%)
9	PEG	B	608	-	6,6,6	0.58	0	5,5,5	0.64	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	605	-	5,5,5	0.34	0	5,5,5	0.39	0
4	GOL	A	606	-	5,5,5	0.41	0	5,5,5	0.93	0
2	HEM	B	602	1	50,50,50	1.47	8 (16%)	67,82,82	1.37	10 (14%)
3	NDP	A	602	-	51,52,52	2.60	11 (21%)	71,80,80	1.46	10 (14%)
4	GOL	C	606	-	5,5,5	0.70	0	5,5,5	0.44	0
4	GOL	D	601	-	5,5,5	0.47	0	5,5,5	0.94	0
3	NDP	B	603	-	51,52,52	2.52	11 (21%)	71,80,80	1.49	12 (16%)
9	PEG	C	608	-	6,6,6	0.53	0	5,5,5	1.04	1 (20%)
4	GOL	D	602	-	5,5,5	0.20	0	5,5,5	0.62	0
4	GOL	C	603	-	5,5,5	0.28	0	5,5,5	0.37	0
4	GOL	A	604	-	5,5,5	0.30	0	5,5,5	0.39	0
4	GOL	D	603	-	5,5,5	0.28	0	5,5,5	0.71	0
5	SO4	A	608	-	4,4,4	0.40	0	6,6,6	0.12	0
4	GOL	B	607	-	5,5,5	0.66	0	5,5,5	0.70	0
5	SO4	A	607	-	4,4,4	0.39	0	6,6,6	0.09	0
3	NDP	D	605	-	51,52,52	2.64	11 (21%)	71,80,80	1.46	10 (14%)
4	GOL	A	603	-	5,5,5	0.34	0	5,5,5	0.48	0
4	GOL	B	604	-	5,5,5	0.27	0	5,5,5	0.72	0
4	GOL	A	605	-	5,5,5	0.44	0	5,5,5	0.46	0
4	GOL	C	604	-	5,5,5	0.37	0	5,5,5	0.53	0
4	GOL	C	607	-	5,5,5	0.44	0	5,5,5	0.60	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	HEM	A	601	1	-	4/14/54/54	-
2	HEM	D	604	1	-	2/14/54/54	-
3	NDP	C	602	-	-	4/34/77/77	0/5/5/5
4	GOL	C	605	-	-	2/4/4/4	-
4	GOL	D	606	-	-	2/4/4/4	-
4	GOL	B	606	-	-	2/4/4/4	-
2	HEM	C	601	1	-	2/14/54/54	-
9	PEG	B	608	-	-	4/4/4/4	-
4	GOL	B	605	-	-	1/4/4/4	-
4	GOL	A	606	-	-	2/4/4/4	-
2	HEM	B	602	1	-	2/14/54/54	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NDP	A	602	-	-	2/34/77/77	0/5/5/5
4	GOL	C	606	-	-	4/4/4/4	-
4	GOL	D	601	-	-	4/4/4/4	-
3	NDP	B	603	-	-	3/34/77/77	0/5/5/5
9	PEG	C	608	-	-	3/4/4/4	-
4	GOL	D	602	-	-	2/4/4/4	-
4	GOL	C	603	-	-	2/4/4/4	-
4	GOL	A	604	-	-	4/4/4/4	-
4	GOL	D	603	-	-	4/4/4/4	-
4	GOL	B	607	-	-	2/4/4/4	-
3	NDP	D	605	-	-	2/34/77/77	0/5/5/5
4	GOL	A	603	-	-	4/4/4/4	-
4	GOL	B	604	-	-	2/4/4/4	-
4	GOL	A	605	-	-	2/4/4/4	-
4	GOL	C	604	-	-	1/4/4/4	-
4	GOL	C	607	-	-	4/4/4/4	-

All (84) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	603	NDP	P2B-O2B	11.92	1.80	1.59
3	C	602	NDP	P2B-O2B	11.83	1.80	1.59
3	A	602	NDP	P2B-O2B	11.47	1.79	1.59
3	D	605	NDP	P2B-O2B	11.43	1.79	1.59
3	D	605	NDP	PA-O3	10.25	1.70	1.59
3	C	602	NDP	PA-O3	9.46	1.69	1.59
3	A	602	NDP	PA-O3	7.83	1.68	1.59
3	B	603	NDP	PA-O3	7.02	1.67	1.59
2	D	604	HEM	FE-NB	6.55	2.15	1.94
2	C	601	HEM	FE-NC	6.42	2.16	1.95
3	A	602	NDP	PN-O3	6.28	1.66	1.59
3	C	602	NDP	PN-O3	5.41	1.65	1.59
2	A	601	HEM	FE-NB	5.28	2.11	1.94
3	D	605	NDP	PN-O3	4.61	1.64	1.59
2	C	601	HEM	FE-ND	4.59	2.09	1.94
2	A	601	HEM	C1B-NB	-4.47	1.32	1.40
2	D	604	HEM	FE-NC	4.44	2.09	1.95
2	B	602	HEM	FE-NC	4.30	2.09	1.95
2	A	601	HEM	C4D-ND	-4.10	1.33	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	603	NDP	PN-O5D	4.05	1.75	1.59
3	A	602	NDP	PN-O5D	4.00	1.75	1.59
2	A	601	HEM	C1C-C2C	-3.93	1.37	1.45
3	C	602	NDP	C2B-C1B	3.76	1.62	1.53
3	B	603	NDP	C2A-N1A	3.73	1.40	1.33
2	B	602	HEM	FE-NA	3.70	2.07	1.95
2	A	601	HEM	FE-NC	3.54	2.06	1.95
3	A	602	NDP	C2B-C1B	3.50	1.61	1.53
2	A	601	HEM	C3C-C4C	-3.48	1.39	1.46
3	C	602	NDP	PN-O5D	3.42	1.72	1.59
2	A	601	HEM	C4B-NB	-3.40	1.32	1.38
3	B	603	NDP	O2B-C2B	-3.39	1.32	1.44
2	A	601	HEM	C1D-ND	-3.39	1.32	1.38
2	D	604	HEM	FE-ND	3.34	2.05	1.94
3	A	602	NDP	C2A-N1A	3.31	1.39	1.33
3	D	605	NDP	C2B-C1B	3.31	1.61	1.53
2	C	601	HEM	FE-NB	3.30	2.05	1.94
3	D	605	NDP	PN-O5D	3.24	1.72	1.59
2	B	602	HEM	FE-NB	3.05	2.04	1.94
2	D	604	HEM	CAC-C3C	3.04	1.55	1.47
2	A	601	HEM	C1C-NC	-3.03	1.33	1.39
2	D	604	HEM	CMB-C2B	3.02	1.57	1.50
2	D	604	HEM	FE-NA	2.98	2.05	1.95
2	A	601	HEM	O2A-CGA	-2.93	1.21	1.30
3	D	605	NDP	O2B-C2B	-2.87	1.34	1.44
3	A	602	NDP	O2B-C2B	-2.82	1.34	1.44
3	B	603	NDP	O4B-C1B	2.75	1.48	1.42
2	C	601	HEM	CAC-C3C	2.69	1.54	1.47
3	D	605	NDP	O4B-C4B	-2.69	1.39	1.45
3	B	603	NDP	C4A-N3A	2.69	1.39	1.34
2	B	602	HEM	FE-ND	2.68	2.03	1.94
3	D	605	NDP	C5A-C4A	2.65	1.43	1.39
3	C	602	NDP	O2B-C2B	-2.64	1.35	1.44
3	A	602	NDP	C8A-N9A	2.63	1.42	1.37
3	B	603	NDP	O4B-C4B	-2.63	1.39	1.45
3	D	605	NDP	C2A-N1A	2.60	1.38	1.33
3	C	602	NDP	C4A-N3A	2.60	1.39	1.34
3	C	602	NDP	C5A-C4A	2.60	1.43	1.39
3	A	602	NDP	C6A-N6A	2.60	1.40	1.34
3	A	602	NDP	C5A-C4A	2.56	1.43	1.39
2	D	604	HEM	CAB-C3B	2.55	1.54	1.47
3	C	602	NDP	C2A-N1A	2.46	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	NDP	C7N-N7N	2.40	1.40	1.33
3	B	603	NDP	C5A-C4A	2.38	1.43	1.39
3	D	605	NDP	C6A-N6A	2.38	1.40	1.34
3	D	605	NDP	C7N-N7N	2.37	1.40	1.33
2	A	601	HEM	C1A-C2A	-2.36	1.39	1.44
3	C	602	NDP	C7N-N7N	2.34	1.40	1.33
2	C	601	HEM	CMD-C2D	2.34	1.55	1.50
2	B	602	HEM	CAB-C3B	2.30	1.53	1.47
2	D	604	HEM	CMC-C2C	2.29	1.55	1.50
2	C	601	HEM	CAB-C3B	2.27	1.53	1.47
2	A	601	HEM	O2D-CGD	-2.25	1.23	1.30
3	C	602	NDP	PA-O5B	2.25	1.68	1.59
2	B	602	HEM	C2A-C3A	-2.19	1.33	1.38
3	C	602	NDP	C8A-N9A	2.18	1.41	1.37
3	C	602	NDP	O3B-C3B	-2.17	1.37	1.43
2	B	602	HEM	CAC-C3C	2.16	1.53	1.47
2	B	602	HEM	CMC-C2C	2.14	1.55	1.50
3	B	603	NDP	C3B-C2B	2.13	1.57	1.53
2	A	601	HEM	C2A-C3A	-2.05	1.33	1.38
2	D	604	HEM	C2A-C3A	-2.04	1.33	1.38
3	C	602	NDP	C8A-N7A	2.04	1.35	1.31
2	A	601	HEM	C4C-NC	-2.03	1.35	1.39
3	B	603	NDP	C3D-C4D	2.02	1.58	1.53

