



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:30 PM UTC

PDB ID : 1HT5 / pdb_00001ht5
Title : THE 2.75 ANGSTROM RESOLUTION MODEL OF OVINE COX-1 COM-
PLEXED WITH METHYL ESTER FLURBIPROFEN
Authors : Selinsky, B.S.; Gupta, K.; Sharkey, C.T.; Loll, P.J.
Deposited on : 2000-12-28
Resolution : 2.75 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

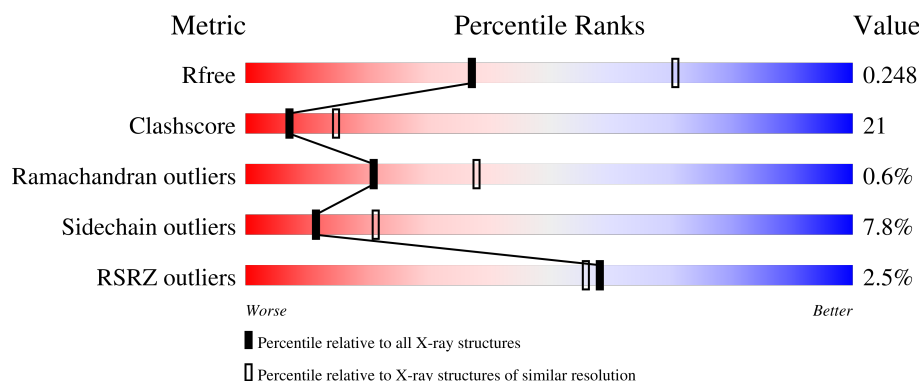
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

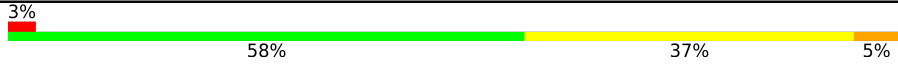
The reported resolution of this entry is 2.75 Å.

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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	551	
1	B	551	

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	NAG	A	1671	-	-	X	-
5	FL2	B	2701	-	-	X	-

2 Entry composition [i](#)

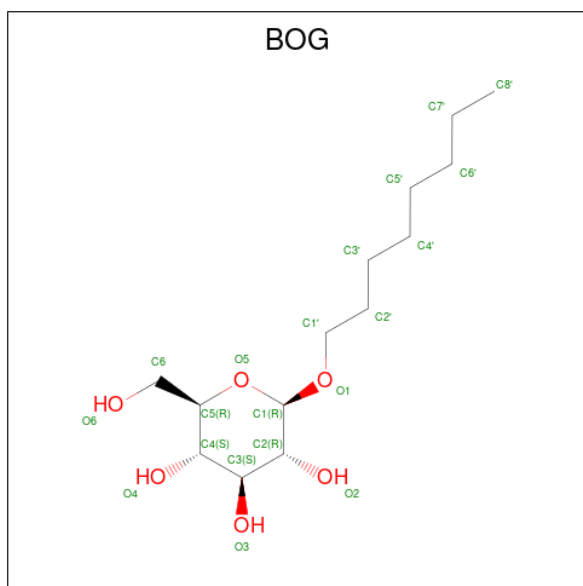
There are 6 unique types of molecules in this entry. The entry contains 9428 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROSTAGLANDIN H2 SYNTHASE-1.

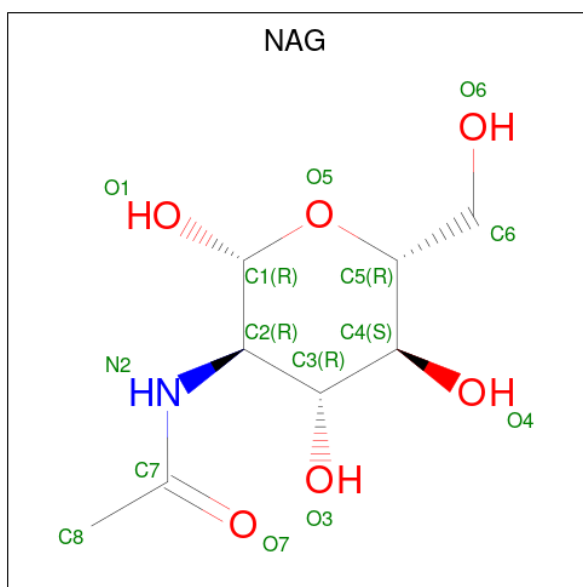
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	551	Total	C	N	O	S	0	0	0
			4477	2903	758	788	28			
1	B	551	Total	C	N	O	S	0	0	0
			4477	2903	758	788	28			

- Molecule 2 is octyl beta-D-glucopyranoside (CCD ID: BOG) (formula: $C_{14}H_{28}O_6$).



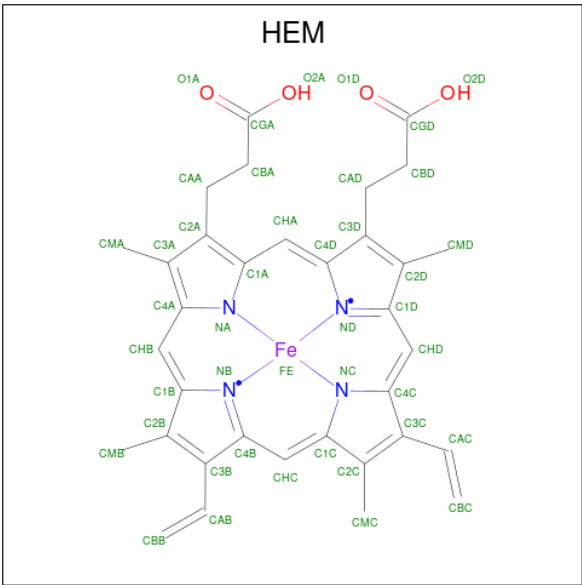
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			20	14	6		
2	B	1	Total	C	O	0	0
			20	14	6		
2	B	1	Total	C	O	0	0
			20	14	6		

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



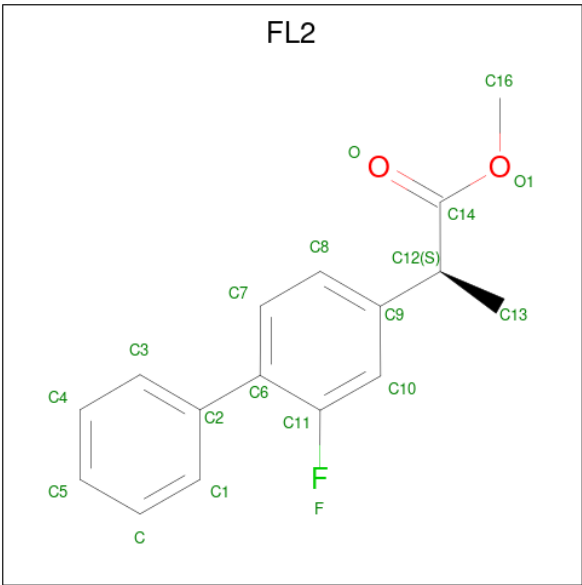
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

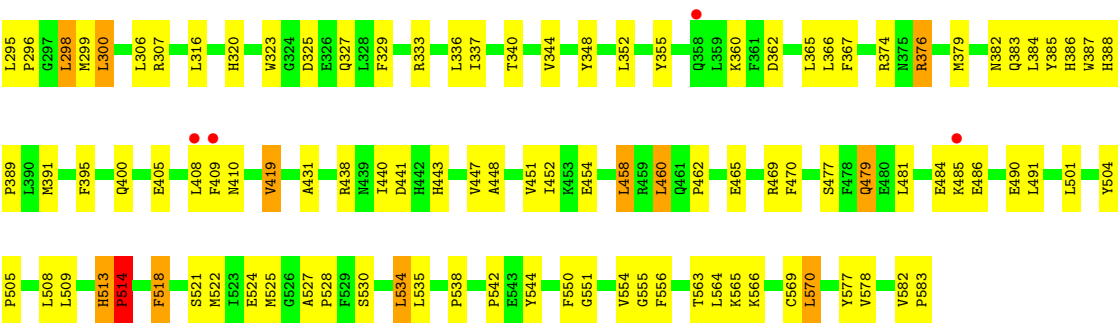
- Molecule 5 is FLURBIPROFEN METHYL ESTER (CCD ID: FL2) (formula: C₁₆H₁₅FO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	F	O	0	0
			19	16	1	2		
5	B	1	Total	C	F	O	0	0
			19	16	1	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	79	Total 79	O 79	0	0
6	B	71	Total 71	O 71	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	98.89Å 209.45Å 223.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.89 – 2.75 50.89 – 2.75	Depositor EDS
% Data completeness (in resolution range)	88.2 (50.89-2.75) 88.2 (50.89-2.75)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 2.69Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.223 , 0.252 0.215 , 0.248	Depositor DCC
R_{free} test set	5438 reflections (9.41%)	wwPDB-VP
Wilson B-factor (Å ²)	42.7	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9428	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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: FL2, BOG, NAG, HEM

