



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:18 AM UTC

PDB ID : 7UFE / pdb_00007ufe
Title : Human CYP3A4 bound to an inhibitor
Authors : Sevrioukova, I.F.
Deposited on : 2022-03-22
Resolution : 2.40 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

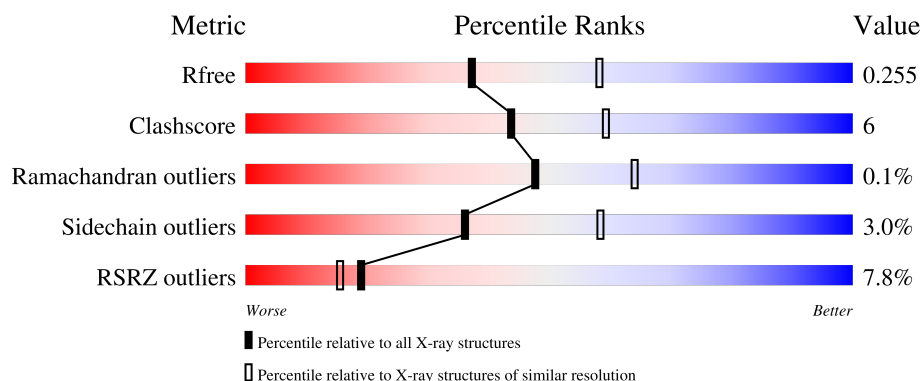
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	487	5% 78% 16% • 6%
1	B	487	9% 75% 16% • 8%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7517 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 Cytochrome P450 3A4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	460	Total	C	N	O	S	0	2	0
			3706	2411	612	659	24			
1	B	446	Total	C	N	O	S	0	0	0
			3576	2329	588	635	24			

There are 52 discrepancies between the modelled and reference sequences:

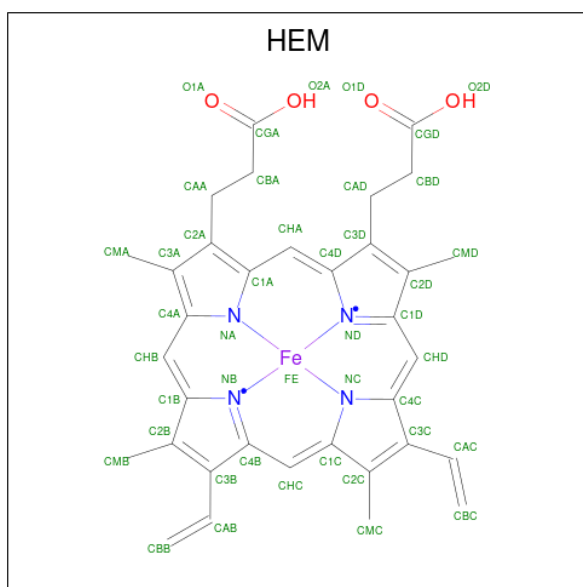
Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	LEU	deletion	UNP P08684
A	?	-	ILE	deletion	UNP P08684
A	?	-	PRO	deletion	UNP P08684
A	?	-	ASP	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	ALA	deletion	UNP P08684
A	?	-	MET	deletion	UNP P08684
A	?	-	GLU	deletion	UNP P08684
A	?	-	THR	deletion	UNP P08684
A	?	-	TRP	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	ALA	deletion	UNP P08684
A	?	-	VAL	deletion	UNP P08684
A	?	-	SER	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	VAL	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	?	-	LEU	deletion	UNP P08684
A	421	ALA	LYS	engineered mutation	UNP P08684
A	424	ALA	LYS	engineered mutation	UNP P08684
A	504	HIS	-	expression tag	UNP P08684
A	505	HIS	-	expression tag	UNP P08684
A	506	HIS	-	expression tag	UNP P08684

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Chain	Residue	Modelled	Actual	Comment	Reference
A	507	HIS	-	expression tag	UNP P08684
B	?	-	LEU	deletion	UNP P08684
B	?	-	ILE	deletion	UNP P08684
B	?	-	PRO	deletion	UNP P08684
B	?	-	ASP	deletion	UNP P08684
B	?	-	LEU	deletion	UNP P08684
B	?	-	ALA	deletion	UNP P08684
B	?	-	MET	deletion	UNP P08684
B	?	-	GLU	deletion	UNP P08684
B	?	-	THR	deletion	UNP P08684
B	?	-	TRP	deletion	UNP P08684
B	?	-	LEU	deletion	UNP P08684
B	?	-	LEU	deletion	UNP P08684
B	?	-	LEU	deletion	UNP P08684
B	?	-	ALA	deletion	UNP P08684
B	?	-	VAL	deletion	UNP P08684
B	?	-	SER	deletion	UNP P08684
B	?	-	LEU	deletion	UNP P08684
B	?	-	VAL	deletion	UNP P08684
B	?	-	LEU	deletion	UNP P08684
B	?	-	LEU	deletion	UNP P08684
B	421	ALA	LYS	engineered mutation	UNP P08684
B	424	ALA	LYS	engineered mutation	UNP P08684
B	504	HIS	-	expression tag	UNP P08684
B	505	HIS	-	expression tag	UNP P08684
B	506	HIS	-	expression tag	UNP P08684
B	507	HIS	-	expression tag	UNP P08684

- 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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



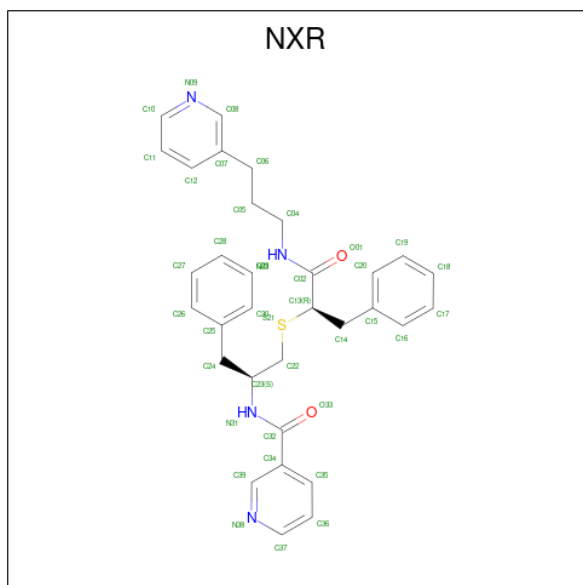
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0

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

- Molecule 4 is N-[(2S)-1- { [(2R)-1-oxo-3-phenyl-1- { [3-(pyridin-3-yl)propyl]amino } propan-2-yl]sulfanyl } -3-phenylpropan-2-yl]pyridine-3-carboxamide (CCD ID: NXR) (formula: $C_{32}H_{34}N_4O_2S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 39	C 32	N 4	O 2	S 1	0	0
4	B	1	Total 39	C 32	N 4	O 2	S 1	0	0