All (97) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	602	NDP	O3-PA-O1A	-5.05	95.52	110.70
2	C	601	HEM	CAD-CBD-CGD	-4.52	101.67	113.67
3	D	605	NDP	O3-PA-O1A	-4.43	97.39	110.70
2	A	601	HEM	CHC-C4B-NB	4.27	129.02	124.42
3	B	603	NDP	O2B-P2B-O1X	-4.27	94.12	109.33
3	B	603	NDP	O3-PA-O1A	-4.24	97.95	110.70
3	A	602	NDP	O3-PA-O1A	-4.21	98.05	110.70
2	D	604	HEM	C3B-C4B-NB	-4.14	106.49	109.47
3	A	602	NDP	O2B-P2B-O1X	-3.88	95.50	109.33
3	D	605	NDP	O2B-P2B-O1X	-3.48	96.93	109.33
2	D	604	HEM	C2D-C1D-ND	-3.35	106.04	109.90
2	D	604	HEM	C1B-NB-C4B	3.28	109.09	105.21
2	C	601	HEM	C3B-C4B-NB	-3.26	107.12	109.47
3	D	605	NDP	O2N-PN-O3	3.26	116.09	107.27
2	D	604	HEM	C2A-C1A-NA	-3.19	106.62	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	602	NDP	O2B-P2B-O1X	-3.17	98.03	109.33
2	D	604	HEM	C4D-ND-C1D	3.14	108.93	105.21
2	A	601	HEM	C1B-NB-C4B	3.12	108.91	105.21
2	B	602	HEM	CAD-CBD-CGD	-3.11	105.41	113.67
2	B	602	HEM	C4D-ND-C1D	3.01	108.77	105.21
2	A	601	HEM	CHD-C1D-ND	2.98	127.63	124.42
2	D	604	HEM	CHB-C4A-NA	2.97	129.25	123.86
2	B	602	HEM	C3D-C4D-ND	-2.95	106.94	110.17
2	C	601	HEM	C1B-NB-C4B	2.94	108.69	105.21
2	C	601	HEM	CMA-C3A-C4A	-2.93	120.96	125.42
3	C	602	NDP	C2A-N1A-C6A	-2.91	113.95	118.73
2	D	604	HEM	C3A-C4A-NA	-2.85	105.57	110.14
3	D	605	NDP	O5D-PN-O1N	-2.78	97.93	108.94
3	C	602	NDP	O5D-PN-O1N	-2.71	98.19	108.94
3	A	602	NDP	O3X-P2B-O2X	2.67	117.82	107.80
3	B	603	NDP	O5D-PN-O1N	-2.66	98.41	108.94
3	C	602	NDP	P2B-O2B-C2B	-2.63	116.41	123.43
2	A	601	HEM	O2A-CGA-O1A	-2.63	116.57	123.33
2	C	601	HEM	C4D-ND-C1D	2.61	108.30	105.21
2	C	601	HEM	C4A-C3A-C2A	2.59	109.79	106.82
2	D	604	HEM	C1D-C2D-C3D	2.58	109.69	106.98
3	C	602	NDP	O2N-PN-O1N	2.57	124.41	112.44
2	D	604	HEM	C4A-NA-C1A	2.57	110.01	105.82
3	C	602	NDP	PN-O5D-C5D	-2.57	106.64	121.35
2	C	601	HEM	CMB-C2B-C1B	-2.56	121.03	125.03
3	A	602	NDP	PA-O5B-C5B	-2.54	106.79	121.35
2	C	601	HEM	C2A-C1A-NA	-2.54	107.34	110.15
3	D	605	NDP	O3X-P2B-O2X	2.52	117.27	107.80
3	C	602	NDP	PA-O5B-C5B	-2.52	106.92	121.35
3	A	602	NDP	P2B-O2B-C2B	-2.51	116.72	123.43
2	A	601	HEM	CHD-C4C-NC	2.51	127.19	124.45
2	A	601	HEM	C4C-NC-C1C	2.51	109.91	105.82
2	B	602	HEM	C2A-C1A-NA	-2.51	107.37	110.15
2	A	601	HEM	CHA-C4D-ND	2.50	127.46	124.37
3	B	603	NDP	O3X-P2B-O2X	2.47	117.07	107.80
3	A	602	NDP	C2A-N1A-C6A	-2.47	114.68	118.73
3	A	602	NDP	O7N-C7N-N7N	-2.45	117.39	122.89
3	D	605	NDP	PA-O5B-C5B	-2.45	107.29	121.35
3	D	605	NDP	P2B-O2B-C2B	-2.45	116.89	123.43
2	D	604	HEM	C4A-C3A-C2A	2.45	109.62	106.82
3	A	602	NDP	O3-PN-O1N	2.44	118.05	110.70
3	B	603	NDP	P2B-O2B-C2B	-2.44	116.92	123.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	605	NDP	C2A-N1A-C6A	-2.44	114.72	118.73
2	A	601	HEM	CHB-C1B-NB	2.44	127.38	124.37
2	B	602	HEM	CMB-C2B-C1B	-2.42	121.25	125.03
3	C	602	NDP	O3X-P2B-O2X	2.39	116.78	107.80
2	D	604	HEM	CBA-CAA-C2A	-2.38	105.95	112.53
2	C	601	HEM	O2D-CGD-CBD	2.37	121.49	114.00
2	D	604	HEM	C4C-NC-C1C	2.31	109.58	105.82
2	D	604	HEM	C4B-C3B-C2B	2.30	109.39	107.28
2	C	601	HEM	C3D-C4D-ND	-2.30	107.65	110.17
3	C	602	NDP	N3A-C4A-N9A	2.29	131.07	127.17
3	B	603	NDP	PA-O5B-C5B	-2.27	108.36	121.35
2	B	602	HEM	C4C-C3C-C2C	2.23	108.75	106.81
2	C	601	HEM	CHB-C1B-NB	2.23	127.12	124.37
3	D	605	NDP	PN-O5D-C5D	-2.21	108.70	121.35
2	C	601	HEM	C4A-NA-C1A	2.19	109.40	105.82
2	B	602	HEM	O1D-CGD-CBD	-2.19	116.14	123.09
2	A	601	HEM	O2D-CGD-CBD	2.19	120.93	114.00
2	A	601	HEM	C4D-ND-C1D	2.17	107.78	105.21
2	B	602	HEM	C4C-NC-C1C	2.17	109.36	105.82
3	B	603	NDP	N3A-C4A-N9A	2.17	130.85	127.17
3	D	605	NDP	O7N-C7N-N7N	-2.15	118.07	122.89
3	B	603	NDP	C3N-C2N-N1N	-2.14	120.06	123.20
9	C	608	PEG	C3-O2-C2	2.13	122.56	113.26
3	C	602	NDP	C5A-C4A-N3A	-2.12	123.80	126.72
2	B	602	HEM	CHD-C4C-C3C	2.11	128.77	125.21
3	C	602	NDP	O2A-PA-O1A	2.10	122.22	112.44
2	C	601	HEM	O1D-CGD-CBD	-2.10	116.43	123.09
3	B	603	NDP	C2A-N1A-C6A	-2.10	115.28	118.73
3	B	603	NDP	O3-PN-O1N	2.09	116.99	110.70
2	A	601	HEM	CHA-C1A-NA	2.08	127.64	123.86
3	B	603	NDP	O2N-PN-O5D	-2.07	98.17	107.57
2	C	601	HEM	C3B-C2B-C1B	2.07	107.96	106.41
2	C	601	HEM	C3A-C4A-NA	-2.07	106.83	110.14
3	A	602	NDP	O5D-PN-O1N	-2.05	100.81	108.94
3	B	603	NDP	O2N-PN-O1N	2.04	121.95	112.44
2	D	604	HEM	CAD-CBD-CGD	-2.03	108.27	113.67
2	C	601	HEM	C4C-NC-C1C	2.02	109.11	105.82
2	B	602	HEM	O2D-CGD-CBD	2.01	120.34	114.00
3	A	602	NDP	N3A-C2A-N1A	2.01	131.62	128.58
3	C	602	NDP	C5A-C6A-N1A	2.00	122.59	117.51

There are no chirality outliers.