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	0.45	0/4615	0.97	24/6264 (0.4%)
1	B	0.46	0/4615	0.99	17/6264 (0.3%)
All	All	0.45	0/9230	0.98	41/12528 (0.3%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	224	LEU	N-CA-C	-9.80	97.69	110.33
1	A	224	LEU	N-CA-C	-8.85	98.74	110.24
1	B	287	VAL	N-CA-C	8.24	121.11	113.20
1	A	287	VAL	N-CA-C	7.75	121.08	113.71
1	B	148	TYR	N-CA-C	-7.63	98.19	110.32
1	B	362	ASP	CA-C-N	7.49	128.16	119.47
1	B	362	ASP	C-N-CA	7.49	128.16	119.47
1	A	148	TYR	N-CA-C	-7.47	98.44	110.32
1	A	185	ARG	N-CA-C	-6.93	102.63	112.13
1	A	419	VAL	N-CA-C	6.63	117.13	110.23
1	A	569	CYS	N-CA-C	6.61	120.93	112.34
1	A	208	GLN	N-CA-C	-6.54	105.46	113.50
1	B	569	CYS	N-CA-C	6.27	119.78	111.75
1	A	356	PHE	N-CA-C	-6.23	105.64	113.18
1	A	197	MET	N-CA-C	-6.19	104.53	111.28
1	B	208	GLN	N-CA-C	-5.98	105.07	112.90
1	B	419	VAL	N-CA-C	5.96	116.14	110.42
1	A	467	ARG	N-CA-C	-5.91	104.75	111.07
1	A	269	ALA	CA-C-N	5.85	127.15	119.84
1	A	269	ALA	C-N-CA	5.85	127.15	119.84
1	A	408	LEU	N-CA-C	5.82	123.20	110.80
1	B	73	GLU	N-CA-C	-5.64	103.47	110.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	185	ARG	N-CA-C	-5.58	104.49	112.13
1	A	61	ARG	N-CA-C	5.57	119.25	111.90
1	B	238	LEU	N-CA-C	5.56	117.02	111.07
1	A	73	GLU	N-CA-C	-5.55	102.27	110.48
1	B	431	ALA	N-CA-C	-5.54	101.83	110.42
1	A	238	LEU	N-CA-C	5.52	116.98	111.07
1	A	71	ILE	CA-C-N	5.50	125.41	119.85
1	A	71	ILE	C-N-CA	5.50	125.41	119.85
1	A	45	GLY	N-CA-C	-5.45	105.11	112.14
1	B	251	LYS	N-CA-C	5.41	118.53	110.52
1	A	251	LYS	N-CA-C	5.32	118.39	110.52
1	A	457	VAL	N-CA-C	-5.26	105.59	110.53
1	B	337	ILE	N-CA-C	-5.22	105.41	110.42
1	A	135	ASP	N-CA-C	-5.21	105.16	112.25
1	B	513	HIS	N-CA-C	-5.19	102.48	110.01
1	A	292	PHE	N-CA-C	5.13	119.03	112.87
1	B	180	ARG	N-CA-C	5.12	116.94	111.36
1	B	486	GLU	N-CA-C	5.09	116.52	110.97
1	A	580	PHE	N-CA-C	-5.05	106.90	113.16

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	4477	0	4386	201	0
1	B	4477	0	4386	185	0
2	A	20	0	28	8	0
2	B	40	0	56	6	0
3	A	70	0	65	11	0
3	B	70	0	65	9	0
4	A	43	0	30	5	0
4	B	43	0	30	8	0
5	A	19	0	15	4	0
5	B	19	0	15	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	79	0	0	0	0
6	B	71	0	0	0	0
All	All	9428	0	9076	380	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 21.