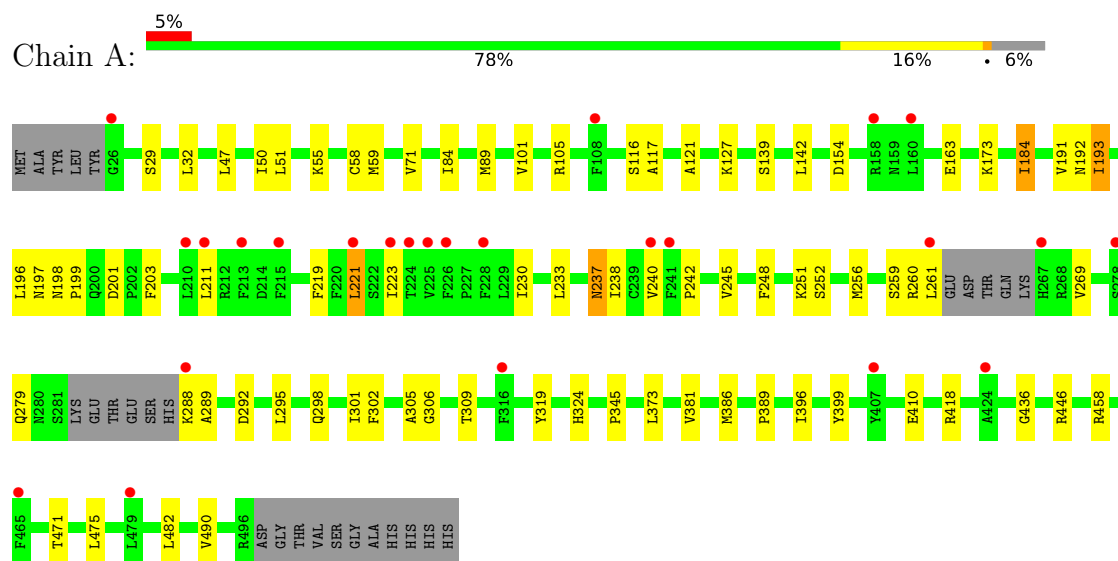
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	38	Total O 38 38	0	0
5	B	15	Total O 15 15	0	0

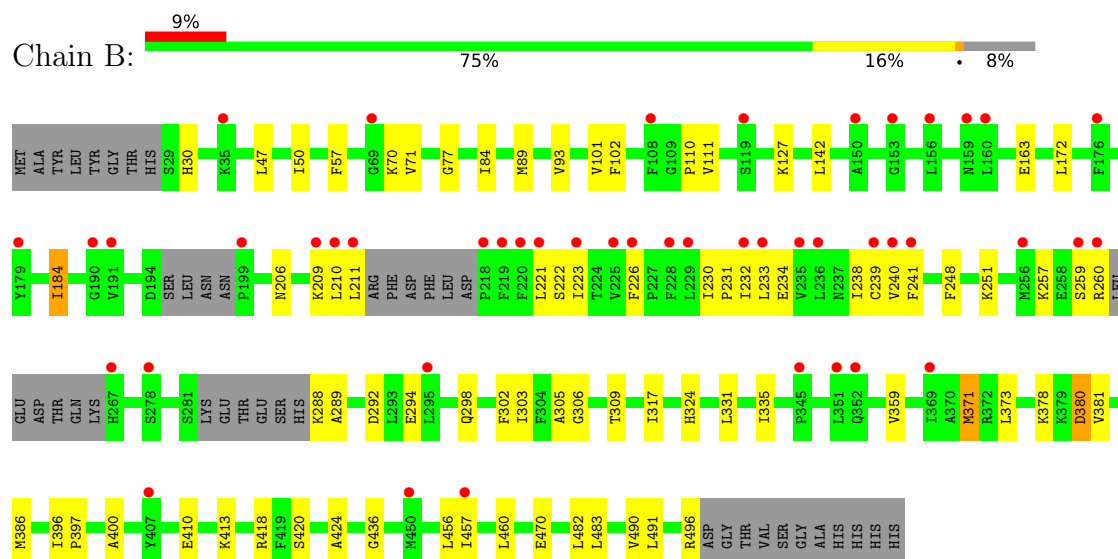
3 Residue-property plots [i](#)

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: Cytochrome P450 3A4



• Molecule 1: Cytochrome P450 3A4



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	153.62Å 96.75Å 93.87Å 90.00° 124.21° 90.00°	Depositor
Resolution (Å)	48.38 – 2.40 48.38 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.4 (48.38-2.40) 98.3 (48.38-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 2.39Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, R_{free}	0.221 , 0.253 0.225 , 0.255	Depositor DCC
R_{free} test set	2119 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	70.9	Xtriage
Anisotropy	0.262	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 60.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7517	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% 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: HEM, NXR, GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.13	0/3802	0.32	0/5143
1	B	0.11	0/3661	0.28	0/4948
All	All	0.12	0/7463	0.30	0/10091

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	3706	0	3792	42	0
1	B	3576	0	3667	44	0
2	A	43	0	30	5	0
2	B	43	0	30	4	0
3	A	12	0	16	1	0
3	B	6	0	8	2	0
4	A	39	0	0	0	0
4	B	39	0	0	1	0
5	A	38	0	0	0	0
5	B	15	0	0	0	0
All	All	7517	0	7543	91	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 6.