All (72) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	603	NDP	C5D-O5D-PN-O3
4	A	603	GOL	O1-C1-C2-C3
4	A	604	GOL	O1-C1-C2-C3
4	A	604	GOL	C1-C2-C3-O3
4	A	605	GOL	O1-C1-C2-C3
4	A	606	GOL	C1-C2-C3-O3
4	B	604	GOL	O1-C1-C2-C3
4	B	607	GOL	O1-C1-C2-O2
4	B	607	GOL	O1-C1-C2-C3
4	C	606	GOL	O1-C1-C2-C3
4	C	607	GOL	O1-C1-C2-O2
4	C	607	GOL	O1-C1-C2-C3
4	C	607	GOL	C1-C2-C3-O3
4	D	601	GOL	O1-C1-C2-C3
4	D	602	GOL	O1-C1-C2-C3
4	D	603	GOL	O1-C1-C2-C3
4	D	603	GOL	C1-C2-C3-O3
4	D	606	GOL	O1-C1-C2-C3
4	B	604	GOL	O1-C1-C2-O2
4	B	606	GOL	O2-C2-C3-O3
4	C	603	GOL	O2-C2-C3-O3
4	D	601	GOL	O1-C1-C2-O2
9	B	608	PEG	O1-C1-C2-O2
9	B	608	PEG	O2-C3-C4-O4
4	A	603	GOL	C1-C2-C3-O3
4	B	605	GOL	O1-C1-C2-C3
4	B	606	GOL	C1-C2-C3-O3
4	C	603	GOL	C1-C2-C3-O3
4	C	605	GOL	C1-C2-C3-O3
4	C	606	GOL	C1-C2-C3-O3
4	D	601	GOL	C1-C2-C3-O3
4	A	603	GOL	O1-C1-C2-O2
4	A	604	GOL	O2-C2-C3-O3
4	A	605	GOL	O1-C1-C2-O2
4	A	606	GOL	O2-C2-C3-O3
4	C	606	GOL	O2-C2-C3-O3
4	D	602	GOL	O1-C1-C2-O2
4	D	603	GOL	O1-C1-C2-O2
4	D	606	GOL	O1-C1-C2-O2
9	C	608	PEG	O1-C1-C2-O2
4	A	604	GOL	O1-C1-C2-O2
4	C	606	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
4	C	607	GOL	O2-C2-C3-O3
4	D	601	GOL	O2-C2-C3-O3
3	B	603	NDP	O4D-C1D-N1N-C6N
3	D	605	NDP	O4D-C1D-N1N-C6N
4	A	603	GOL	O2-C2-C3-O3
3	C	602	NDP	O4D-C1D-N1N-C6N
2	A	601	HEM	C2C-C3C-CAC-CBC
9	C	608	PEG	C4-C3-O2-C2
3	A	602	NDP	O4D-C1D-N1N-C6N
3	B	603	NDP	PA-O3-PN-O1N
3	C	602	NDP	PA-O3-PN-O1N
9	B	608	PEG	C4-C3-O2-C2
4	D	603	GOL	O2-C2-C3-O3
4	C	604	GOL	C1-C2-C3-O3
9	C	608	PEG	C1-C2-O2-C3
9	B	608	PEG	C1-C2-O2-C3
2	D	604	HEM	CAD-CBD-CGD-O2D
2	D	604	HEM	CAD-CBD-CGD-O1D
4	C	605	GOL	O2-C2-C3-O3
3	C	602	NDP	O4D-C4D-C5D-O5D
2	A	601	HEM	CAD-CBD-CGD-O2D
2	B	602	HEM	CAD-CBD-CGD-O2D
2	A	601	HEM	CAD-CBD-CGD-O1D
2	B	602	HEM	CAD-CBD-CGD-O1D
3	C	602	NDP	O4B-C4B-C5B-O5B
2	C	601	HEM	CAD-CBD-CGD-O2D
2	A	601	HEM	C4C-C3C-CAC-CBC
2	C	601	HEM	CAD-CBD-CGD-O1D
3	D	605	NDP	PA-O3-PN-O2N
3	A	602	NDP	PA-O3-PN-O2N

There are no ring outliers.

12 monomers are involved in 31 short contacts:

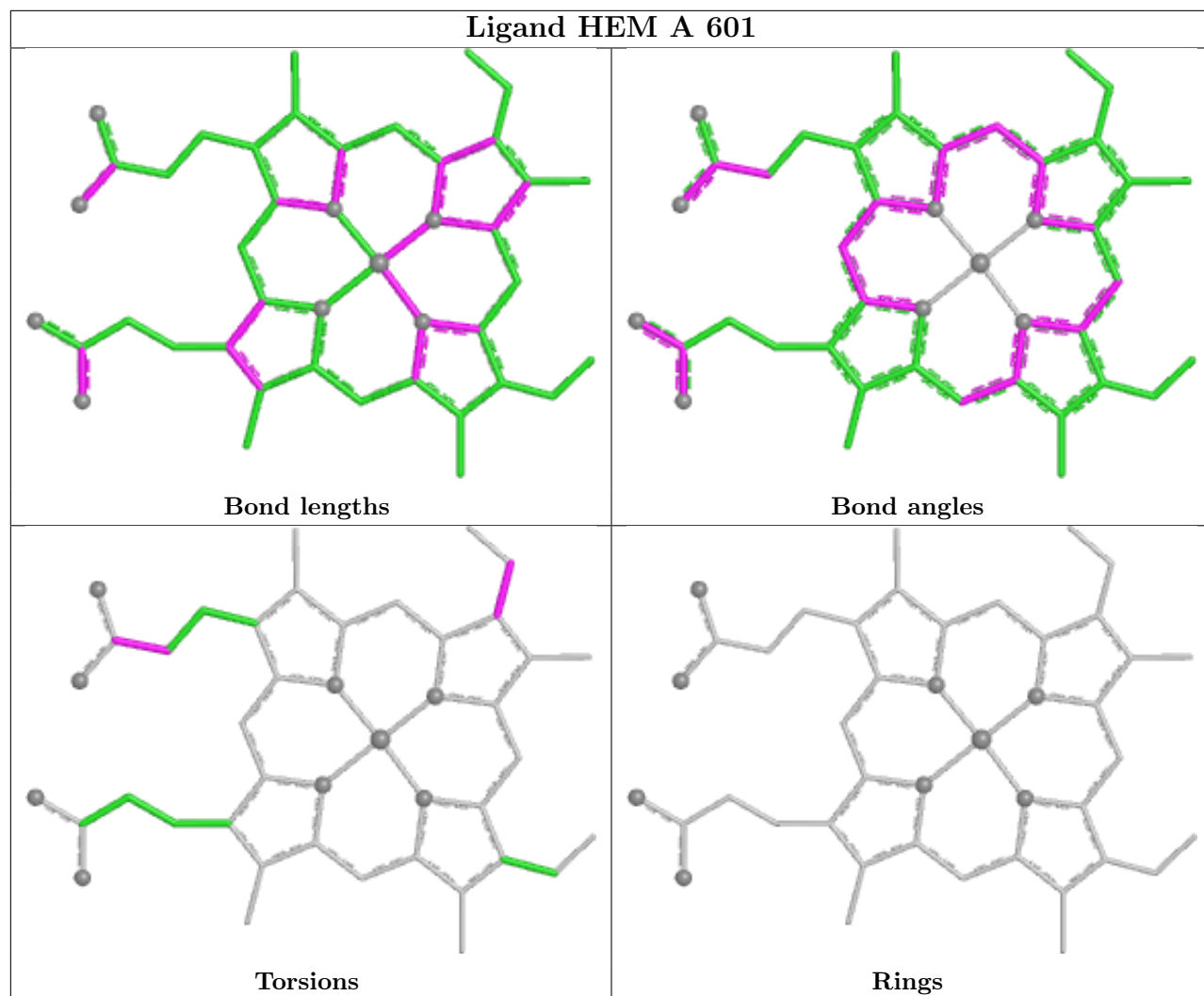
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	HEM	4	0
2	D	604	HEM	3	0
4	C	605	GOL	3	0
2	C	601	HEM	2	0
9	B	608	PEG	1	0
4	A	606	GOL	1	0
2	B	602	HEM	1	0