All (380) 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:B:144:ASN:HD21	3:B:2671:NAG:C1	1.80	0.94
1:B:410:ASN:HD21	3:B:2681:NAG:C1	1.81	0.94
1:A:391:MET:HG3	4:A:601:HEM:HAB	1.47	0.92
1:A:163:MET:HE3	1:A:171:LEU:HD21	1.51	0.90
1:B:185:ARG:HH21	1:B:438:ARG:NH1	1.70	0.90
1:A:410:ASN:HD21	3:A:1681:NAG:C1	1.85	0.90
1:B:291:VAL:HG13	1:B:294:LEU:HD12	1.54	0.88
1:A:144:ASN:HD21	3:A:1671:NAG:C1	1.87	0.87
1:A:400:GLN:HE21	1:A:401:ASP:N	1.74	0.85
1:B:180:ARG:HD3	1:B:490:GLU:OE2	1.77	0.83
1:A:120:ARG:HG2	5:A:1701:FL2:H161	1.59	0.83
1:B:294:LEU:HD22	1:B:409:PHE:CE2	2.12	0.83
1:B:49:ARG:HG3	1:B:49:ARG:HH11	1.46	0.81
1:A:513:HIS:HE2	1:A:520:GLU:HB3	1.47	0.80
1:B:185:ARG:HH21	1:B:438:ARG:HH11	1.30	0.80
1:B:563:THR:HG22	1:B:565:LYS:H	1.48	0.78
1:B:295:LEU:HD12	1:B:298:LEU:HD22	1.63	0.78
1:A:140:GLU:CD	1:A:144:ASN:HD22	1.91	0.77
1:B:185:ARG:NH2	1:B:438:ARG:HH11	1.82	0.77
1:A:400:GLN:HE21	1:A:401:ASP:H	1.32	0.77
1:B:144:ASN:ND2	3:B:2671:NAG:C1	2.48	0.77
1:B:295:LEU:HD11	4:B:601:HEM:HBB2	1.65	0.76
1:B:454:GLU:HG2	1:B:458:LEU:HD22	1.68	0.75
1:A:144:ASN:ND2	3:A:1671:NAG:C1	2.49	0.75
1:A:400:GLN:NE2	1:A:401:ASP:H	1.85	0.74
1:B:208:GLN:NE2	1:B:230:LEU:H	1.84	0.74
1:B:150:ARG:HD3	1:B:152:LEU:O	1.87	0.73
1:A:582:VAL:HG22	1:A:583:PRO:HD2	1.71	0.71
1:B:113:MET:HE3	1:B:117:LEU:HD13	1.72	0.71
1:B:243:GLN:HG3	1:B:270:PRO:HD2	1.72	0.71
1:A:563:THR:HG22	1:A:565:LYS:H	1.55	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:ILE:HD12	2:A:1802:BOG:H62	1.73	0.71
1:A:208:GLN:NE2	1:A:230:LEU:H	1.89	0.70
1:A:563:THR:HG22	1:A:565:LYS:N	2.06	0.70
1:B:272:LEU:HD13	1:B:273:MET:N	2.07	0.69
1:A:113:MET:HE3	1:A:117:LEU:HD13	1.73	0.69
1:B:389:PRO:HG3	1:B:440:ILE:HG12	1.74	0.69
1:A:165:THR:HG22	1:A:166:LYS:HD3	1.74	0.69
1:A:462:PRO:HG2	1:A:465:GLU:HG2	1.74	0.69
1:A:374:ARG:NH1	1:B:374:ARG:HB3	2.08	0.69
1:B:150:ARG:HB3	1:B:379:MET:HE3	1.75	0.68
1:B:237:ASN:ND2	1:B:240:ARG:H	1.92	0.68
1:A:295:LEU:HD11	4:A:601:HEM:HBB2	1.74	0.68
1:B:120:ARG:HG2	5:B:2701:FL2:H161	1.74	0.68
1:A:67:PRO:C	1:A:68:ASN:HD22	2.02	0.68
1:B:295:LEU:CD1	4:B:601:HEM:HBB2	2.24	0.67
1:A:114:ARG:HD3	1:A:365:LEU:O	1.95	0.66
1:B:116:VAL:HG13	5:B:2701:FL2:H162	1.77	0.66
1:B:563:THR:HG22	1:B:565:LYS:N	2.10	0.66
1:A:91:PHE:O	1:A:95:HIS:HD2	1.78	0.66
1:A:447:VAL:O	1:A:451:VAL:HG23	1.96	0.66
1:B:273:MET:HE2	1:B:285:MET:O	1.95	0.65
1:A:276:PRO:HD2	1:A:279:ILE:HD12	1.78	0.65
1:B:185:ARG:HG2	1:B:185:ARG:HH11	1.61	0.65
1:A:280:PRO:HB2	1:A:282:GLN:OE1	1.96	0.65
1:A:344:VAL:O	1:A:348:TYR:HB3	1.97	0.65
1:B:120:ARG:NH1	2:B:2802:BOG:H1	2.12	0.64
1:A:504:TYR:HB3	1:A:505:PRO:HD3	1.79	0.64
1:A:237:ASN:ND2	1:A:239:GLU:HG2	2.13	0.64
1:A:320:HIS:HD2	1:B:49:ARG:O	1.81	0.64
1:A:150:ARG:HD3	1:A:152:LEU:O	1.98	0.63
1:A:482:THR:O	1:A:484:GLU:N	2.31	0.63
1:A:208:GLN:HE21	1:A:230:LEU:H	1.47	0.63
1:A:527:ALA:HB3	1:A:528:PRO:HD3	1.81	0.63
1:A:513:HIS:NE2	1:A:520:GLU:HB3	2.14	0.63
1:B:123:LEU:HD23	1:B:470:PHE:HD1	1.64	0.62
1:A:49:ARG:HG3	1:A:49:ARG:HH11	1.63	0.62
1:B:391:MET:HE2	4:B:601:HEM:HBB1	1.81	0.62
1:B:504:TYR:HB3	1:B:505:PRO:HD3	1.81	0.62
1:B:86:PRO:HG3	2:B:2802:BOG:H61	1.82	0.62
1:A:483:GLY:H	1:A:511:LYS:HB3	1.65	0.62
1:B:447:VAL:O	1:B:451:VAL:HG23	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:410:ASN:ND2	3:B:2681:NAG:C1	2.60	0.61
1:A:294:LEU:HD22	1:A:409:PHE:CE2	2.36	0.61
1:B:49:ARG:HH11	1:B:49:ARG:CG	2.12	0.61
1:B:116:VAL:O	1:B:120:ARG:HB2	2.00	0.61
1:B:554:VAL:HG13	1:B:555:GLY:N	2.16	0.61
1:A:582:VAL:CG2	1:A:583:PRO:HD2	2.31	0.61
1:B:213:SER:OG	1:B:215:LYS:HG2	2.01	0.61
1:B:387:TRP:HB2	4:B:601:HEM:HAC	1.83	0.61
1:A:410:ASN:ND2	3:A:1681:NAG:C1	2.61	0.60
1:A:172:PRO:HB2	1:A:177:LEU:HD22	1.82	0.60
1:A:180:ARG:O	1:A:438:ARG:NH1	2.33	0.60
1:B:120:ARG:HE	5:B:2701:FL2:C16	2.15	0.60
1:A:116:VAL:O	1:A:120:ARG:HB2	2.02	0.60
1:B:276:PRO:HD2	1:B:279:ILE:HD12	1.84	0.60
1:B:566:LYS:O	1:B:570:LEU:HB2	2.02	0.59
1:A:306:LEU:HD23	1:A:306:LEU:C	2.26	0.59
1:B:185:ARG:CZ	1:B:438:ARG:HD3	2.32	0.59
1:B:237:ASN:ND2	1:B:239:GLU:HG2	2.17	0.59
1:A:387:TRP:HB2	4:A:601:HEM:HAC	1.83	0.59
1:A:230:LEU:HG	1:A:233:ILE:HD12	1.86	0.58
1:B:294:LEU:HD22	1:B:409:PHE:HE2	1.65	0.58
1:A:382:ASN:OD1	1:A:386:HIS:HE1	1.87	0.58
1:B:120:ARG:HE	5:B:2701:FL2:H161	1.69	0.58
1:B:156:PRO:HB2	1:B:159:CYS:SG	2.43	0.58
1:B:518:PHE:CD1	1:B:522:MET:HG2	2.37	0.58
1:B:294:LEU:O	1:B:295:LEU:HG	2.04	0.58
1:A:86:PRO:HG3	2:A:1802:BOG:H61	1.84	0.57
1:A:504:TYR:CZ	1:A:508:LEU:HD11	2.39	0.57
1:A:294:LEU:O	1:A:295:LEU:HG	2.04	0.57
1:B:582:VAL:CG2	1:B:583:PRO:HD2	2.34	0.57
1:B:116:VAL:CG1	5:B:2701:FL2:H162	2.34	0.57
1:B:185:ARG:NE	1:B:438:ARG:HD3	2.19	0.57
1:A:293:GLY:HA2	1:A:299:MET:HE2	1.87	0.57
1:A:518:PHE:CD1	1:A:522:MET:HG2	2.40	0.57
1:A:200:PHE:CE1	1:A:391:MET:HE1	2.39	0.57
1:B:89:ILE:HD12	2:B:2802:BOG:H62	1.87	0.57
1:A:102:PHE:O	1:A:106:THR:HG23	2.05	0.56
1:B:389:PRO:HG3	1:B:440:ILE:CG1	2.35	0.56
1:A:323:TRP:CE3	1:A:327:GLN:HG2	2.41	0.56
1:A:389:PRO:HG3	1:A:440:ILE:HG12	1.86	0.56
1:A:49:ARG:O	1:B:320:HIS:HD2	1.89	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:479:GLN:CD	1:B:479:GLN:H	2.13	0.56
1:A:295:LEU:HD11	4:A:601:HEM:CBB	2.35	0.56
1:A:183:LEU:HD13	1:A:184:ARG:O	2.07	0.55
1:A:379:MET:SD	1:A:458:LEU:HG	2.46	0.55
1:A:200:PHE:CZ	1:A:391:MET:HE1	2.41	0.55
1:B:366:LEU:HD12	1:B:535:LEU:HD12	1.86	0.55
1:A:275:TYR:CE2	1:A:284:GLN:HB3	2.41	0.55
1:A:120:ARG:NH1	2:A:1802:BOG:H1	2.21	0.55
1:A:162:PRO:O	1:A:163:MET:HE2	2.07	0.55
1:B:77:TRP:CZ2	2:B:2801:BOG:H7'2	2.41	0.55
1:B:132:ILE:HD13	1:B:458:LEU:HD12	1.89	0.55
1:B:462:PRO:HG2	1:B:465:GLU:HG2	1.89	0.55
1:B:140:GLU:CD	1:B:144:ASN:HD22	2.15	0.54
1:A:88:PHE:CZ	1:A:92:LEU:HD11	2.43	0.54
1:A:208:GLN:NE2	1:A:228:VAL:HA	2.23	0.54
1:A:384:LEU:C	1:A:384:LEU:HD23	2.32	0.54
1:A:504:TYR:OH	1:A:508:LEU:HD11	2.08	0.54
1:A:524:GLU:OE1	2:A:1802:BOG:H5	2.08	0.54
1:A:197:MET:HE3	1:A:301:TYR:OH	2.08	0.54
1:A:388:HIS:N	1:A:389:PRO:CD	2.71	0.54
1:B:527:ALA:HB3	1:B:528:PRO:HD3	1.90	0.54
1:A:240:ARG:NH1	1:A:271:VAL:HG13	2.22	0.53
1:A:563:THR:HB	1:A:566:LYS:HG3	1.89	0.53
1:B:577:TYR:CE2	1:B:583:PRO:HD3	2.44	0.53
1:A:374:ARG:HH12	1:B:374:ARG:HB3	1.71	0.53
1:A:482:THR:C	1:A:484:GLU:H	2.17	0.53
1:A:554:VAL:HG13	1:A:555:GLY:N	2.23	0.53
1:A:179:ARG:HH11	1:A:179:ARG:HG2	1.74	0.52
1:A:306:LEU:HD23	1:A:306:LEU:O	2.09	0.52
1:A:352:LEU:HD11	1:A:387:TRP:HH2	1.74	0.52
1:B:340:THR:O	1:B:344:VAL:HG23	2.08	0.52
1:B:550:PHE:O	1:B:555:GLY:HA3	2.08	0.52
1:A:140:GLU:OE2	1:A:144:ASN:ND2	2.42	0.52
1:B:123:LEU:HD23	1:B:470:PHE:CD1	2.43	0.52
1:B:208:GLN:NE2	1:B:228:VAL:HA	2.25	0.52
1:A:34:ASN:HB3	1:A:37:CYS:SG	2.49	0.52
1:A:294:LEU:HD22	1:A:409:PHE:CD2	2.45	0.52
1:B:63:GLY:O	1:B:73:GLU:HG3	2.10	0.52
1:A:144:ASN:OD1	3:A:1671:NAG:C1	2.58	0.52
3:A:1671:NAG:H61	3:B:1672:NAG:HN2	1.75	0.52
1:B:293:GLY:HA2	1:B:299:MET:CE	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:ARG:NH2	1:B:438:ARG:NH1	2.43	0.51
1:B:208:GLN:HE21	1:B:230:LEU:H	1.56	0.51
1:B:477:SER:OG	1:B:479:GLN:HG2	2.10	0.51
1:A:180:ARG:HH21	1:A:490:GLU:CD	2.18	0.51
1:A:180:ARG:HD3	1:A:490:GLU:OE2	2.10	0.51
1:A:213:SER:OG	1:A:215:LYS:HG2	2.11	0.51
1:B:114:ARG:HD3	1:B:365:LEU:O	2.09	0.51
1:B:223:ALA:C	1:B:224:LEU:O	2.46	0.51
1:B:293:GLY:HA2	1:B:299:MET:HE2	1.92	0.51
1:B:391:MET:HE2	4:B:601:HEM:CBB	2.40	0.51
1:B:49:ARG:CG	1:B:49:ARG:NH1	2.71	0.51
1:A:203:GLN:HA	4:A:601:HEM:HBC2	1.93	0.50
1:B:142:PHE:C	1:B:376:ARG:HH22	2.19	0.50
1:B:344:VAL:O	1:B:348:TYR:HB3	2.11	0.50
1:B:491:LEU:HD11	1:B:509:LEU:HD13	1.92	0.50
1:A:269:ALA:O	1:A:271:VAL:N	2.45	0.50
1:B:513:HIS:O	1:B:514:PRO:C	2.53	0.50
1:A:144:ASN:CG	3:A:1671:NAG:C1	2.83	0.50
1:B:140:GLU:OE2	1:B:144:ASN:ND2	2.40	0.50
1:B:300:LEU:HD21	1:B:419:VAL:HG13	1.94	0.50
1:B:163:MET:HE2	1:B:163:MET:HA	1.93	0.50
1:A:123:LEU:O	1:A:469:ARG:NH2	2.45	0.50
1:B:126:SER:HA	1:B:127:PRO:C	2.35	0.50
1:B:513:HIS:HB3	1:B:514:PRO:HD2	1.93	0.50
1:A:403:SER:OG	1:A:405:GLU:HG2	2.12	0.49
1:B:35:PRO:HB2	1:B:55:TYR:CD2	2.47	0.49
1:B:582:VAL:HG23	1:B:583:PRO:HD2	1.94	0.49
1:A:400:GLN:HE21	1:A:400:GLN:CA	2.25	0.49
1:A:400:GLN:HE21	1:A:400:GLN:C	2.19	0.49
1:A:482:THR:C	1:A:484:GLU:N	2.71	0.49
1:B:49:ARG:HG3	1:B:49:ARG:NH1	2.23	0.49
1:A:291:VAL:HG12	1:A:291:VAL:O	2.12	0.49
1:A:77:TRP:CZ2	1:A:81:THR:HG21	2.48	0.49
1:A:352:LEU:HD11	1:A:387:TRP:CH2	2.48	0.49
1:B:295:LEU:HD12	1:B:298:LEU:CD2	2.36	0.49
1:A:117:LEU:HD12	1:A:531:LEU:HD13	1.95	0.48
1:A:179:ARG:HG2	1:A:179:ARG:NH1	2.28	0.48