All (91) 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)
2:A:601:HEM:HHD	2:A:601:HEM:HBC2	1.68	0.76
2:B:601:HEM:HBC2	2:B:601:HEM:HHD	1.66	0.76
1:A:184:ILE:HG13	1:A:306:GLY:HA3	1.72	0.71
1:A:248:PHE:HA	1:A:251:LYS:HE2	1.73	0.71
1:B:184:ILE:HG13	1:B:306:GLY:HA3	1.74	0.69
1:B:222:SER:O	1:B:226:PHE:N	2.25	0.69
1:A:223:ILE:HG23	1:A:230:ILE:HD11	1.76	0.68
1:A:154:ASP:OD1	1:A:458:ARG:NH1	2.28	0.66
1:B:470:GLU:HG3	1:B:490:VAL:HG21	1.81	0.63
1:B:248:PHE:HA	1:B:251:LYS:HE2	1.80	0.63
1:B:410:GLU:O	1:B:418:ARG:NH2	2.32	0.62
1:A:47:LEU:HD22	1:A:50:ILE:HD11	1.82	0.61
1:B:239:CYS:SG	1:B:240:VAL:N	2.73	0.61
1:B:324:HIS:NE2	3:B:602:GOL:O3	2.35	0.60
1:B:206:ASN:HA	1:B:209:LYS:HE2	1.84	0.59
1:B:233:LEU:HB3	1:B:238:ILE:HB	1.85	0.58
1:A:116:SER:O	1:A:298:GLN:NE2	2.31	0.58
1:B:482:LEU:H	1:B:482:LEU:HD23	1.70	0.57
1:A:50:ILE:HD13	1:A:221:LEU:HD21	1.88	0.56
1:A:139:SER:HB3	1:A:446:ARG:HH22	1.72	0.55
1:B:302:PHE:CD2	2:B:601:HEM:HBC1	2.42	0.54
1:B:380:ASP:OD1	1:B:380:ASP:N	2.38	0.53
2:A:601:HEM:HBB2	2:A:601:HEM:HMB2	1.91	0.53
1:B:172:LEU:HD11	1:B:491:LEU:HD12	1.89	0.53
1:A:471:THR:OG1	1:A:490:VAL:O	2.22	0.53
1:A:55:LYS:HB3	1:A:59:MET:HB2	1.91	0.52
1:B:331:LEU:HD22	1:B:359:VAL:HG21	1.91	0.52
2:B:601:HEM:HBB2	2:B:601:HEM:HMB2	1.89	0.52
1:B:317:ILE:HD13	1:B:456:LEU:HD11	1.91	0.51
1:A:302:PHE:CD2	2:A:601:HEM:HBC1	2.45	0.51
1:A:482:LEU:H	1:A:482:LEU:HD23	1.76	0.51
1:B:436:GLY:HA3	2:B:601:HEM:HBA1	1.93	0.51
1:A:410:GLU:O	1:A:418:ARG:NH2	2.39	0.50
1:A:242:PRO:HG2	1:A:245:VAL:HB	1.92	0.50
1:B:397:PRO:HB2	1:B:400:ALA:HB3	1.93	0.50
1:A:324:HIS:NE2	3:A:603:GOL:O2	2.37	0.50
1:B:110:PRO:HD2	1:B:240:VAL:HG13	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:GLN:NE2	1:A:292:ASP:OD1	2.39	0.48
1:A:32:LEU:HD11	1:A:389:PRO:HG2	1.95	0.48
1:A:84:ILE:HB	1:A:89:MET:HE3	1.94	0.48
3:B:602:GOL:O3	3:B:602:GOL:O1	2.19	0.48
1:B:373:LEU:HB2	1:B:396:ILE:HB	1.97	0.47
1:B:84:ILE:HB	1:B:89:MET:HE3	1.97	0.47
1:A:436:GLY:HA3	2:A:601:HEM:HBA1	1.97	0.46
1:A:269:VAL:O	1:A:269:VAL:HG12	2.14	0.46
1:A:373:LEU:HB2	1:A:396:ILE:HB	1.98	0.46
1:B:101:VAL:HG21	1:B:381:VAL:HG11	1.98	0.46
1:B:50:ILE:HD13	1:B:221:LEU:HD11	1.98	0.46
1:B:239:CYS:SG	1:B:241:PHE:N	2.89	0.45
1:B:259:SER:HA	1:B:260:ARG:HA	1.56	0.45
1:B:70:LYS:HG3	1:B:71:VAL:HG23	1.99	0.45
1:B:257:LYS:NZ	1:B:292:ASP:OD1	2.50	0.44
1:A:219:PHE:O	1:A:223:ILE:HG12	2.17	0.44
1:B:420:SER:O	1:B:424:ALA:N	2.45	0.44
1:A:219:PHE:HE2	1:A:240:VAL:HG12	1.82	0.44
1:A:101:VAL:HG21	1:A:381:VAL:HG11	1.98	0.44
1:B:288:LYS:N	1:B:289:ALA:HA	2.32	0.44
1:B:71:VAL:HG21	1:B:386:MET:HE1	2.00	0.44
1:B:57:PHE:HB2	1:B:371:MET:HE3	1.99	0.43
1:A:193:ILE:H	1:A:193:ILE:HG13	1.71	0.43
1:A:345:PRO:HD3	1:A:458:ARG:HD3	2.01	0.43
1:A:259:SER:HA	1:A:260:ARG:C	2.43	0.43
1:A:305:ALA:O	1:A:309:THR:OG1	2.32	0.43
1:A:252:SER:O	1:A:256:MET:N	2.51	0.43
1:A:198:ASN:N	1:A:199:PRO:HD3	2.34	0.43
1:B:101:VAL:HA	1:B:378:LYS:HG2	2.00	0.42
1:B:410:GLU:HB3	1:B:413:LYS:HG3	2.01	0.42
1:A:58:CYS:HB3	1:A:399:TYR:CE2	2.54	0.42
1:A:191:VAL:HG21	1:A:256:MET:HG3	2.01	0.42
1:B:210:LEU:HD13	4:B:603:NXR:C18	2.50	0.42
1:B:231:PRO:HA	1:B:234:GLU:HB2	2.02	0.42
1:B:30:HIS:NE2	1:B:77:GLY:O	2.52	0.42
1:B:184:ILE:HG21	1:B:303:ILE:HA	2.02	0.41
1:B:230:ILE:HB	1:B:231:PRO:HD3	2.02	0.41
1:B:335:ILE:HD13	1:B:457:ILE:HA	2.01	0.41
1:B:47:LEU:HD22	1:B:50:ILE:HD11	2.02	0.41
1:A:117:ALA:O	1:A:121:ALA:N	2.49	0.41
1:B:371:MET:SD	1:B:483:LEU:HB2	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:ALA:O	1:B:309:THR:OG1	2.30	0.41
1:B:460:LEU:O	1:B:496:ARG:NH1	2.54	0.41
1:A:237:ASN:O	1:A:237:ASN:ND2	2.48	0.41
1:A:117:ALA:HB1	1:A:301:ILE:HG13	2.03	0.41
1:B:93:VAL:HG13	1:B:102:PHE:CG	2.56	0.41
1:B:294:GLU:O	1:B:298:GLN:HG2	2.20	0.41
1:A:319:TYR:CZ	1:A:475:LEU:HB2	2.55	0.41
1:A:233:LEU:HB3	1:A:238:ILE:HB	2.02	0.40
1:A:58:CYS:HB3	1:A:399:TYR:CD2	2.56	0.40
1:A:201:ASP:OD2	1:A:203:PHE:HB2	2.22	0.40
1:A:288:LYS:HB3	1:A:289:ALA:HA	2.03	0.40
1:A:105:ARG:HH21	2:A:601:HEM:HAA1	1.86	0.40
1:A:71:VAL:HG21	1:A:386:MET:HE1	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	456/487 (94%)	426 (93%)	29 (6%)	1 (0%)	43	58
1	B	436/487 (90%)	407 (93%)	29 (7%)	0	100	100
All	All	892/974 (92%)	833 (93%)	58 (6%)	1 (0%)	48	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	197	ASN