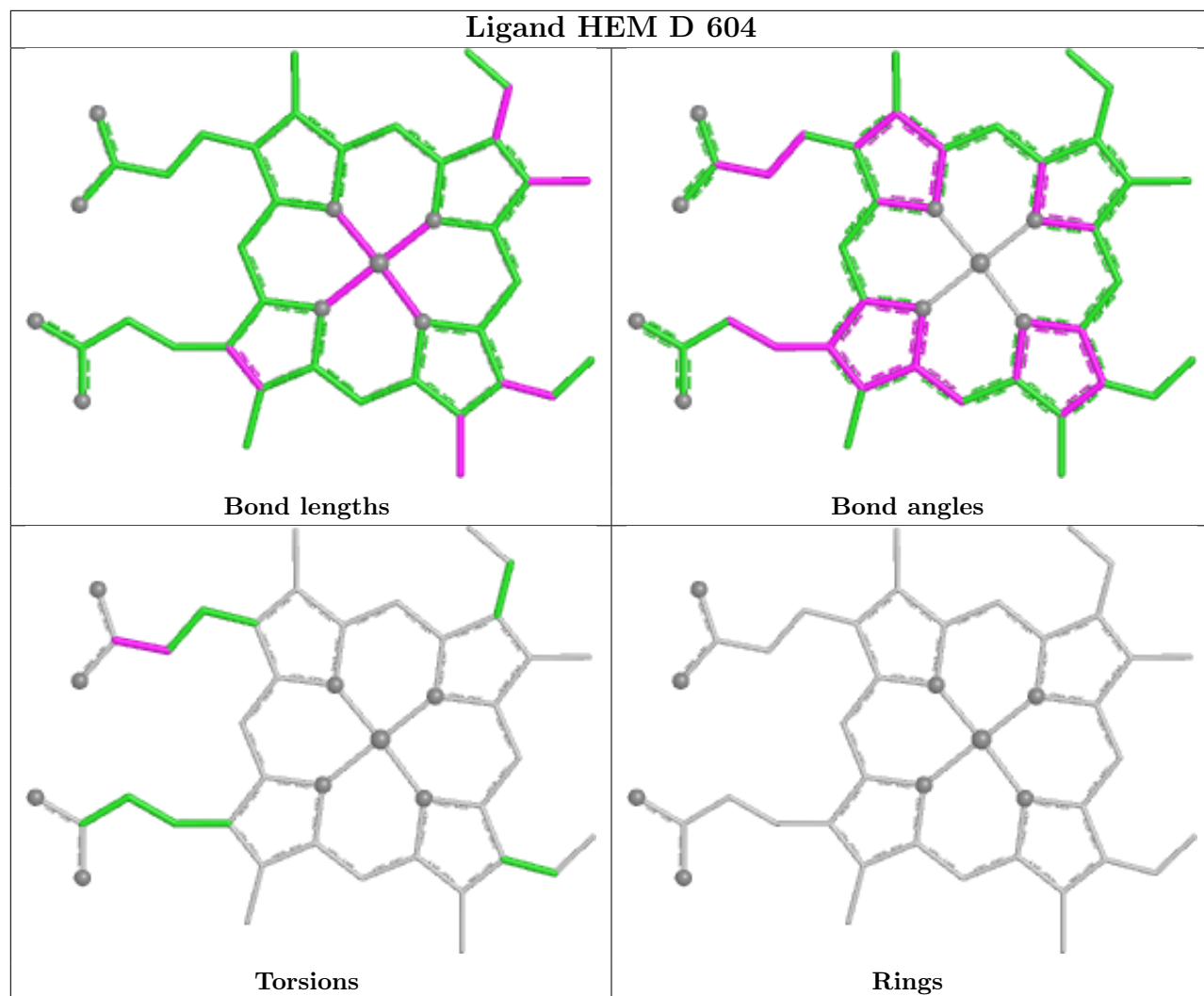
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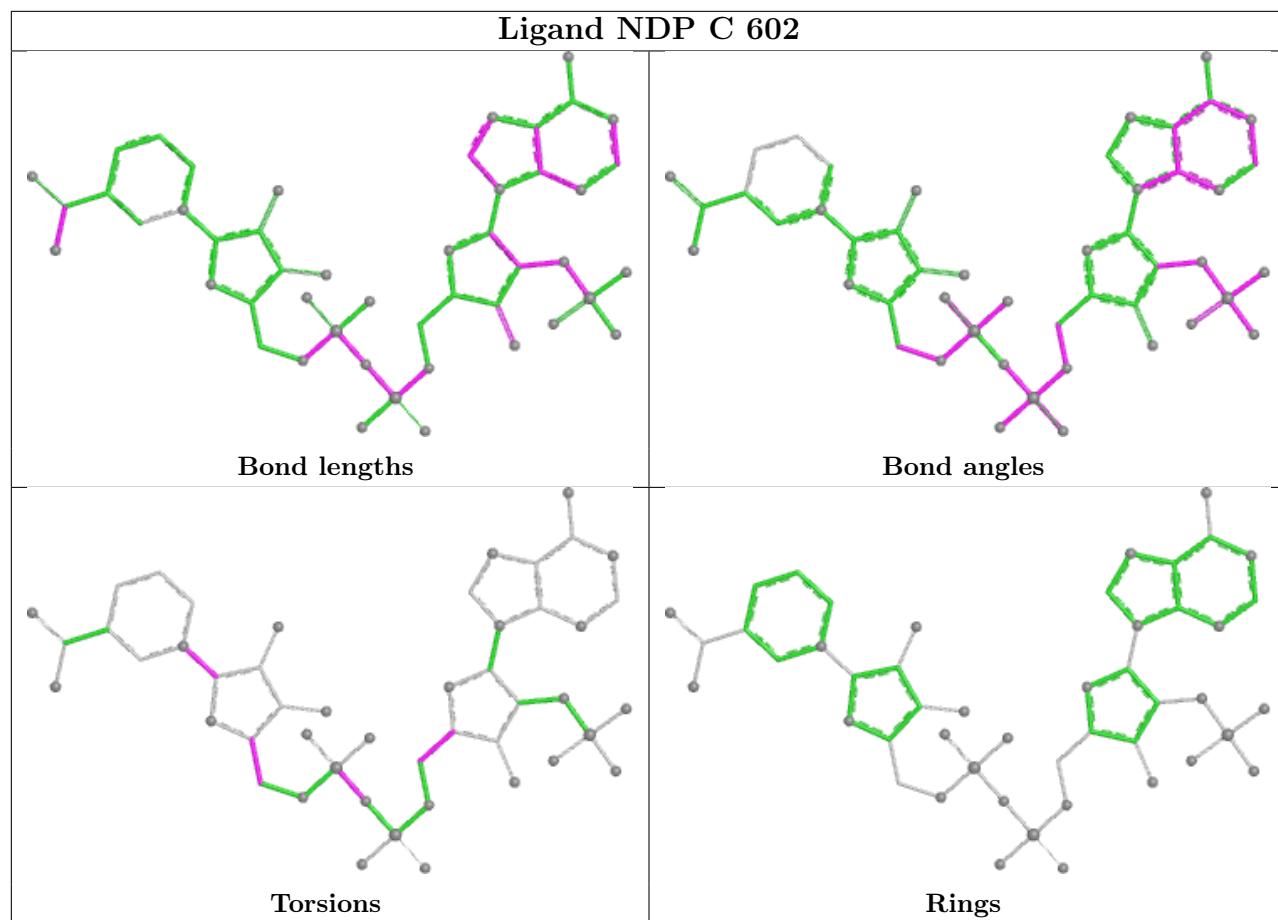
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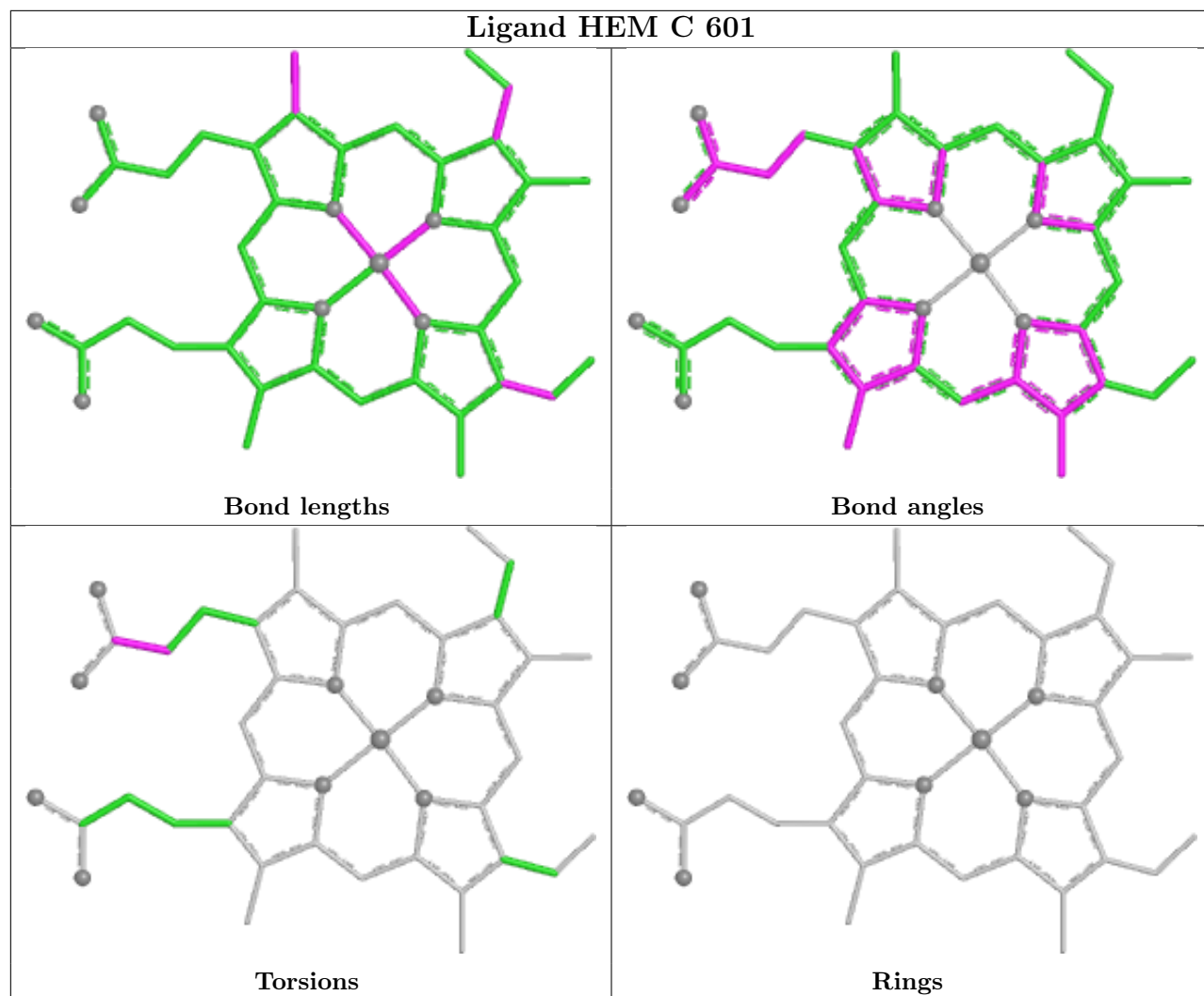
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	601	GOL	6	0
9	C	608	PEG	3	0
4	C	603	GOL	1	0
4	D	603	GOL	1	0
4	C	604	GOL	5	0