1:B:391:MET:HG3	4:B:601:HEM:HAB	1.95	0.48
1:A:151:ILE:CG2	1:A:469:ARG:NH1	2.76	0.48
1:B:150:ARG:NH1	1:B:154:SER:HA	2.27	0.48
1:B:162:PRO:O	1:B:163:MET:HE2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:GLN:OE1	1:B:243:GLN:HA	2.13	0.48
1:B:291:VAL:HG12	1:B:291:VAL:O	2.13	0.48
1:A:49:ARG:HH11	1:A:49:ARG:CG	2.25	0.48
1:A:261:VAL:O	1:A:307:ARG:NH1	2.41	0.48
1:B:386:HIS:CE1	4:B:601:HEM:HAD2	2.47	0.48
1:A:35:PRO:CG	1:A:55:TYR:HB3	2.44	0.48
1:A:68:ASN:HD22	1:A:68:ASN:N	2.11	0.48
1:B:388:HIS:CE1	1:B:447:VAL:HG11	2.49	0.48
1:A:88:PHE:O	1:A:91:PHE:HB3	2.13	0.48
1:A:491:LEU:HD11	1:A:509:LEU:HD13	1.95	0.48
1:A:497:ASP:OD2	1:A:498:ILE:N	2.47	0.48
1:B:388:HIS:N	1:B:389:PRO:CD	2.76	0.48
1:A:84:PRO:HG2	1:A:89:ILE:HD11	1.95	0.48
1:B:70:THR:O	1:B:72:PRO:HD3	2.14	0.48
1:A:374:ARG:HB3	1:B:374:ARG:NH1	2.29	0.48
1:A:160:PRO:HD2	1:A:164:GLY:O	2.13	0.48
1:B:116:VAL:HG13	5:B:2701:FL2:C16	2.43	0.48
1:A:134:HIS:HD2	1:A:138:SER:OG	1.96	0.47
1:A:86:PRO:HA	2:A:1802:BOG:H61	1.96	0.47
1:B:99:LEU:O	1:B:103:VAL:HG23	2.14	0.47
1:A:391:MET:HE3	1:A:395:PHE:CE2	2.49	0.47
1:A:400:GLN:NE2	1:A:401:ASP:N	2.48	0.47
1:B:352:LEU:HD11	1:B:387:TRP:CH2	2.48	0.47
1:A:183:LEU:HD13	1:A:183:LEU:C	2.39	0.47
1:B:291:VAL:O	1:B:291:VAL:CG1	2.61	0.47
1:A:120:ARG:HE	5:A:1701:FL2:C16	2.27	0.47
1:A:120:ARG:HG2	5:A:1701:FL2:C16	2.37	0.47
1:A:479:GLN:HG3	1:A:485:LYS:HZ2	1.79	0.47
1:A:77:TRP:CE2	1:A:81:THR:HG21	2.49	0.47
1:A:142:PHE:CE2	1:B:538:PRO:HG3	2.49	0.47
1:A:513:HIS:HB3	1:A:514:PRO:HD2	1.96	0.47
1:B:240:ARG:NH1	1:B:271:VAL:HG13	2.30	0.47
1:B:384:LEU:C	1:B:384:LEU:HD23	2.39	0.47
1:A:145:VAL:HG13	1:A:145:VAL:O	2.15	0.47
1:A:340:THR:O	1:A:344:VAL:HG23	2.15	0.47
1:A:566:LYS:O	1:A:570:LEU:HB2	2.15	0.47
1:B:272:LEU:HD13	1:B:272:LEU:C	2.39	0.46
1:B:93:LEU:HD13	1:B:355:TYR:CE1	2.50	0.46
1:A:180:ARG:NH2	1:A:490:GLU:OE1	2.48	0.46
1:A:420:GLU:HG3	1:A:572:THR:HB	1.97	0.46
1:A:530:SER:O	1:A:534:LEU:HD22	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:ARG:NH2	1:B:360:LYS:HB2	2.31	0.46
1:A:234:TYR:CE2	1:A:333:ARG:HG3	2.51	0.46
1:A:294:LEU:HD22	1:A:409:PHE:HE2	1.81	0.46
1:A:320:HIS:CE1	1:A:551:GLY:O	2.69	0.46
1:A:452:ILE:O	1:A:455:SER:HB3	2.16	0.46
1:B:74:ILE:HG23	1:B:75:TRP:N	2.31	0.46
1:B:144:ASN:HD21	3:B:2671:NAG:C2	2.29	0.46
1:A:144:ASN:HD21	3:A:1671:NAG:C2	2.27	0.46
1:B:163:MET:HE3	1:B:171:LEU:HD21	1.98	0.46
1:A:96:GLY:O	1:A:97:ARG:C	2.57	0.46
1:B:530:SER:O	1:B:534:LEU:HD22	2.16	0.46
1:A:151:ILE:HG22	1:A:469:ARG:NH1	2.30	0.46
1:A:230:LEU:HD21	1:A:336:LEU:HB3	1.98	0.46
1:A:156:PRO:HD2	1:A:159:CYS:SG	2.56	0.45
1:B:382:ASN:OD1	1:B:386:HIS:HE1	1.98	0.45
1:B:538:PRO:HB2	1:B:544:TYR:CE1	2.51	0.45
1:B:183:LEU:HD22	1:B:184:ARG:N	2.30	0.45
1:B:206:THR:HB	1:B:210:PHE:CD2	2.51	0.45
1:A:306:LEU:C	1:A:306:LEU:CD2	2.90	0.45
1:A:162:PRO:C	1:A:163:MET:HE2	2.41	0.45
1:A:342:LYS:HG2	1:A:562:ALA:HB3	1.99	0.45
1:A:391:MET:HE3	1:A:395:PHE:HE2	1.81	0.45
1:A:513:HIS:HB2	1:A:516:SER:OG	2.16	0.45
1:A:520:GLU:HG3	1:A:521:SER:N	2.31	0.45
1:B:43:HIS:O	1:B:44:GLN:HB2	2.16	0.45
1:A:206:THR:C	1:A:208:GLN:H	2.25	0.45
1:A:513:HIS:O	1:A:514:PRO:C	2.60	0.45
1:A:367:PHE:CE2	1:A:542:PRO:HG3	2.52	0.45
1:B:256:MET:O	1:B:257:LEU:HD23	2.17	0.45
1:A:353:SER:HB2	1:A:355:TYR:CE2	2.52	0.45
1:B:102:PHE:O	1:B:106:THR:HG23	2.17	0.45
1:B:144:ASN:CG	3:B:2671:NAG:C1	2.90	0.45
1:B:290:GLU:H	1:B:290:GLU:CD	2.25	0.45
1:A:88:PHE:O	1:A:92:LEU:HD13	2.17	0.44
1:A:367:PHE:CD2	1:A:542:PRO:HG3	2.52	0.44
1:B:234:TYR:CE2	1:B:333:ARG:HG3	2.51	0.44
1:B:38:TYR:O	1:B:39:TYR:C	2.58	0.44
1:B:39:TYR:N	1:B:40:PRO:CD	2.79	0.44
1:B:323:TRP:CE3	1:B:327:GLN:HG2	2.52	0.44
1:B:504:TYR:CZ	1:B:508:LEU:HD11	2.53	0.44
1:B:521:SER:O	1:B:525:MET:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:VAL:CG1	5:A:1701:FL2:H162	2.47	0.44
1:A:120:ARG:HH12	1:A:524:GLU:CD	2.26	0.44
1:A:554:VAL:HG13	1:A:555:GLY:H	1.82	0.44
1:B:210:PHE:CE1	1:B:382:ASN:HA	2.52	0.44
1:A:479:GLN:CD	1:A:479:GLN:H	2.25	0.44
1:B:266:VAL:HG12	1:B:266:VAL:O	2.17	0.44
1:A:237:ASN:HD22	1:A:239:GLU:HG2	1.83	0.44
1:B:408:LEU:HB3	1:B:409:PHE:CD1	2.53	0.44
1:B:77:TRP:CZ2	2:B:2801:BOG:H5'2	2.53	0.44
1:B:185:ARG:HG2	1:B:185:ARG:NH1	2.32	0.44
1:A:375:ASN:HD22	1:A:532:LYS:HG3	1.83	0.44
1:A:98:TRP:CG	1:A:99:LEU:N	2.86	0.43
1:B:384:LEU:HG	1:B:522:MET:SD	2.58	0.43
1:A:323:TRP:CD2	1:A:327:GLN:HG2	2.54	0.43
1:A:55:TYR:CD1	1:A:55:TYR:C	2.97	0.43
1:A:569:CYS:HA	1:A:572:THR:OG1	2.18	0.43
1:B:180:ARG:O	1:B:438:ARG:NH1	2.51	0.43
1:B:185:ARG:NH2	1:B:438:ARG:HD3	2.33	0.43
1:B:237:ASN:HD21	1:B:240:ARG:H	1.66	0.43
1:B:280:PRO:HA	1:B:281:PRO:HD3	1.91	0.43
1:B:306:LEU:C	1:B:306:LEU:HD23	2.44	0.43
1:B:130:TYR:HB3	1:B:134:HIS:O	2.18	0.43
1:A:159:CYS:HB3	1:A:164:GLY:O	2.18	0.43
1:B:74:ILE:HG23	1:B:75:TRP:CD1	2.53	0.43
1:B:91:PHE:O	1:B:95:HIS:HD2	2.01	0.43
1:A:74:ILE:HG23	1:A:75:TRP:N	2.34	0.43
1:A:322:THR:OG1	1:B:49:ARG:HG2	2.18	0.43
1:A:367:PHE:HE1	1:B:61:ARG:HH21	1.65	0.43
1:B:441:ASP:OD2	1:B:443:HIS:CD2	2.72	0.43
1:A:223:ALA:O	1:A:224:LEU:C	2.59	0.43
1:A:245:ARG:NH1	1:A:326:GLU:OE2	2.48	0.43
1:A:275:TYR:CD2	1:A:284:GLN:HG2	2.53	0.43
1:B:230:LEU:CD2	1:B:336:LEU:HB3	2.48	0.43
1:A:202:ALA:O	1:A:206:THR:HG23	2.18	0.43
1:A:246:LEU:O	1:A:246:LEU:HD23	2.19	0.43
1:A:290:GLU:CD	1:A:290:GLU:H	2.27	0.43
1:A:544:TYR:O	1:A:546:LYS:N	2.50	0.43
1:B:448:ALA:O	1:B:452:ILE:HG13	2.19	0.43
1:A:68:ASN:OD1	3:A:1661:NAG:C1	2.67	0.43
1:A:454:GLU:HG2	1:A:458:LEU:HD22	2.01	0.43
1:B:230:LEU:HA	1:B:232:HIS:CE1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:479:GLN:HG3	1:A:485:LYS:NZ	2.34	0.42
1:B:352:LEU:HD22	1:B:518:PHE:CZ	2.53	0.42
1:A:391:MET:HB3	1:A:404:TYR:OH	2.18	0.42
1:B:554:VAL:CG1	1:B:555:GLY:N	2.80	0.42
1:A:458:LEU:HB3	1:A:460:LEU:HD13	2.01	0.42
1:A:86:PRO:CA	2:A:1802:BOG:H61	2.49	0.42
1:A:54:ARG:HG3	1:A:54:ARG:HH11	1.84	0.42
1:B:206:THR:HB	1:B:210:PHE:HD2	1.84	0.42
1:A:43:HIS:O	1:A:44:GLN:HB2	2.20	0.42
1:A:113:MET:HE3	1:A:117:LEU:CD1	2.47	0.42
1:A:287:VAL:HG11	1:A:302:ALA:HB1	2.00	0.42
1:A:49:ARG:CG	1:A:49:ARG:NH1	2.83	0.42
1:A:394:SER:HB2	1:A:402:TYR:O	2.20	0.42
1:B:250:GLY:HA2	1:B:325:ASP:OD1	2.19	0.42
1:B:367:PHE:CZ	1:B:542:PRO:HA	2.55	0.42
1:A:582:VAL:HG22	1:A:583:PRO:CD	2.46	0.42
1:A:206:THR:HB	1:A:210:PHE:CD2	2.55	0.42
1:A:295:LEU:HB2	1:A:298:LEU:HD22	2.01	0.42
3:A:1671:NAG:O4	3:B:1672:NAG:C1	2.68	0.42
1:B:504:TYR:OH	1:B:508:LEU:HD11	2.19	0.42
1:A:39:TYR:N	1:A:40:PRO:CD	2.83	0.41
1:B:172:PRO:HB2	1:B:177:LEU:HD22	2.02	0.41
1:B:280:PRO:HB2	1:B:282:GLN:CD	2.45	0.41
1:B:485:LYS:HD3	1:B:485:LYS:HA	1.90	0.41
1:B:144:ASN:OD1	3:B:2671:NAG:C1	2.68	0.41
1:A:64:TYR:CE1	1:A:76:THR:HG21	2.54	0.41
1:B:59:CYS:O	1:B:60:THR:C	2.64	0.41
1:B:79:ARG:HG2	1:B:79:ARG:HH11	1.85	0.41
1:A:196:LEU:HD21	1:A:392:PRO:HD3	2.02	0.41
1:A:262:TYR:HB3	1:A:285:MET:HE1	2.02	0.41
1:B:210:PHE:HB3	4:B:601:HEM:HBD1	2.02	0.41
1:B:481:LEU:HD22	1:B:501:LEU:CD2	2.50	0.41
1:A:251:LYS:HB3	1:A:310:ASN:OD1	2.21	0.41
1:B:55:TYR:CD1	1:B:55:TYR:C	2.99	0.41
1:B:145:VAL:HG13	1:B:145:VAL:O	2.19	0.41
1:B:173:ASP:OD2	1:B:175:GLU:HB3	2.21	0.41
1:B:320:HIS:HE1	1:B:551:GLY:O	2.04	0.41
1:A:344:VAL:O	1:A:349:VAL:HG23	2.21	0.41
1:B:176:PHE:CZ	1:B:180:ARG:HG3	2.56	0.41
1:B:238:LEU:HD23	1:B:238:LEU:HA	1.88	0.41
1:B:269:ALA:O	1:B:271:VAL:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ARG:NH2	1:A:490:GLU:CD	2.79	0.41
1:A:214:GLY:H	1:A:215:LYS:HZ1	1.68	0.41
1:B:240:ARG:NH1	1:B:290:GLU:HG3	2.36	0.41
1:B:292:PHE:CD1	1:B:298:LEU:HG	2.56	0.41
1:B:295:LEU:O	1:B:296:PRO:C	2.63	0.41
1:B:391:MET:CE	1:B:395:PHE:HE2	2.33	0.41
1:A:364:GLU:HG2	1:A:367:PHE:CE1	2.56	0.41
1:B:68:ASN:HD22	1:B:68:ASN:N	2.18	0.41
1:B:527:ALA:HB1	5:B:2701:FL2:H163	2.02	0.41
1:A:115:LEU:HD23	2:A:1802:BOG:H5'1	2.03	0.40
1:A:213:SER:HB3	1:A:216:MET:HB2	2.02	0.40
1:A:238:LEU:HA	1:A:238:LEU:HD23	1.80	0.40
1:A:91:PHE:O	1:A:95:HIS:CD2	2.65	0.40
1:B:383:GLN:HG3	1:B:460:LEU:HD21	2.04	0.40
1:A:320:HIS:HE1	1:A:551:GLY:O	2.04	0.40
1:A:406:GLN:HA	3:A:1681:NAG:C8	2.52	0.40
1:B:116:VAL:HG22	2:B:2802:BOG:H2'1	2.02	0.40
1:B:120:ARG:HH12	1:B:524:GLU:CD	2.29	0.40
1:B:245:ARG:HD2	1:B:329:PHE:CZ	2.57	0.40
1:A:86:PRO:CG	2:A:1802:BOG:H61	2.50	0.40
1:B:77:TRP:CE3	1:B:78:LEU:HD23	2.56	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	549/551 (100%)	509 (93%)	36 (7%)	4 (1%)	18	34
1	B	549/551 (100%)	504 (92%)	42 (8%)	3 (0%)	24	43
All	All	1098/1102 (100%)	1013 (92%)	78 (7%)	7 (1%)	21	38