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	420/441 (95%)	405 (96%)	15 (4%)	31	52
1	B	405/441 (92%)	395 (98%)	10 (2%)	42	64
All	All	825/882 (94%)	800 (97%)	25 (3%)	36	58

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	SER
1	A	51	LEU
1	A	127	LYS
1	A	142	LEU
1	A	163	GLU
1	A	173	LYS
1	A	184	ILE
1	A	192	ASN
1	A	193	ILE
1	A	196	LEU
1	A	211	LEU
1	A	221	LEU
1	A	237	ASN
1	A	261	LEU
1	A	295	LEU
1	B	111	VAL
1	B	127	LYS
1	B	142	LEU
1	B	163	GLU
1	B	184	ILE
1	B	211	LEU
1	B	223	ILE
1	B	232	ILE
1	B	371	MET
1	B	380	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	280	ASN
1	A	451	ASN
1	A	472	GLN
1	B	484	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

7 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	$\# Z > 2$
2	HEM	A	601	1,4	50,50,50	1.43	7 (14%)	67,82,82	1.18	4 (5%)
3	GOL	A	603	-	5,5,5	0.37	0	5,5,5	0.38	0
3	GOL	B	602	-	5,5,5	0.38	0	5,5,5	0.48	0
4	NXR	A	604	2	41,42,42	1.48	3 (7%)	47,53,53	1.43	6 (12%)
4	NXR	B	603	2	41,42,42	1.48	4 (9%)	47,53,53	1.45	6 (12%)
3	GOL	A	602	-	5,5,5	0.37	0	5,5,5	0.28	0
2	HEM	B	601	1,4	50,50,50	1.44	7 (14%)	67,82,82	1.15	5 (7%)

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	-	2/14/54/54	-
3	GOL	A	603	-	-	0/4/4/4	-
3	GOL	B	602	-	-	2/4/4/4	-
4	NXR	A	604	2	-	3/32/32/32	0/4/4/4
4	NXR	B	603	2	-	9/32/32/32	0/4/4/4
3	GOL	A	602	-	-	2/4/4/4	-
2	HEM	B	601	1,4	-	1/14/54/54	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	604	NXR	C32-N31	5.54	1.47	1.34
4	B	603	NXR	C02-N03	5.38	1.46	1.33
4	B	603	NXR	C32-N31	5.35	1.46	1.34
4	A	604	NXR	C02-N03	5.12	1.45	1.33
2	A	601	HEM	FE-NB	3.40	2.05	1.94
2	B	601	HEM	FE-NB	3.38	2.05	1.94
2	A	601	HEM	FE-NC	3.31	2.06	1.95
2	B	601	HEM	FE-NC	3.29	2.06	1.95
2	B	601	HEM	FE-ND	3.23	2.04	1.94
2	A	601	HEM	FE-NA	3.19	2.05	1.95
2	A	601	HEM	FE-ND	3.14	2.04	1.94
2	B	601	HEM	FE-NA	3.14	2.05	1.95
2	A	601	HEM	CAB-C3B	3.12	1.55	1.47
2	B	601	HEM	CAB-C3B	3.11	1.55	1.47
2	A	601	HEM	CAC-C3C	2.98	1.55	1.47
2	B	601	HEM	CAC-C3C	2.96	1.55	1.47
4	B	603	NXR	C22-S21	-2.35	1.79	1.82
2	B	601	HEM	CMB-C2B	2.14	1.55	1.50
2	A	601	HEM	CMB-C2B	2.10	1.55	1.50
4	B	603	NXR	C13-S21	-2.10	1.81	1.83
4	A	604	NXR	C13-S21	-2.05	1.81	1.83

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	603	NXR	C34-C32-N31	4.48	125.35	117.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	604	NXR	C34-C32-N31	4.14	124.72	117.04
4	B	603	NXR	C24-C23-N31	3.83	117.64	110.34
4	A	604	NXR	O01-C02-N03	-3.08	116.45	122.98
4	B	603	NXR	O33-C32-N31	-3.07	116.64	122.47
4	A	604	NXR	C35-C34-C39	-2.83	114.46	117.61
4	A	604	NXR	O33-C32-N31	-2.78	117.19	122.47
4	B	603	NXR	C35-C34-C39	-2.68	114.63	117.61
4	A	604	NXR	C15-C14-C13	2.55	117.01	113.35
4	B	603	NXR	O01-C02-N03	-2.48	117.74	122.98
2	A	601	HEM	C4D-ND-C1D	2.48	108.14	105.21
2	B	601	HEM	C4D-ND-C1D	2.47	108.13	105.21
4	A	604	NXR	C23-N31-C32	2.30	126.90	122.56
4	B	603	NXR	C04-N03-C02	2.26	126.60	122.55
2	A	601	HEM	C1B-NB-C4B	2.23	107.85	105.21
2	B	601	HEM	C1B-NB-C4B	2.23	107.84	105.21
2	A	601	HEM	C3D-C4D-ND	-2.20	107.76	110.17
2	A	601	HEM	C2A-C1A-NA	-2.15	107.76	110.15
2	B	601	HEM	C3B-C2B-C1B	2.11	108.00	106.41
2	B	601	HEM	C3D-C4D-ND	-2.10	107.87	110.17
2	B	601	HEM	C2A-C1A-NA	-2.08	107.84	110.15

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	GOL	O1-C1-C2-C3
3	B	602	GOL	O1-C1-C2-C3
4	B	603	NXR	C02-C13-S21-C22
4	B	603	NXR	C14-C13-S21-C22
4	B	603	NXR	S21-C22-C23-C24
4	B	603	NXR	S21-C22-C23-N31
4	A	604	NXR	N03-C04-C05-C06
3	A	602	GOL	O1-C1-C2-O2
4	A	604	NXR	C04-C05-C06-C07
4	B	603	NXR	N03-C04-C05-C06
3	B	602	GOL	O1-C1-C2-O2
2	B	601	HEM	C4C-C3C-CAC-CBC
4	B	603	NXR	C23-C24-C25-C30
4	B	603	NXR	C23-C24-C25-C26
4	B	603	NXR	N31-C23-C24-C25
2	A	601	HEM	C4C-C3C-CAC-CBC
4	A	604	NXR	C02-C13-C14-C15