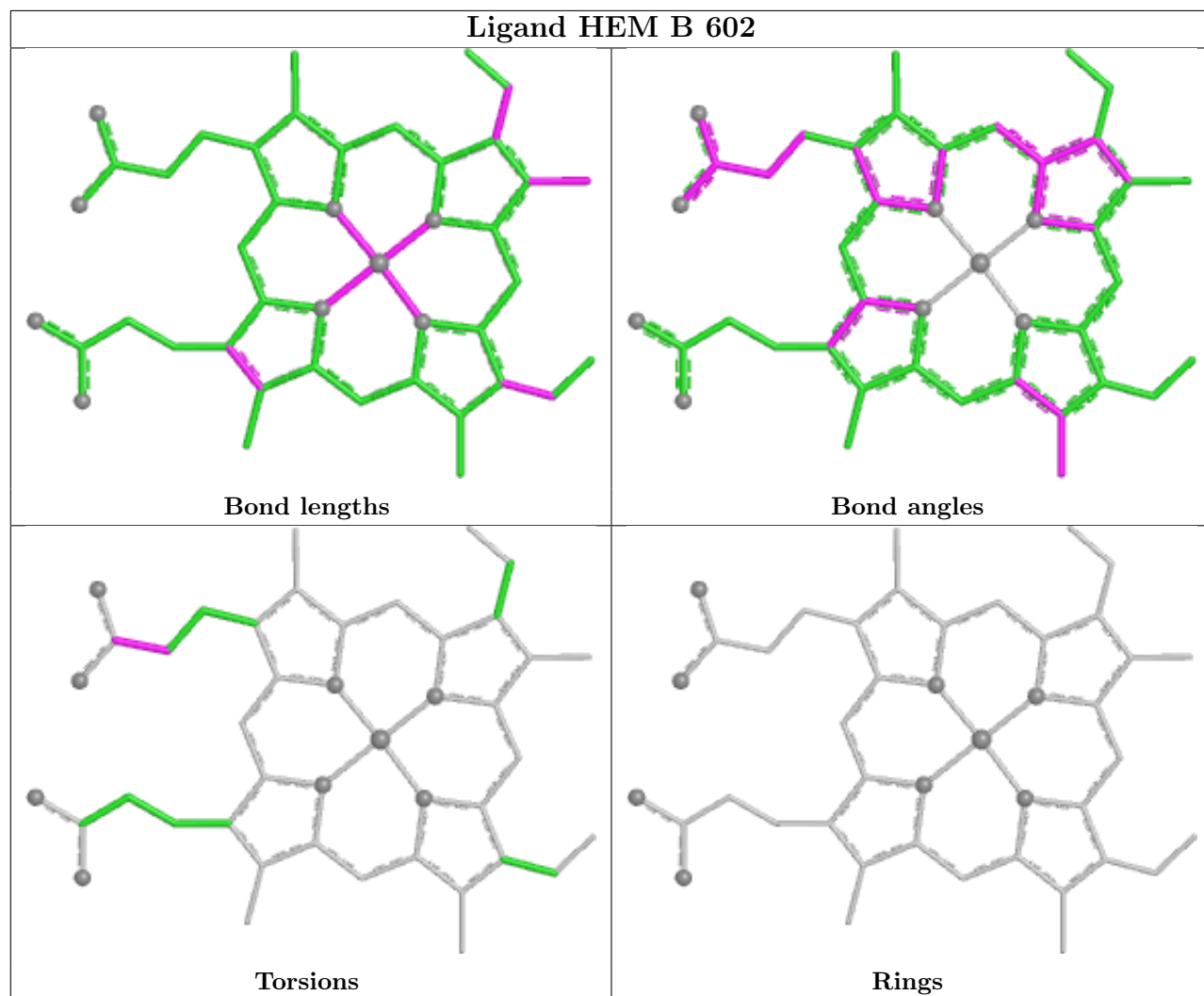
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