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	514	PRO
1	B	514	PRO
1	A	483	GLY
1	A	270	PRO
1	B	290	GLU
1	A	157	ARG
1	B	270	PRO

5.3.2 Protein sidechains [i](#)

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	486/486 (100%)	452 (93%)	34 (7%)	14	26
1	B	486/486 (100%)	444 (91%)	42 (9%)	10	18
All	All	972/972 (100%)	896 (92%)	76 (8%)	11	21

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	ARG
1	A	117	LEU
1	A	120	ARG
1	A	145	VAL
1	A	165	THR
1	A	171	LEU
1	A	215	LYS
1	A	232	HIS
1	A	238	LEU
1	A	244	LEU
1	A	246	LEU
1	A	252	LEU
1	A	271	VAL
1	A	277	ARG
1	A	289	GLN
1	A	298	LEU

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Mol	Chain	Res	Type
1	A	300	LEU
1	A	307	ARG
1	A	316	LEU
1	A	376	ARG
1	A	385	TYR
1	A	400	GLN
1	A	405	GLU
1	A	438	ARG
1	A	458	LEU
1	A	460	LEU
1	A	469	ARG
1	A	479	GLN
1	A	484	GLU
1	A	514	PRO
1	A	518	PHE
1	A	534	LEU
1	A	564	LEU
1	A	578	VAL
1	B	33	VAL
1	B	49	ARG
1	B	61	ARG
1	B	70	THR
1	B	97	ARG
1	B	117	LEU
1	B	120	ARG
1	B	143	SER
1	B	145	VAL
1	B	146	SER
1	B	171	LEU
1	B	183	LEU
1	B	190	ASP
1	B	238	LEU
1	B	243	GLN
1	B	244	LEU
1	B	246	LEU
1	B	248	LYS
1	B	252	LEU
1	B	271	VAL
1	B	272	LEU
1	B	289	GLN
1	B	298	LEU
1	B	300	LEU

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Mol	Chain	Res	Type
1	B	307	ARG
1	B	316	LEU
1	B	376	ARG
1	B	385	TYR
1	B	400	GLN
1	B	405	GLU
1	B	458	LEU
1	B	460	LEU
1	B	469	ARG
1	B	479	GLN
1	B	484	GLU
1	B	514	PRO
1	B	518	PHE
1	B	534	LEU
1	B	556	PHE
1	B	564	LEU
1	B	570	LEU
1	B	578	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	HIS
1	A	68	ASN
1	A	95	HIS
1	A	134	HIS
1	A	144	ASN
1	A	208	GLN
1	A	237	ASN
1	A	255	GLN
1	A	320	HIS
1	A	358	GLN
1	A	370	GLN
1	A	375	ASN
1	A	386	HIS
1	A	400	GLN
1	A	443	HIS
1	A	446	HIS
1	A	479	GLN
1	A	571	ASN
1	B	42	GLN
1	B	68	ASN