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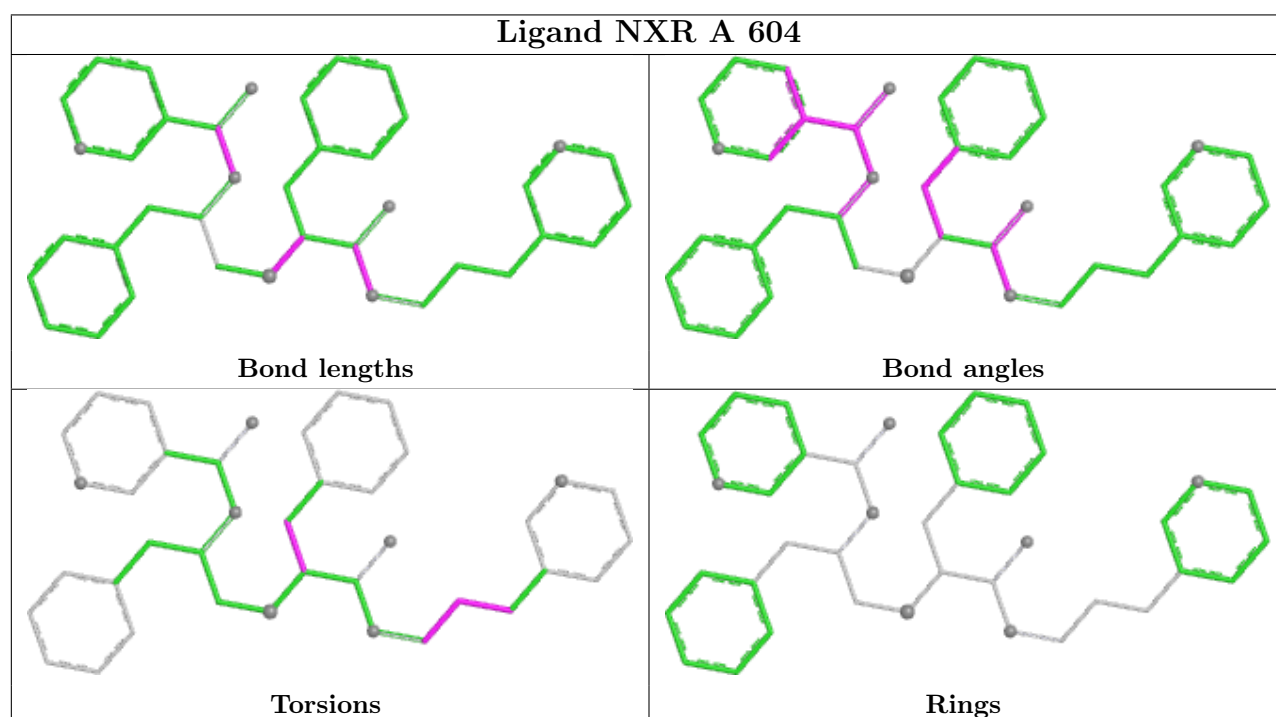
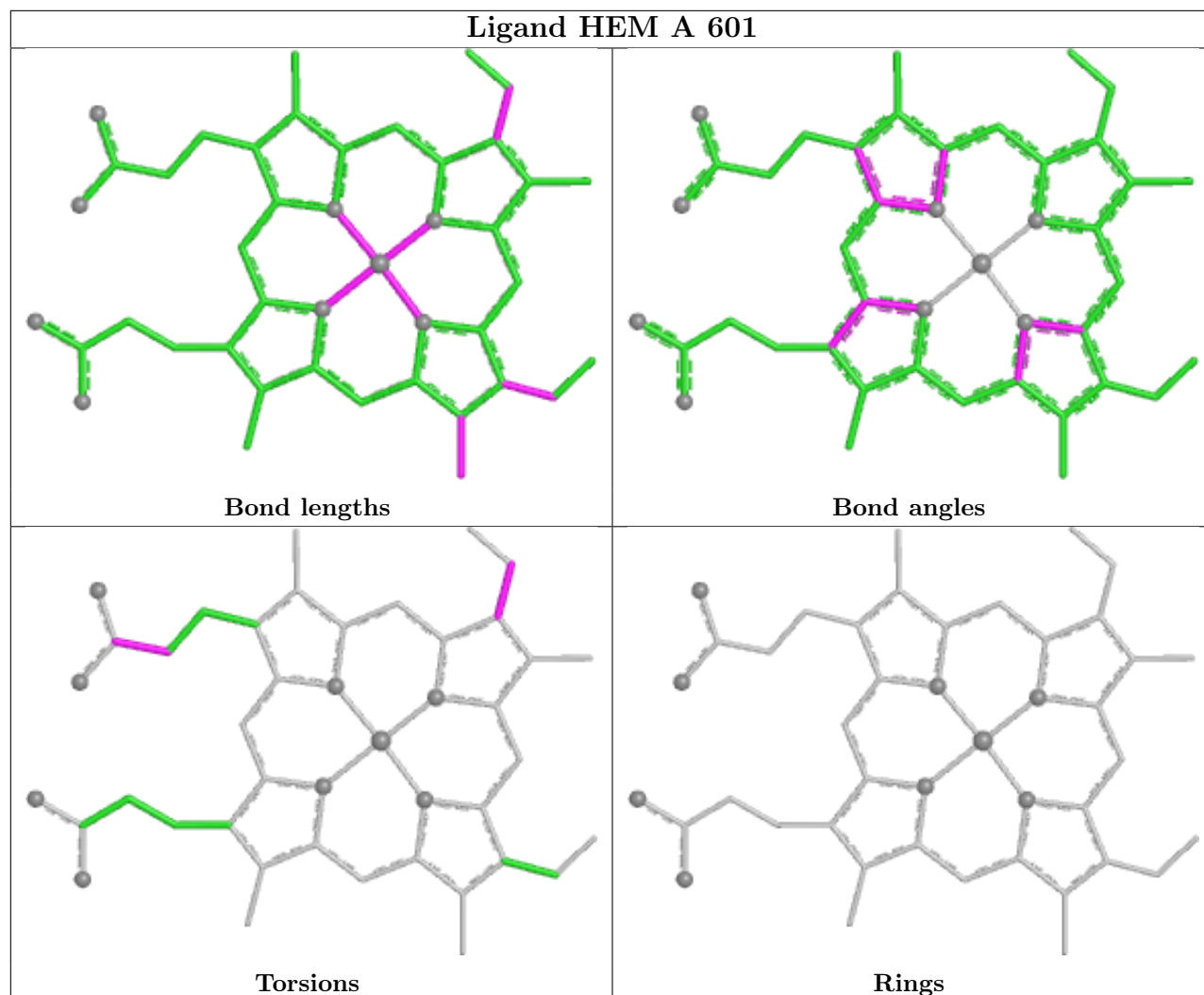
Mol	Chain	Res	Type	Atoms
4	B	603	NXR	C04-C05-C06-C07
2	A	601	HEM	CAD-CBD-CGD-O2D