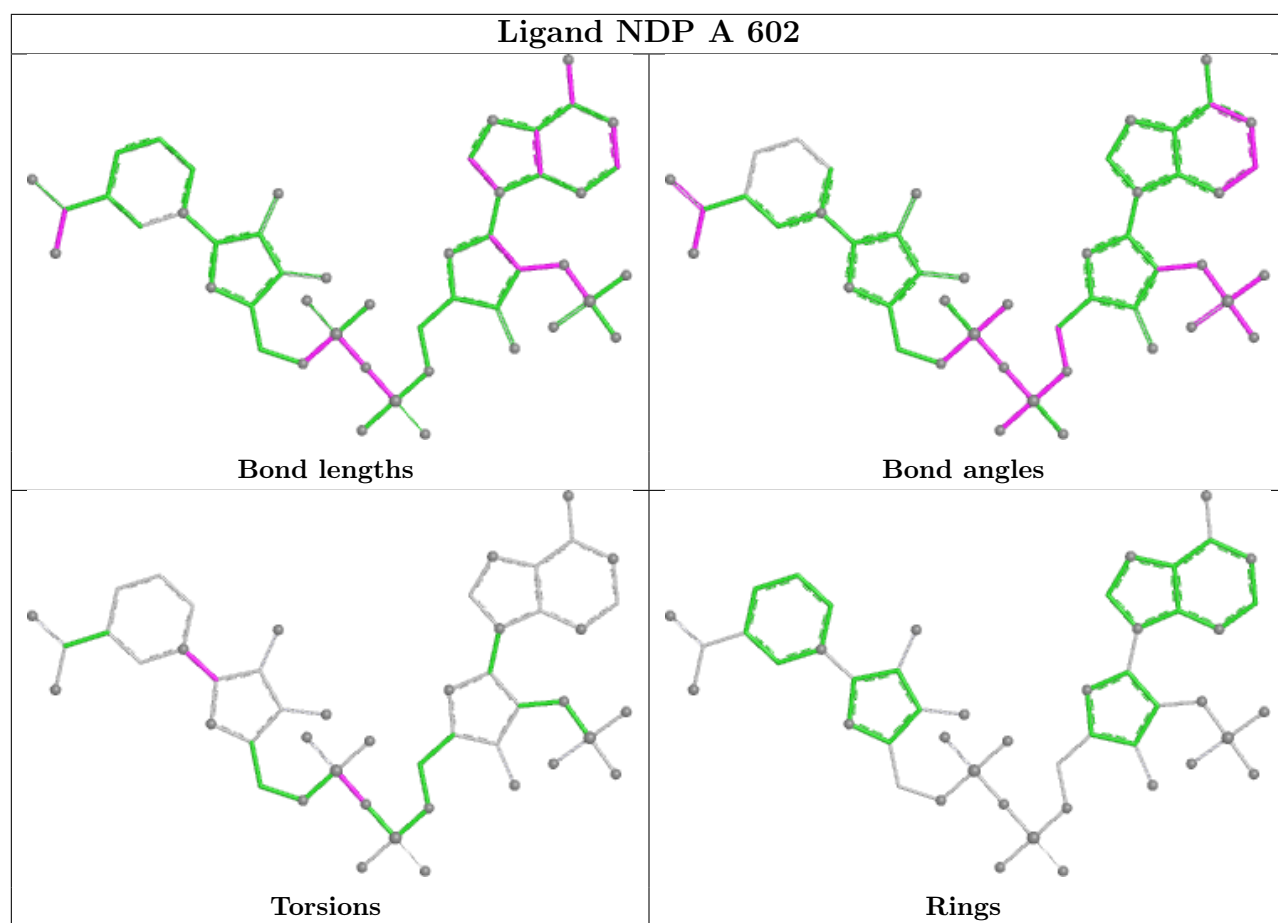


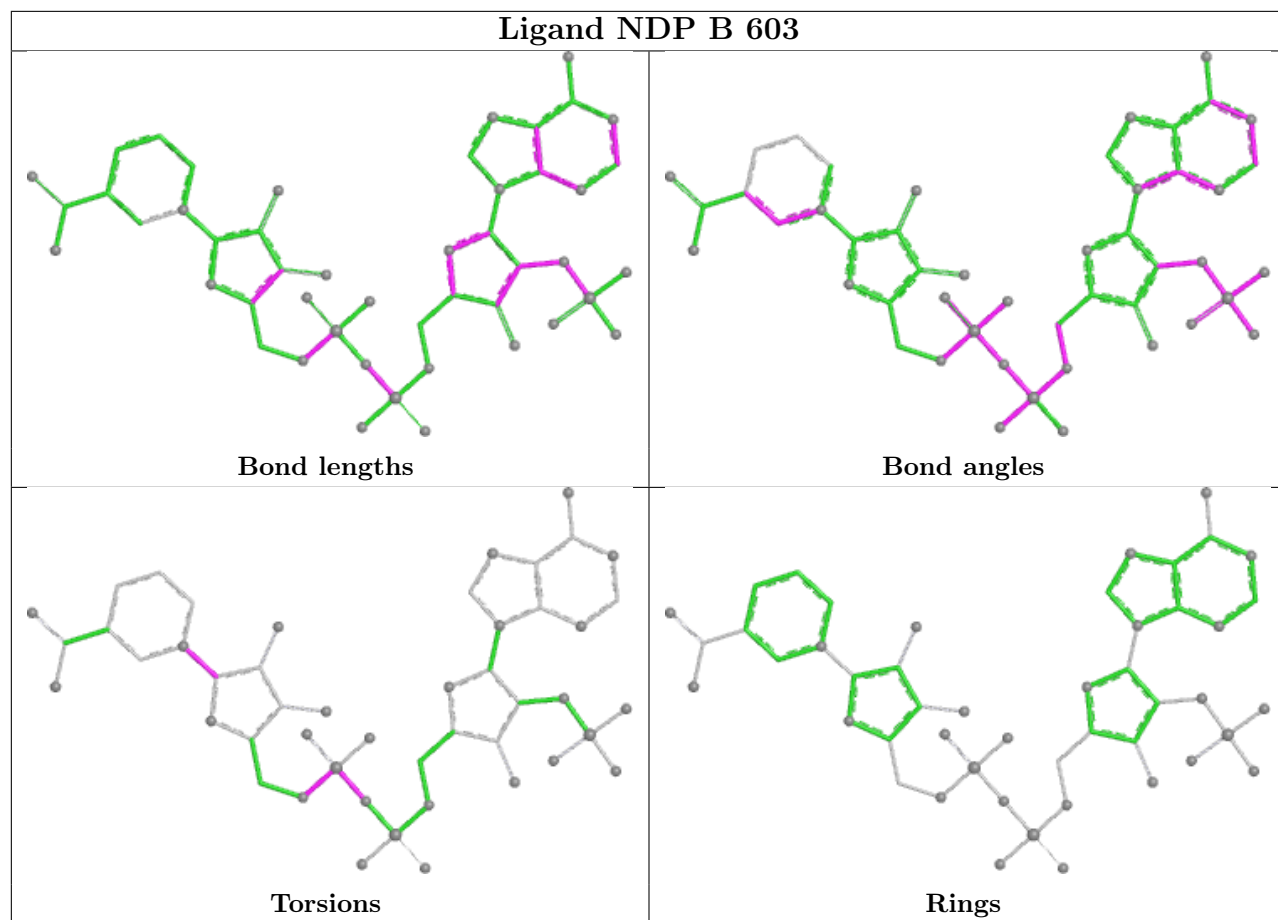


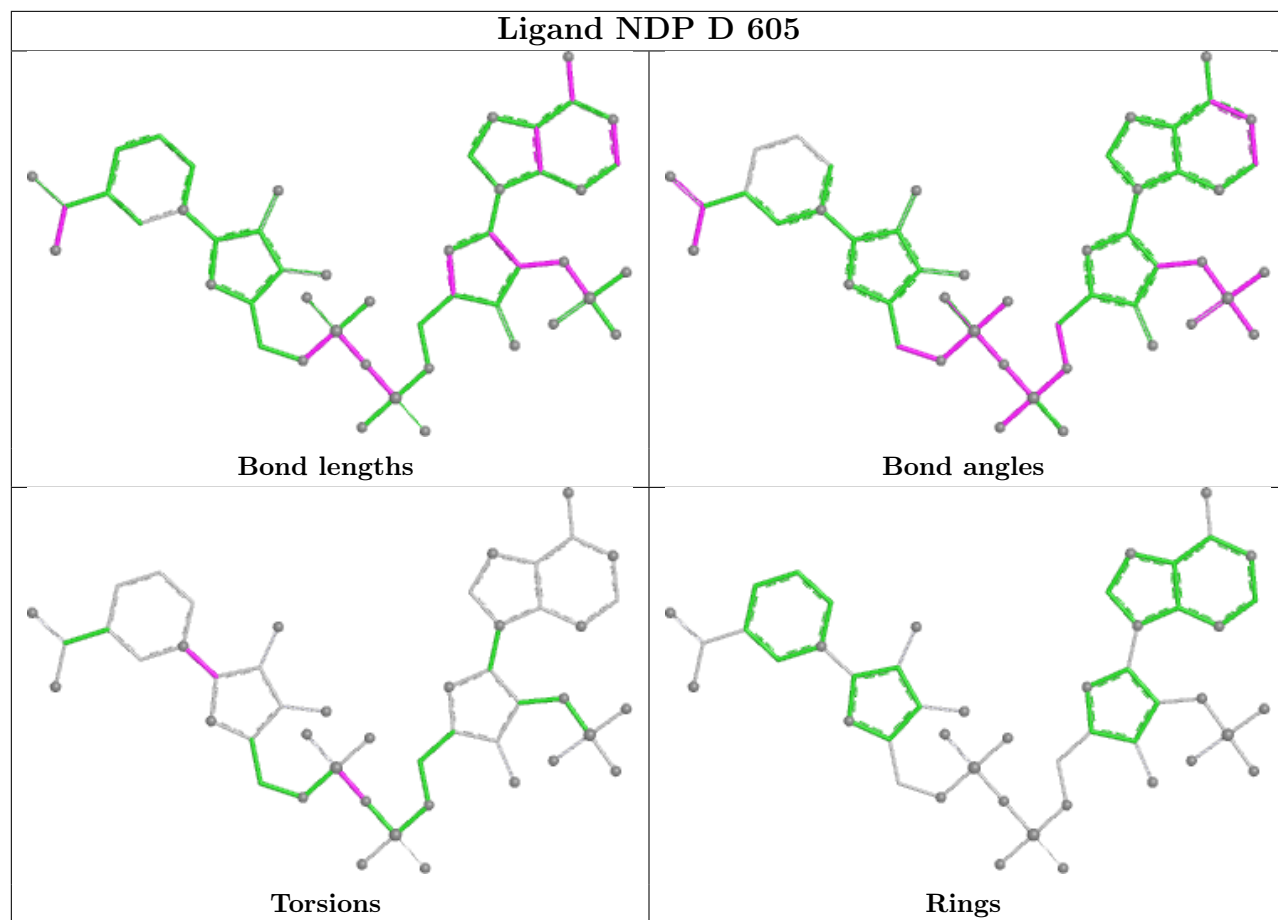












5.7 Other polymers [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	502/512 (98%)	-0.53	5 (0%) 79 80	20, 29, 53, 92	0
1	B	502/512 (98%)	-0.56	2 (0%) 88 89	19, 28, 50, 85	0
1	C	502/512 (98%)	-0.49	5 (0%) 79 80	14, 30, 49, 88	4 (0%)
1	D	502/512 (98%)	-0.55	3 (0%) 85 86	13, 30, 49, 91	5 (0%)
All	All	2008/2048 (98%)	-0.53	15 (0%) 84 85	13, 29, 50, 92	9 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	408	ILE	3.8
1	B	408	ILE	3.4
1	C	504	GLY	3.3
1	D	408	ILE	3.2
1	C	408	ILE	2.9
1	D	504	GLY	2.7
1	A	413	HIS	2.6
1	C	413	HIS	2.6
1	B	504	GLY	2.5
1	D	413	HIS	2.4
1	A	504	GLY	2.4
1	C	409	GLN	2.1
1	A	407	GLN	2.1
1	A	406	ALA	2.0
1	C	400	CYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	A	604	6/6	0.71	0.19	45,50,59,68	0
4	GOL	B	605	6/6	0.76	0.17	62,67,74,74	0
4	GOL	D	603	6/6	0.76	0.21	61,62,66,67	0
4	GOL	A	606	6/6	0.77	0.21	59,64,68,69	0
9	PEG	B	608	7/7	0.77	0.20	49,57,67,68	0
4	GOL	C	605	6/6	0.80	0.20	39,55,60,63	0
9	PEG	C	608	7/7	0.80	0.19	56,62,69,71	0
4	GOL	B	604	6/6	0.82	0.17	54,62,62,65	0
4	GOL	B	607	6/6	0.82	0.17	53,60,68,81	0
4	GOL	D	606	6/6	0.83	0.15	63,66,73,78	0
4	GOL	D	601	6/6	0.85	0.24	50,60,63,64	0
5	SO4	A	607	5/5	0.86	0.17	34,39,40,70	5
5	SO4	A	608	5/5	0.86	0.18	39,43,52,66	5
4	GOL	B	606	6/6	0.87	0.14	55,62,71,78	0
4	GOL	A	605	6/6	0.87	0.14	40,61,62,70	0
4	GOL	D	602	6/6	0.87	0.21	50,53,59,59	0
4	GOL	C	607	6/6	0.88	0.12	52,55,59,64	0
4	GOL	A	603	6/6	0.89	0.12	51,56,59,68	0
4	GOL	C	604	6/6	0.89	0.13	55,58,65,67	0
4	GOL	C	606	6/6	0.90	0.20	52,55,64,68	0
4	GOL	C	603	6/6	0.91	0.13	45,49,53,54	0
8	NA	B	612	1/1	0.93	0.05	50,50,50,50	0
6	CL	D	607	1/1	0.94	0.10	64,64,64,64	0
7	K	D	608	1/1	0.96	0.20	51,51,51,51	0
7	K	B	601	1/1	0.96	0.29	51,51,51,51	0
8	NA	D	609	1/1	0.96	0.10	49,49,49,49	0
7	K	B	613	1/1	0.96	0.39	62,62,62,62	0
7	K	C	611	1/1	0.96	0.26	57,57,57,57	0
3	NDP	C	602	48/48	0.97	0.06	25,35,45,51	0
7	K	A	611	1/1	0.97	0.15	51,51,51,51	0
3	NDP	D	605	48/48	0.97	0.06	25,32,37,42	0
2	HEM	A	601	43/43	0.97	0.06	18,23,28,32	0

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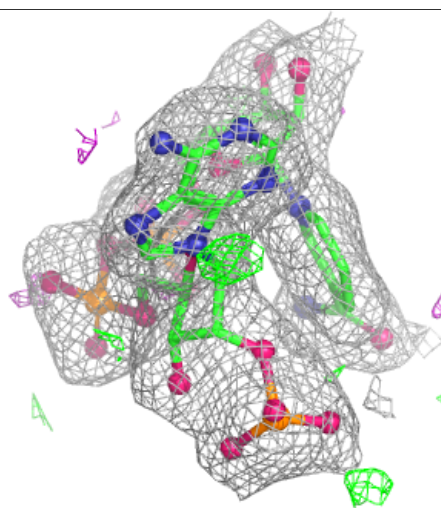
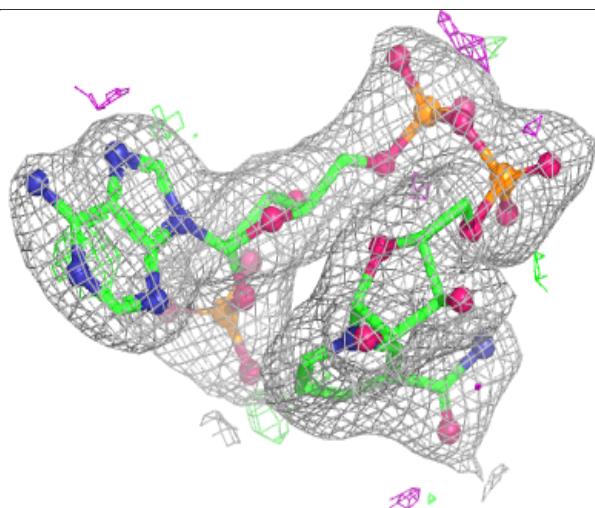
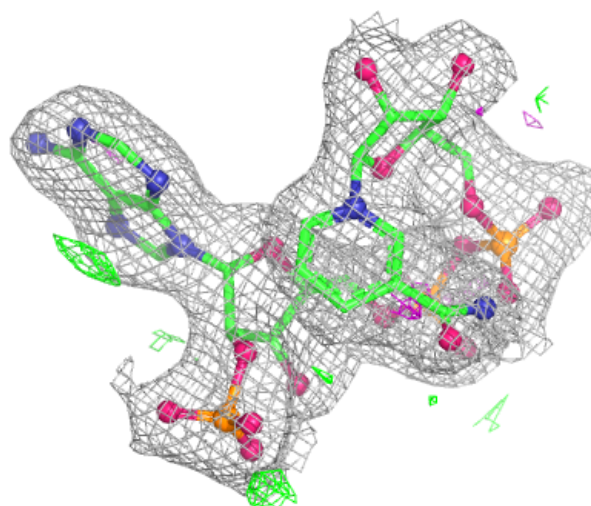
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	CL	A	610	1/1	0.97	0.07	50,50,50,50	0
7	K	C	609	1/1	0.98	0.14	54,54,54,54	0
6	CL	B	610	1/1	0.98	0.08	46,46,46,46	0
3	NDP	A	602	48/48	0.98	0.05	25,30,35,41	0
8	NA	A	612	1/1	0.98	0.08	37,37,37,37	0
3	NDP	B	603	48/48	0.98	0.04	21,28,37,44	0
8	NA	C	610	1/1	0.98	0.04	41,41,41,41	0
2	HEM	B	602	43/43	0.98	0.05	17,23,29,30	0
7	K	B	611	1/1	0.98	0.15	46,46,46,46	0
2	HEM	D	604	43/43	0.98	0.05	16,24,29,34	0
6	CL	B	609	1/1	0.99	0.03	35,35,35,35	0
2	HEM	C	601	43/43	0.99	0.05	17,23,30,34	0
6	CL	A	609	1/1	1.00	0.02	35,35,35,35	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.