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Mol	Chain	Res	Type
1	B	90	HIS
1	B	95	HIS
1	B	134	HIS
1	B	208	GLN
1	B	237	ASN
1	B	255	GLN
1	B	320	HIS
1	B	375	ASN
1	B	386	HIS
1	B	400	GLN
1	B	410	ASN
1	B	442	HIS
1	B	443	HIS
1	B	479	GLN
1	B	557	ASN

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](#)

17 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
5	FL2	B	2701	-	20,20,20	3.17	14 (70%)	27,27,27	1.30	4 (14%)
3	NAG	A	1681	-	14,14,15	1.70	3 (21%)	17,19,21	0.92	0
3	NAG	B	1672	-	14,14,15	1.51	4 (28%)	17,19,21	1.13	1 (5%)
3	NAG	A	2672	-	14,14,15	1.86	6 (42%)	17,19,21	1.12	1 (5%)
2	BOG	B	2802	-	20,20,20	1.55	5 (25%)	25,25,25	0.96	1 (4%)
3	NAG	A	1661	-	14,14,15	1.35	3 (21%)	17,19,21	1.02	1 (5%)
4	HEM	A	601	1	50,50,50	1.44	5 (10%)	67,82,82	1.11	5 (7%)
2	BOG	A	1802	-	20,20,20	1.58	5 (25%)	25,25,25	0.88	2 (8%)
3	NAG	B	2662	-	14,14,15	1.70	5 (35%)	17,19,21	1.02	1 (5%)
3	NAG	A	1671	-	14,14,15	1.18	1 (7%)	17,19,21	0.97	1 (5%)
2	BOG	B	2801	-	20,20,20	1.73	6 (30%)	25,25,25	1.08	1 (4%)
3	NAG	B	2671	-	14,14,15	1.10	1 (7%)	17,19,21	0.94	1 (5%)
3	NAG	A	1662	-	14,14,15	1.66	5 (35%)	17,19,21	1.00	1 (5%)
5	FL2	A	1701	-	20,20,20	3.15	12 (60%)	27,27,27	1.32	4 (14%)
3	NAG	B	2681	-	14,14,15	1.57	3 (21%)	17,19,21	1.03	1 (5%)
3	NAG	B	2661	-	14,14,15	1.40	2 (14%)	17,19,21	1.03	1 (5%)
4	HEM	B	601	1	50,50,50	1.43	5 (10%)	67,82,82	1.15	7 (10%)

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
5	FL2	B	2701	-	-	3/14/14/14	0/2/2/2
3	NAG	A	1681	-	-	4/6/23/26	0/1/1/1
3	NAG	B	1672	-	-	4/6/23/26	0/1/1/1
3	NAG	A	2672	-	-	4/6/23/26	0/1/1/1
2	BOG	B	2802	-	-	4/11/31/31	0/1/1/1
3	NAG	A	1661	-	-	2/6/23/26	0/1/1/1
4	HEM	A	601	1	-	4/14/54/54	-
2	BOG	A	1802	-	-	4/11/31/31	0/1/1/1
3	NAG	B	2662	-	-	2/6/23/26	0/1/1/1
3	NAG	A	1671	-	-	4/6/23/26	0/1/1/1
2	BOG	B	2801	-	-	1/11/31/31	0/1/1/1
3	NAG	B	2671	-	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	1662	-	-	2/6/23/26	0/1/1/1
5	FL2	A	1701	-	-	3/14/14/14	0/2/2/2
3	NAG	B	2681	-	-	4/6/23/26	0/1/1/1
3	NAG	B	2661	-	-	0/6/23/26	0/1/1/1
4	HEM	B	601	1	-	7/14/54/54	-

All (85) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1701	FL2	C7-C6	5.83	1.48	1.40
5	B	2701	FL2	C7-C6	5.73	1.48	1.40
5	A	1701	FL2	C8-C9	5.34	1.47	1.39
5	B	2701	FL2	C8-C9	5.10	1.47	1.39
5	A	1701	FL2	C1-C2	4.74	1.48	1.39
5	B	2701	FL2	C6-C11	4.63	1.47	1.39
4	B	601	HEM	CBC-CAC	4.62	1.52	1.30
4	A	601	HEM	CBB-CAB	4.60	1.52	1.30
4	A	601	HEM	CBC-CAC	4.46	1.51	1.30
4	B	601	HEM	CBB-CAB	4.44	1.51	1.30
5	B	2701	FL2	C1-C2	4.40	1.48	1.39
5	A	1701	FL2	C8-C7	4.36	1.45	1.38
5	A	1701	FL2	C6-C11	4.29	1.46	1.39
5	B	2701	FL2	C8-C7	4.13	1.45	1.38
5	A	1701	FL2	C4-C3	4.02	1.45	1.38
5	B	2701	FL2	C-C1	3.96	1.45	1.38
5	B	2701	FL2	C4-C3	3.96	1.45	1.38
2	B	2801	BOG	O5-C1	3.67	1.51	1.41
5	A	1701	FL2	C-C1	3.57	1.45	1.38
3	B	2662	NAG	C3-C2	3.42	1.59	1.52
3	A	2672	NAG	C1-C2	3.35	1.56	1.52
3	A	1681	NAG	O5-C5	3.34	1.49	1.43
3	A	1662	NAG	C3-C2	3.26	1.59	1.52
2	A	1802	BOG	O5-C1	3.24	1.50	1.41
2	B	2802	BOG	O5-C1	3.16	1.50	1.41
5	B	2701	FL2	C3-C2	3.14	1.45	1.39
5	A	1701	FL2	C3-C2	3.08	1.45	1.39
3	A	1661	NAG	C3-C2	2.96	1.58	1.52
3	B	2661	NAG	C3-C2	2.96	1.58	1.52
2	B	2801	BOG	O5-C5	2.92	1.51	1.44
3	A	1681	NAG	C3-C2	2.91	1.58	1.52
5	A	1701	FL2	C5-C4	2.88	1.44	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2681	NAG	O5-C5	2.87	1.49	1.43
2	B	2802	BOG	O5-C5	2.86	1.51	1.44
3	B	2681	NAG	C4-C5	2.85	1.59	1.53
5	A	1701	FL2	O-C14	2.84	1.28	1.21
5	B	2701	FL2	C5-C4	2.83	1.44	1.38
2	B	2801	BOG	C3'-C2'	-2.81	1.37	1.51
4	A	601	HEM	CMB-C2B	2.81	1.56	1.50
2	A	1802	BOG	C4-C5	-2.77	1.47	1.53
3	A	1681	NAG	C4-C5	2.73	1.58	1.53
3	A	1671	NAG	C4-C5	2.71	1.58	1.53
4	A	601	HEM	CAB-C3B	2.71	1.54	1.47
3	A	2672	NAG	O5-C5	2.71	1.48	1.43
3	B	1672	NAG	C3-C2	2.70	1.58	1.52
2	B	2802	BOG	C3'-C2'	-2.68	1.38	1.51
3	A	2672	NAG	C3-C2	2.68	1.58	1.52
3	B	2662	NAG	O5-C5	2.67	1.48	1.43
4	B	601	HEM	CAC-C3C	2.66	1.54	1.47
4	B	601	HEM	CMB-C2B	2.66	1.56	1.50
5	B	2701	FL2	O-C14	2.65	1.27	1.21
2	A	1802	BOG	C3'-C2'	-2.62	1.38	1.51
5	B	2701	FL2	C9-C12	2.62	1.57	1.52
2	B	2802	BOG	C4-C5	-2.57	1.47	1.53
3	A	2672	NAG	C2-N2	2.56	1.50	1.46
2	A	1802	BOG	O5-C5	2.52	1.50	1.44
3	B	2681	NAG	C3-C2	2.52	1.57	1.52
3	B	2671	NAG	C3-C2	2.50	1.57	1.52
4	A	601	HEM	CAC-C3C	2.50	1.54	1.47
3	B	2662	NAG	C1-C2	2.41	1.55	1.52
2	B	2801	BOG	C4-C5	-2.40	1.47	1.53
5	B	2701	FL2	C10-C9	2.40	1.42	1.39
3	A	1662	NAG	C4-C5	2.37	1.58	1.53
3	A	2672	NAG	C4-C3	2.36	1.58	1.52
3	A	1662	NAG	O5-C5	2.35	1.48	1.43
3	A	1661	NAG	C4-C5	2.32	1.58	1.53
2	B	2801	BOG	C4-C3	2.30	1.58	1.52
3	A	1662	NAG	C1-C2	2.30	1.55	1.52
3	B	2661	NAG	C4-C5	2.29	1.57	1.53
3	A	2672	NAG	C4-C5	2.29	1.57	1.53
4	B	601	HEM	CAB-C3B	2.28	1.53	1.47
3	B	1672	NAG	C4-C5	2.28	1.57	1.53
3	B	1672	NAG	O5-C5	2.24	1.47	1.43
5	B	2701	FL2	C5-C	2.23	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2801	BOG	C3-C2	2.21	1.58	1.52
2	A	1802	BOG	C7'-C6'	-2.21	1.38	1.51
2	B	2802	BOG	C7'-C6'	-2.17	1.38	1.51
3	B	1672	NAG	C4-C3	2.16	1.57	1.52
3	B	2662	NAG	C4-C5	2.13	1.57	1.53
5	B	2701	FL2	C10-C11	2.11	1.41	1.37
5	A	1701	FL2	C5-C	2.11	1.42	1.38
5	A	1701	FL2	C10-C9	2.09	1.42	1.39
3	B	2662	NAG	C4-C3	2.06	1.57	1.52
3	A	1662	NAG	C4-C3	2.04	1.57	1.52
3	A	1661	NAG	O5-C5	2.02	1.47	1.43