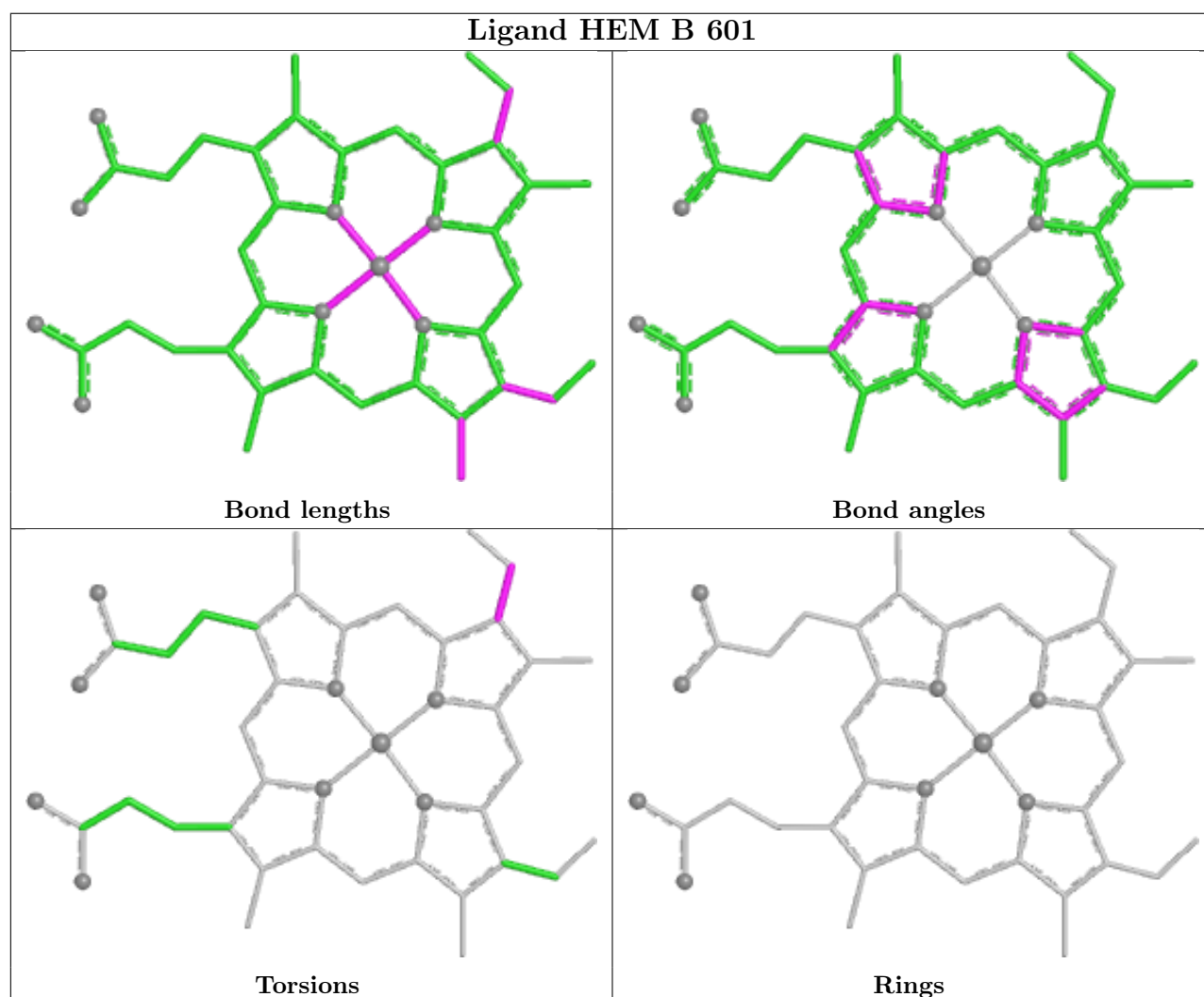
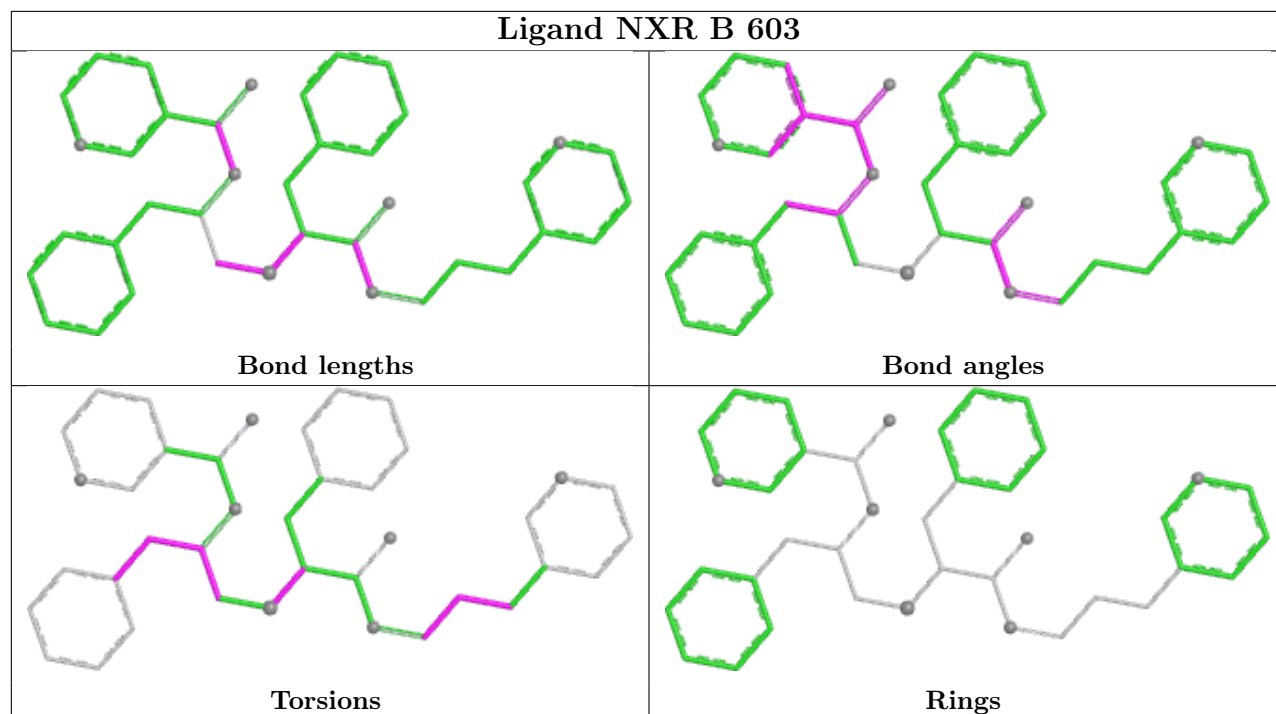
There are no ring outliers.