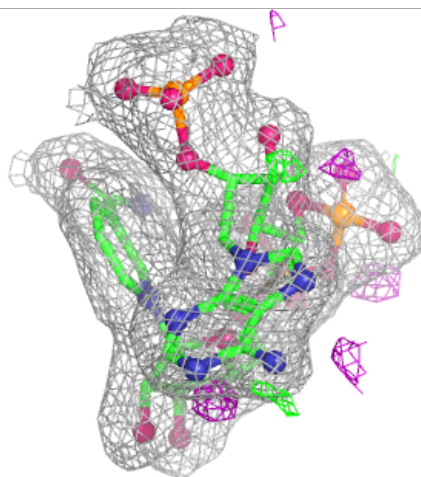
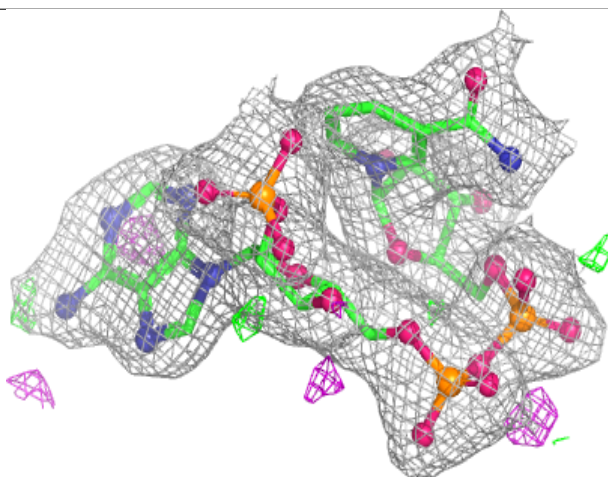
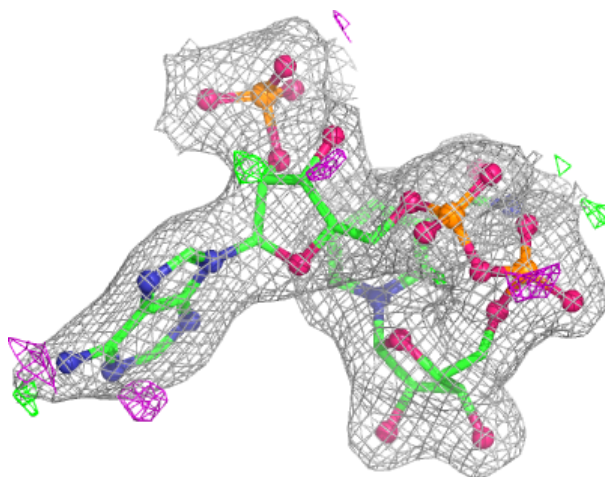
Electron density around NDP C 602:

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



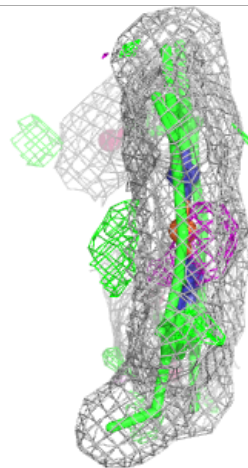
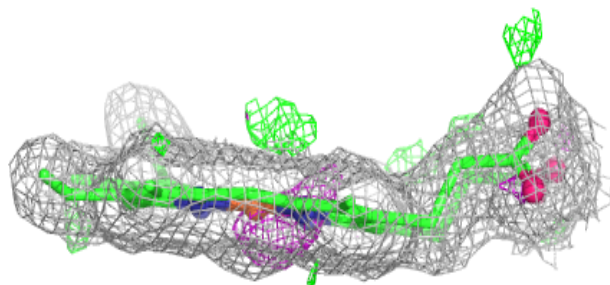
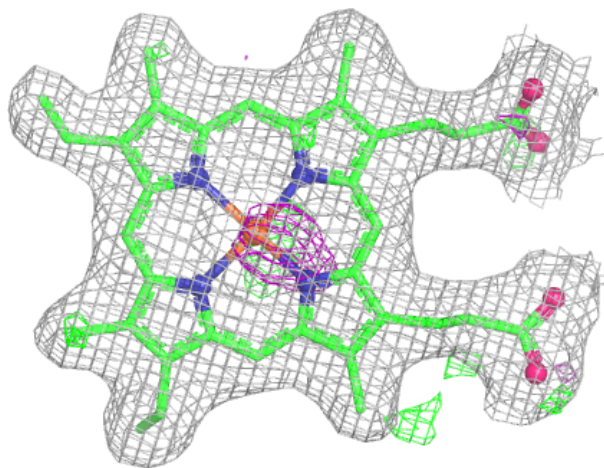
Electron density around NDP D 605:

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



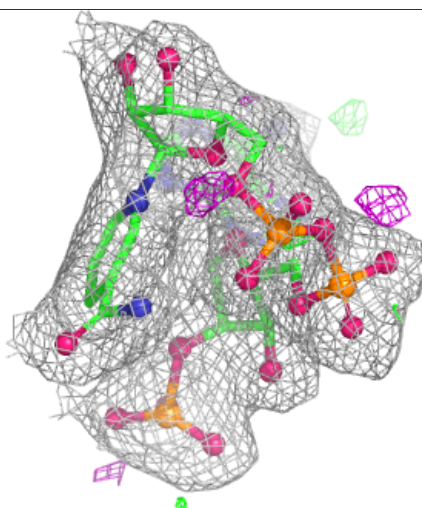
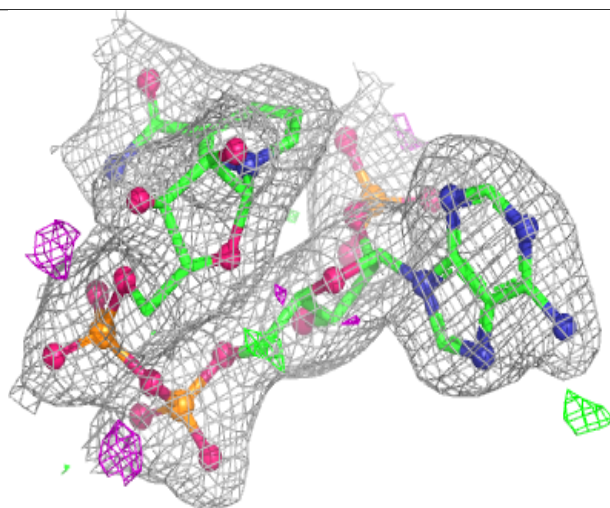
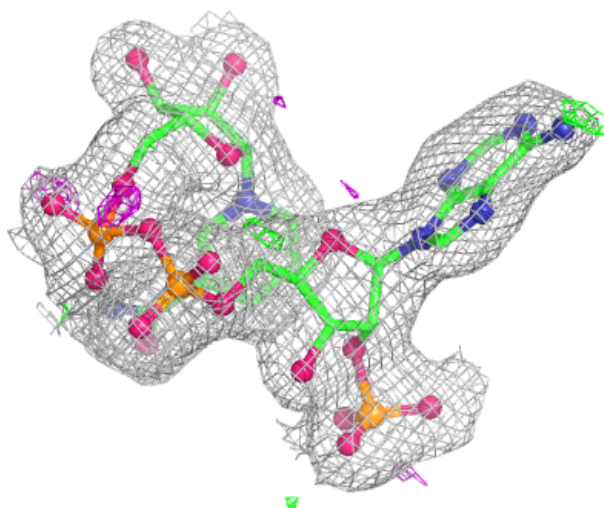
Electron density around HEM A 601:

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



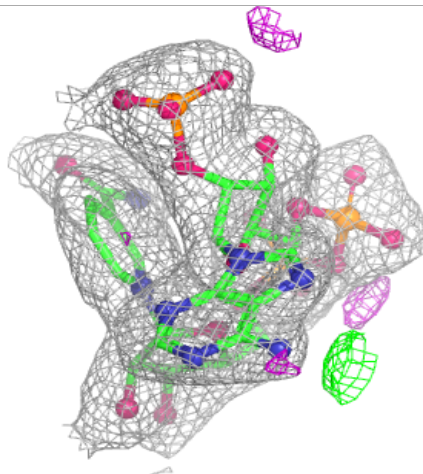
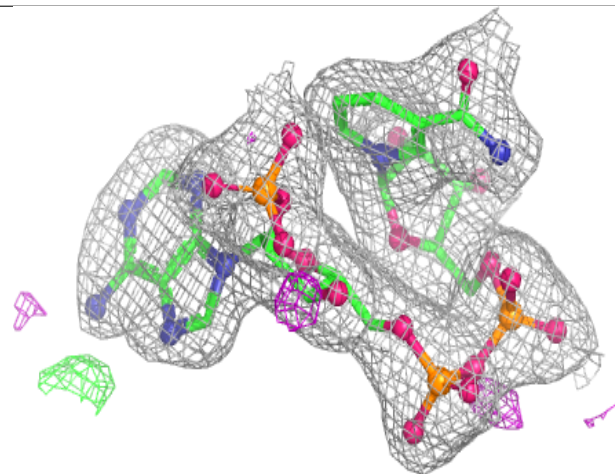
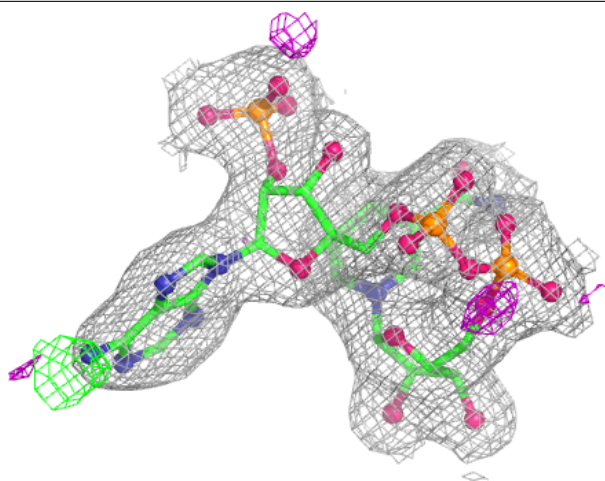
Electron density around NDP A 602:

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



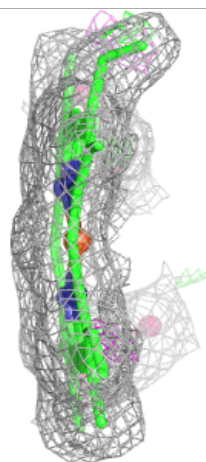
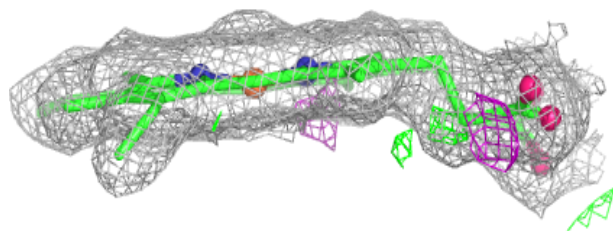
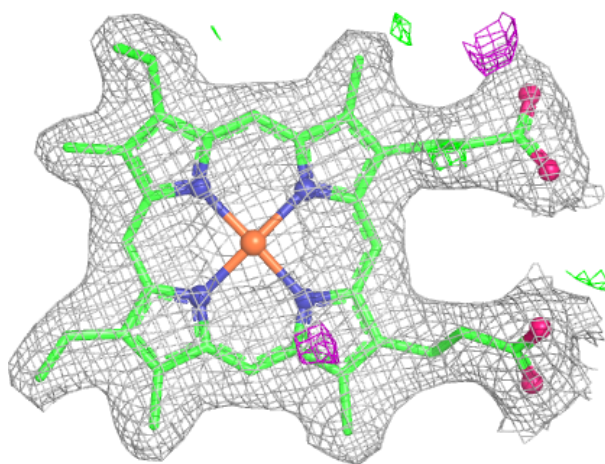
Electron density around NDP B 603:

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



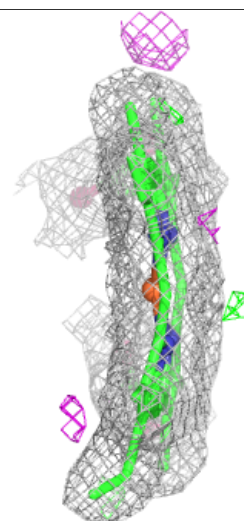
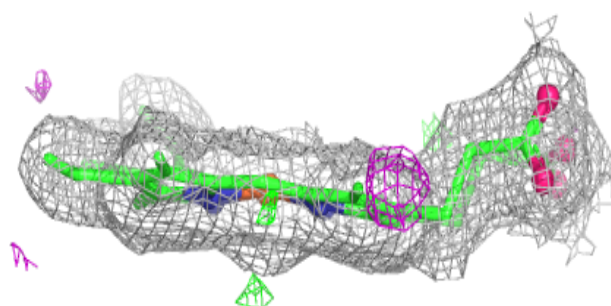
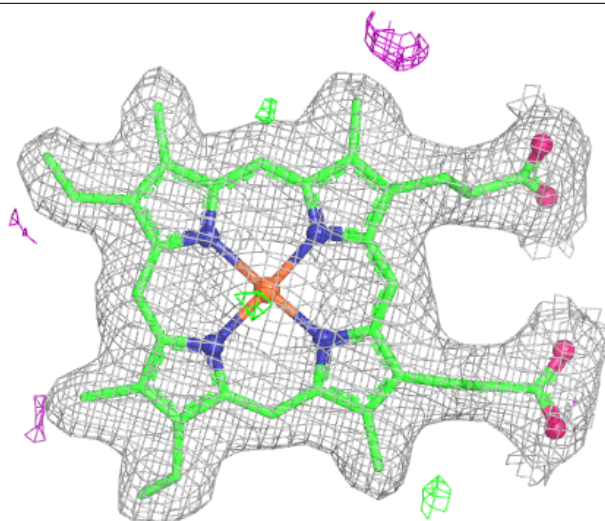
Electron density around HEM B 602:

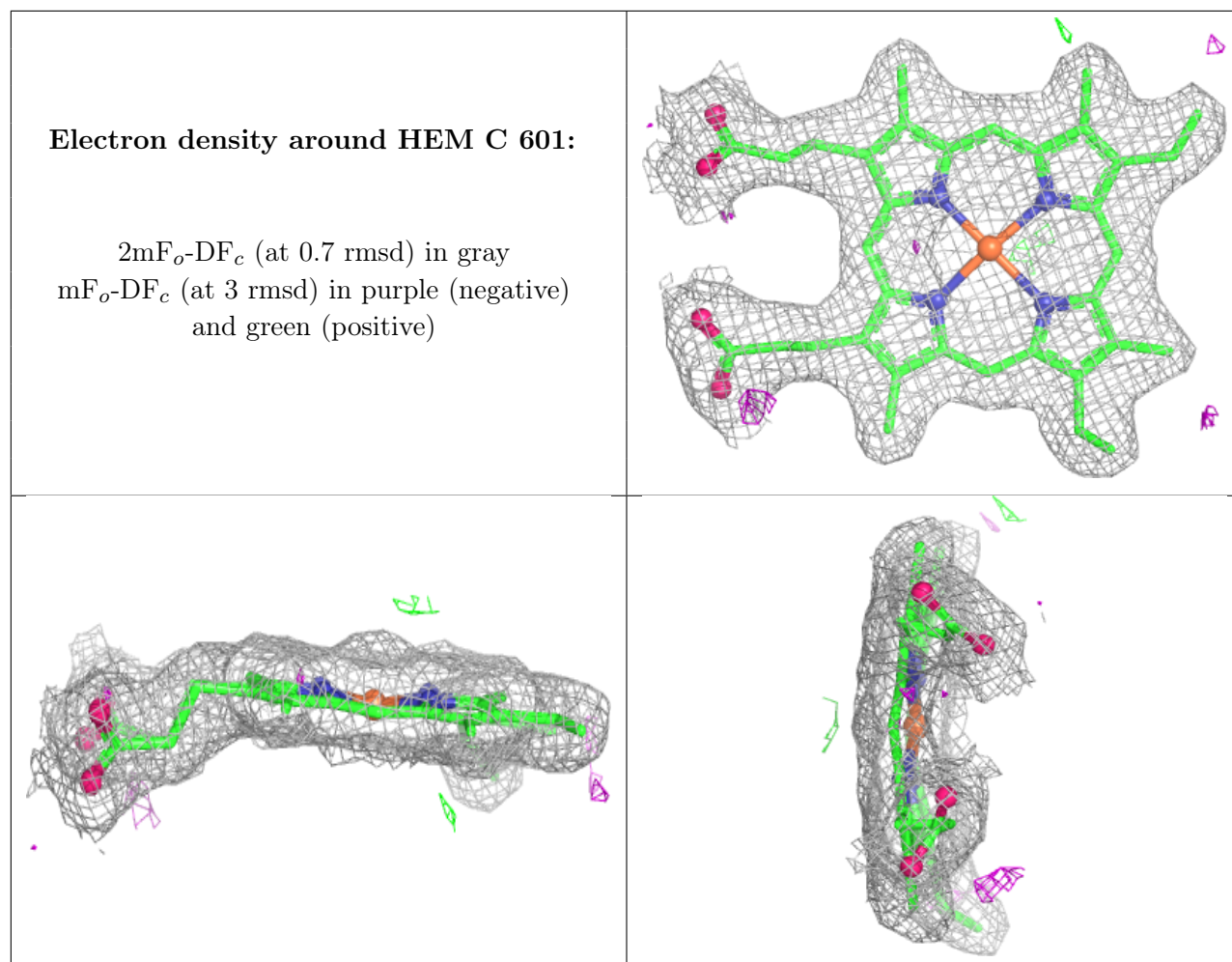
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM D 604:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.