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2801	BOG	C1'-O1-C1	4.44	121.26	113.68
2	B	2802	BOG	C1'-O1-C1	3.79	120.15	113.68
4	A	601	HEM	C2B-C1B-NB	3.58	113.96	109.84
4	B	601	HEM	C2B-C1B-NB	3.39	113.73	109.84
5	B	2701	FL2	C16-O1-C14	2.79	122.25	115.92
5	A	1701	FL2	C16-O1-C14	2.78	122.23	115.92
2	A	1802	BOG	O1-C1-C2	2.76	112.46	108.27
5	A	1701	FL2	O-C14-C12	-2.75	109.95	122.70
5	B	2701	FL2	O-C14-C12	-2.73	110.03	122.70
4	B	601	HEM	C2D-C1D-ND	2.69	113.02	109.90
3	A	2672	NAG	O7-C7-C8	-2.69	117.27	122.05
4	B	601	HEM	CBB-CAB-C3B	-2.67	114.18	127.53
4	A	601	HEM	C3B-C4B-NB	2.55	111.30	109.47
3	B	1672	NAG	O7-C7-C8	-2.53	117.54	122.05
5	A	1701	FL2	C13-C12-C14	2.42	119.27	111.09
3	B	2662	NAG	O7-C7-C8	-2.38	117.82	122.05
2	A	1802	BOG	C1'-O1-C1	2.34	117.68	113.68
5	B	2701	FL2	C13-C12-C14	2.32	118.90	111.09
4	B	601	HEM	C3B-C2B-C1B	-2.31	104.68	106.41
3	A	1662	NAG	O7-C7-C8	-2.27	118.02	122.05
4	B	601	HEM	C1D-C2D-C3D	-2.22	104.64	106.98
4	A	601	HEM	CBB-CAB-C3B	-2.20	116.52	127.53
4	A	601	HEM	C3B-C2B-C1B	-2.20	104.76	106.41
3	B	2661	NAG	O7-C7-C8	-2.16	118.20	122.05
3	A	1671	NAG	C4-C3-C2	-2.15	107.87	111.02
3	B	2671	NAG	O7-C7-C8	-2.11	118.30	122.05
4	B	601	HEM	CBA-CAA-C2A	-2.11	106.70	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	2681	NAG	O7-C7-C8	-2.10	118.31	122.05
4	B	601	HEM	C3B-C4B-NB	2.09	110.97	109.47
5	B	2701	FL2	O1-C14-C12	2.08	129.04	114.23
3	A	1661	NAG	O7-C7-C8	-2.07	118.37	122.05
5	A	1701	FL2	O1-C14-C12	2.05	128.83	114.23
4	A	601	HEM	C1D-C2D-C3D	-2.05	104.83	106.98

There are no chirality outliers.

All (54) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1681	NAG	C8-C7-N2-C2
3	A	1681	NAG	O7-C7-N2-C2
4	A	601	HEM	C2C-C3C-CAC-CBC
4	B	601	HEM	C2B-C3B-CAB-CBB
4	B	601	HEM	C4B-C3B-CAB-CBB
5	A	1701	FL2	C12-C14-O1-C16
5	A	1701	FL2	O-C14-O1-C16
5	B	2701	FL2	C12-C14-O1-C16
5	B	2701	FL2	O-C14-O1-C16
3	B	1672	NAG	C8-C7-N2-C2
3	B	1672	NAG	O7-C7-N2-C2
3	A	2672	NAG	O5-C5-C6-O6
3	B	1672	NAG	O5-C5-C6-O6
3	A	1662	NAG	C8-C7-N2-C2
3	B	2662	NAG	C8-C7-N2-C2
3	B	2681	NAG	C4-C5-C6-O6
3	A	2672	NAG	C4-C5-C6-O6
3	B	1672	NAG	C4-C5-C6-O6
3	B	2681	NAG	O5-C5-C6-O6
3	A	1662	NAG	O7-C7-N2-C2
3	A	2672	NAG	C8-C7-N2-C2
3	B	2662	NAG	O7-C7-N2-C2
3	B	2681	NAG	C8-C7-N2-C2
3	A	1671	NAG	O5-C5-C6-O6
3	A	1671	NAG	C4-C5-C6-O6
2	B	2802	BOG	O1-C1'-C2'-C3'
3	A	1681	NAG	O5-C5-C6-O6
2	A	1802	BOG	O1-C1'-C2'-C3'
3	A	2672	NAG	O7-C7-N2-C2
3	B	2681	NAG	O7-C7-N2-C2
2	A	1802	BOG	C1'-C2'-C3'-C4'

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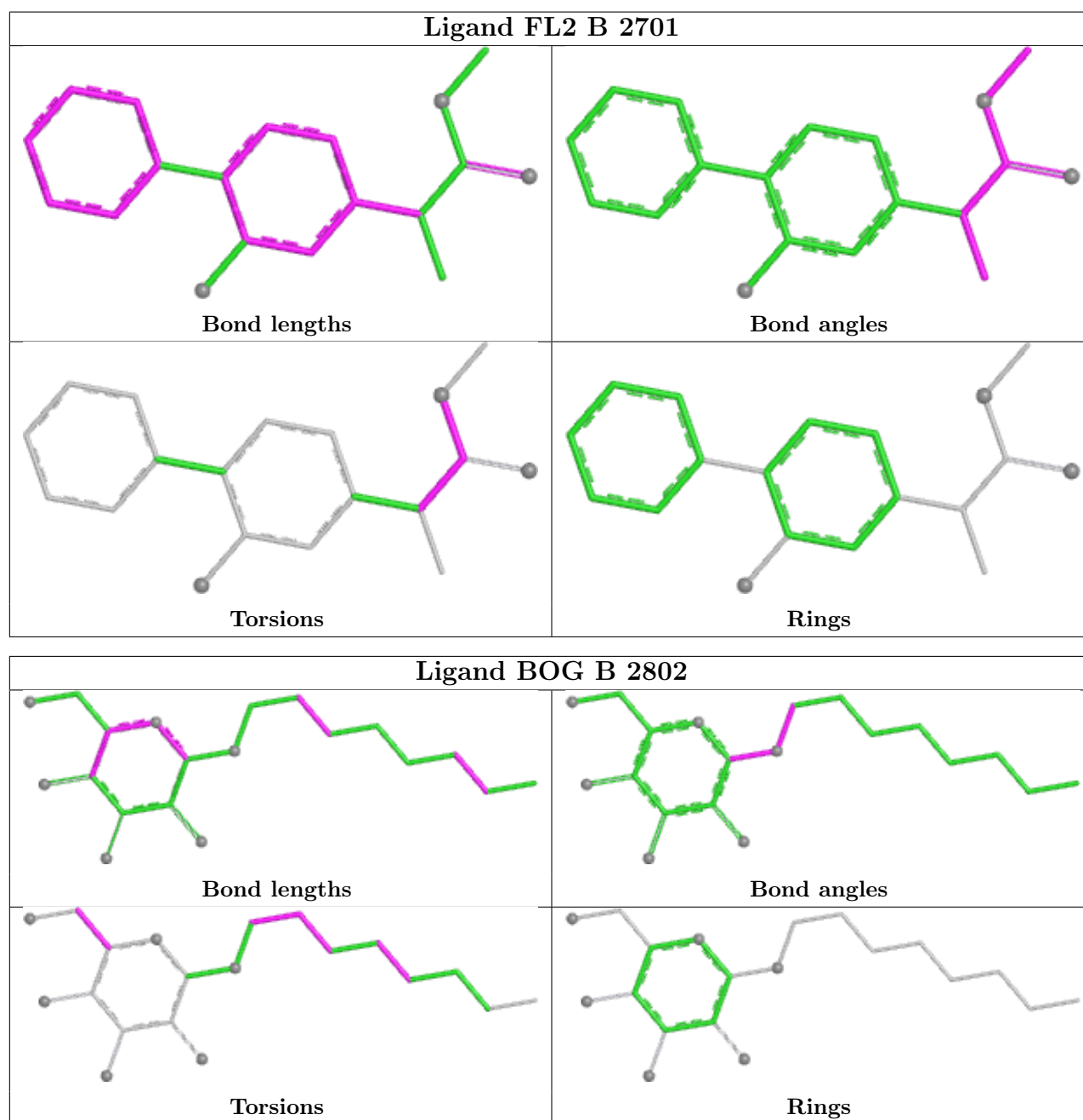
Mol	Chain	Res	Type	Atoms
4	B	601	HEM	C2C-C3C-CAC-CBC
2	B	2802	BOG	C1'-C2'-C3'-C4'
4	A	601	HEM	C4C-C3C-CAC-CBC
3	B	2671	NAG	C4-C5-C6-O6
2	B	2802	BOG	C3'-C4'-C5'-C6'
3	A	1661	NAG	C4-C5-C6-O6
3	B	2671	NAG	O5-C5-C6-O6
2	A	1802	BOG	C4-C5-C6-O6
2	A	1802	BOG	O5-C5-C6-O6
4	A	601	HEM	C2B-C3B-CAB-CBB
4	B	601	HEM	C4C-C3C-CAC-CBC
3	A	1671	NAG	C8-C7-N2-C2
3	A	1661	NAG	O5-C5-C6-O6
5	A	1701	FL2	C13-C12-C14-O1
5	B	2701	FL2	C13-C12-C14-O1
4	B	601	HEM	C3D-CAD-CBD-CGD
3	A	1671	NAG	O7-C7-N2-C2
2	B	2801	BOG	C3'-C4'-C5'-C6'
3	A	1681	NAG	C4-C5-C6-O6
4	A	601	HEM	C4B-C3B-CAB-CBB
4	B	601	HEM	CAD-CBD-CGD-O2D
2	B	2802	BOG	C4-C5-C6-O6
4	B	601	HEM	CAD-CBD-CGD-O1D