5 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	HEM	5	0
3	A	603	GOL	1	0
3	B	602	GOL	2	0
4	B	603	NXR	1	0
2	B	601	HEM	4	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	460/487 (94%)	0.62	25 (5%) 31 28	43, 85, 144, 182	2 (0%)
1	B	446/487 (91%)	0.81	46 (10%) 12 9	54, 108, 164, 198	0
All	All	906/974 (93%)	0.72	71 (7%) 19 15	43, 95, 160, 198	2 (0%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	210	LEU	5.6
1	B	221	LEU	5.3
1	A	226	PHE	5.1
1	B	229	LEU	4.9
1	B	226	PHE	4.5
1	A	215	PHE	4.4
1	B	160	LEU	4.3
1	B	108	PHE	4.2
1	B	260	ARG	3.7
1	A	211	LEU	3.7
1	B	232	ILE	3.7
1	A	225	VAL	3.5
1	A	465	PHE	3.4
1	A	479	LEU	3.3
1	B	156	LEU	3.3
1	B	220	PHE	3.2
1	B	153	GLY	3.2
1	B	223	ILE	3.1
1	B	219	PHE	3.1
1	A	240	VAL	3.1
1	B	240	VAL	3.1
1	B	256	MET	3.0
1	A	241	PHE	2.9
1	B	351	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	213	PHE	2.9
1	A	278	SER	2.8
1	B	211	LEU	2.8
1	B	235	VAL	2.8
1	A	26	GLY	2.7
1	B	190	GLY	2.7
1	B	228	PHE	2.7
1	B	191	VAL	2.7
1	A	108	PHE	2.6
1	B	267	HIS	2.6
1	B	159	ASN	2.6
1	B	209	LYS	2.6
1	B	179	TYR	2.6
1	B	218	PRO	2.6
1	A	228	PHE	2.6
1	B	119	SER	2.6
1	A	267	HIS	2.5
1	A	221	LEU	2.5
1	B	233	LEU	2.4
1	B	236	LEU	2.4
1	B	369	ILE	2.4
1	B	199	PRO	2.4
1	B	176	PHE	2.4
1	B	457	ILE	2.4
1	A	160	LEU	2.3
1	B	241	PHE	2.3
1	B	345	PRO	2.3
1	B	259	SER	2.3
1	A	316	PHE	2.3
1	A	158	ARG	2.3
1	A	424	ALA	2.2
1	B	352	GLN	2.2
1	B	407	TYR	2.2
1	B	210	LEU	2.2
1	B	239	CYS	2.2
1	A	288	LYS	2.2
1	B	450	MET	2.1
1	A	224	THR	2.1
1	A	223	ILE	2.1
1	A	407	TYR	2.1
1	B	225	VAL	2.1
1	A	261	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	69	GLY	2.1
1	B	295	LEU	2.1
1	B	278	SER	2.1
1	B	150	ALA	2.0
1	B	35	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

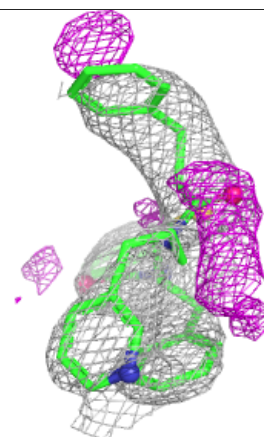
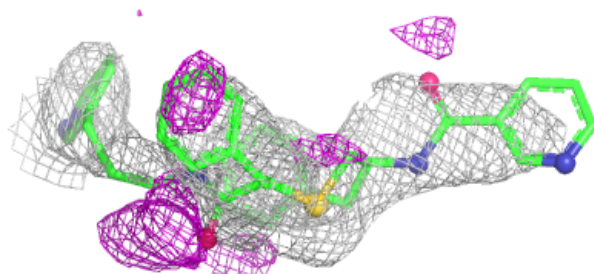
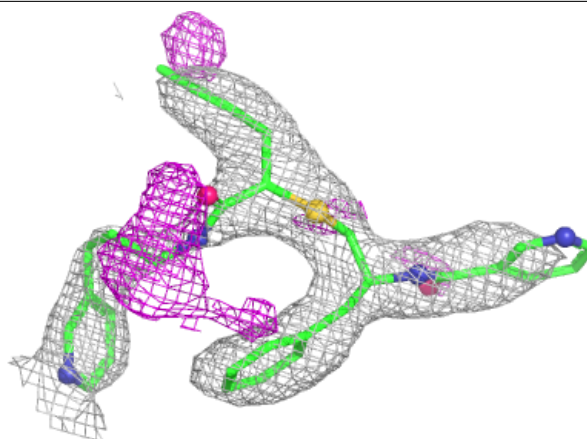
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	B	602	6/6	0.75	0.15	122,127,130,131	0
3	GOL	A	603	6/6	0.83	0.15	97,112,118,119	0
3	GOL	A	602	6/6	0.87	0.13	91,106,107,108	0
4	NXR	B	603	39/39	0.87	0.17	71,98,135,136	0
4	NXR	A	604	39/39	0.92	0.14	51,74,116,120	0
2	HEM	B	601	43/43	0.98	0.07	52,62,80,88	0
2	HEM	A	601	43/43	0.98	0.08	37,49,60,80	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.

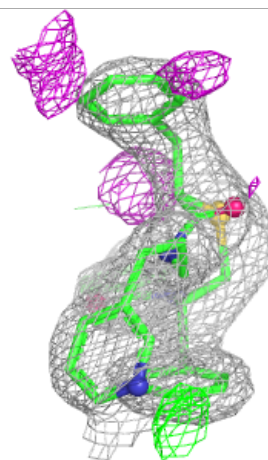
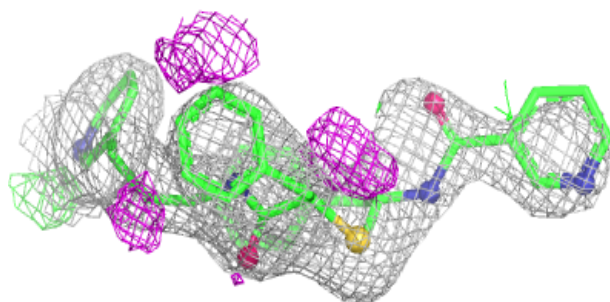
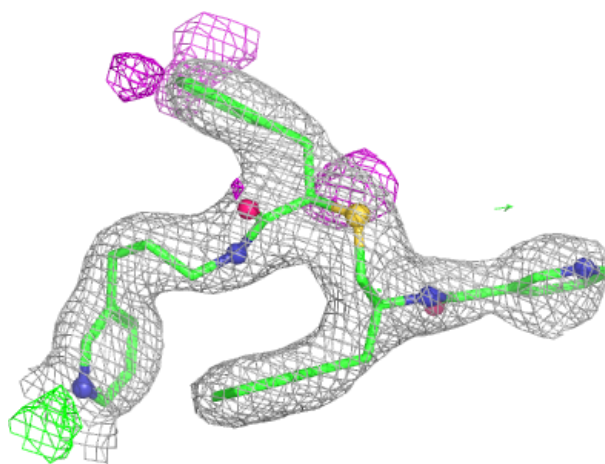
Electron density around NXR B 603:

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



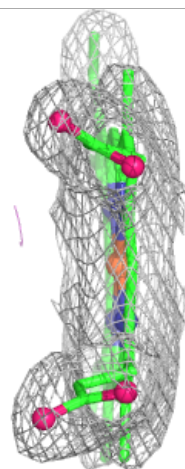
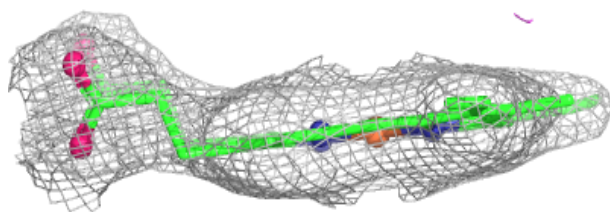
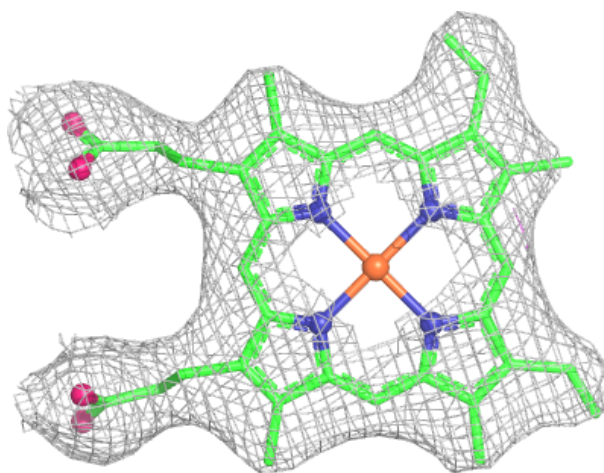
Electron density around NXR A 604:

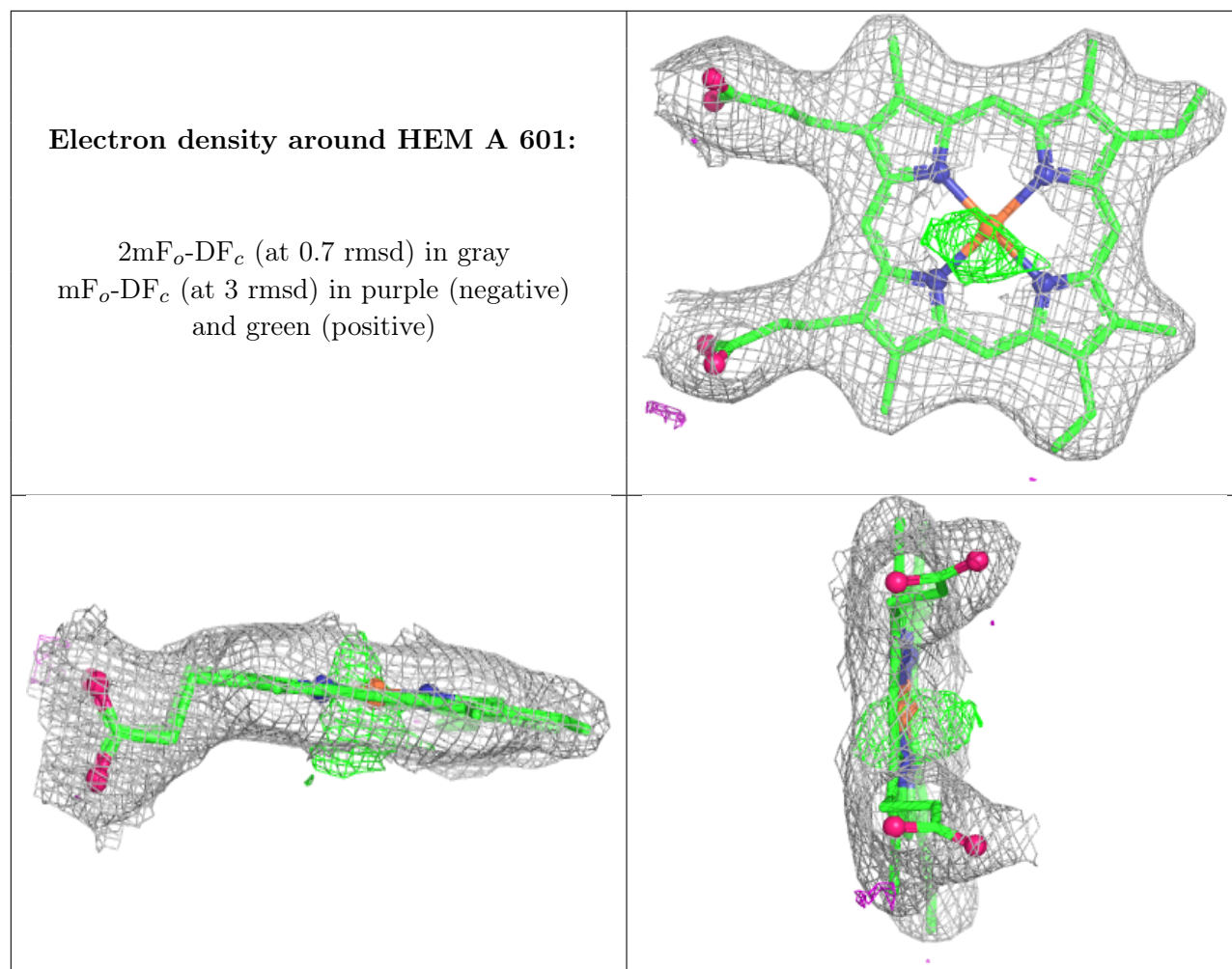
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM B 601:

$2mF_o-DF_c$ (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.