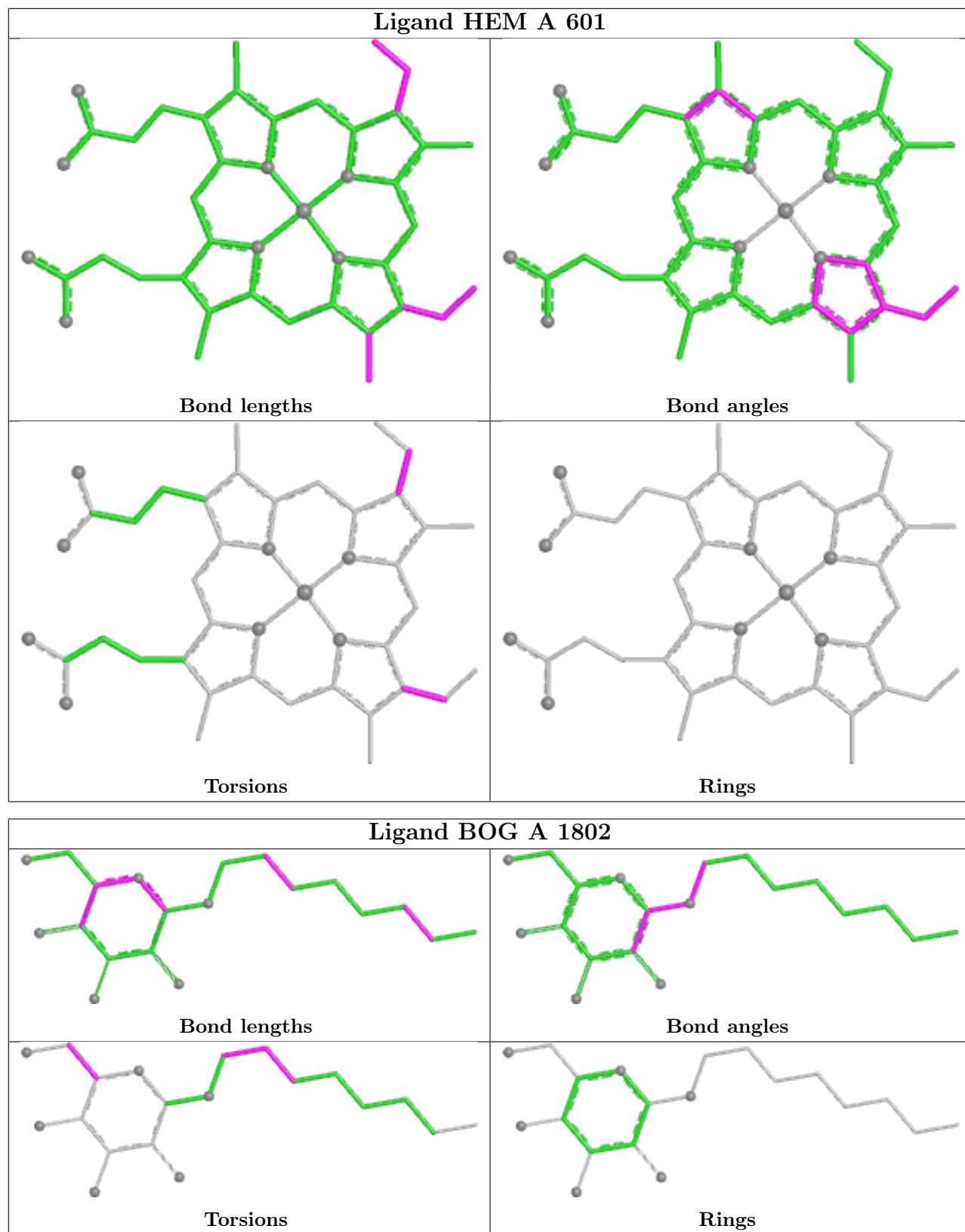
There are no ring outliers.

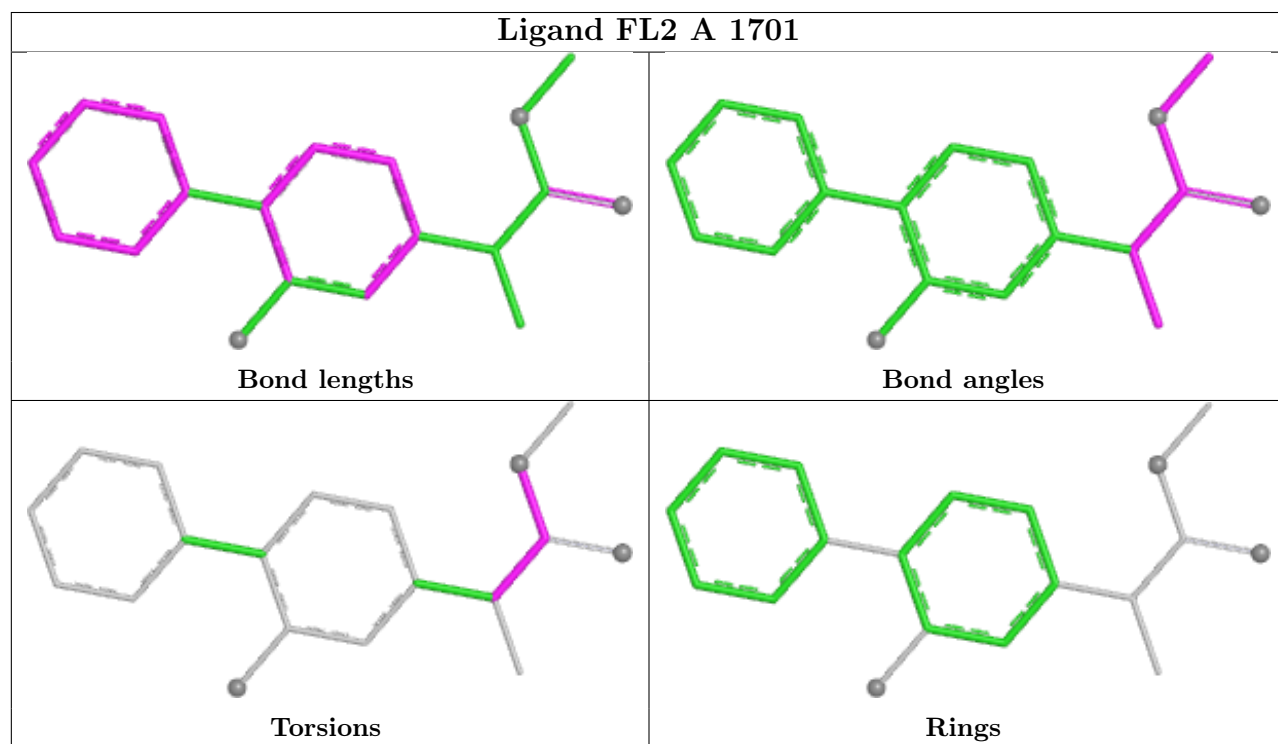
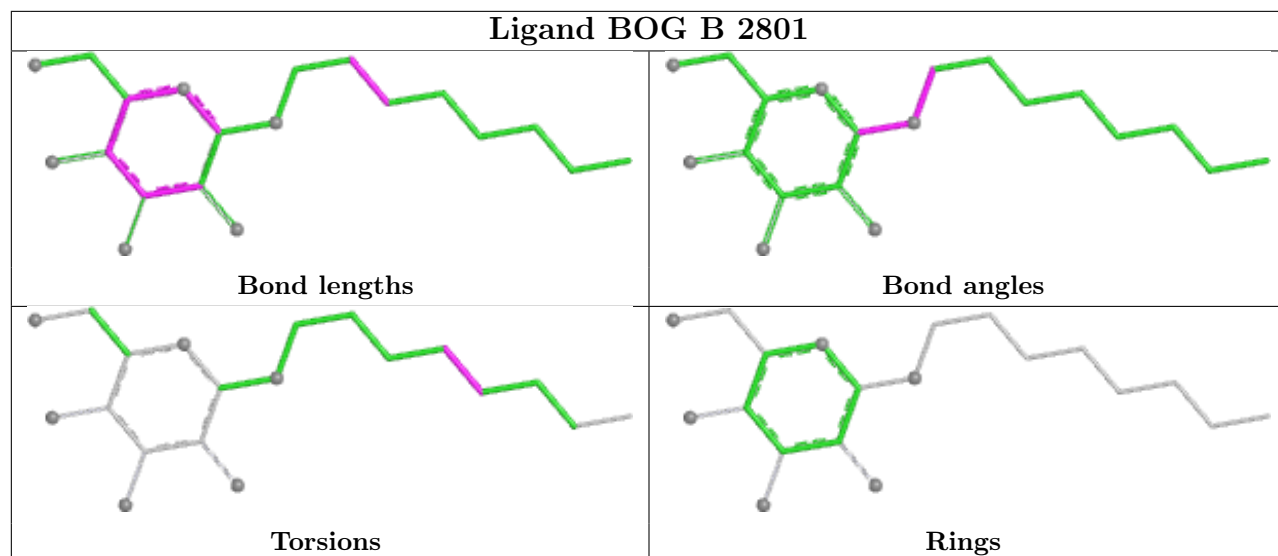
13 monomers are involved in 56 short contacts:

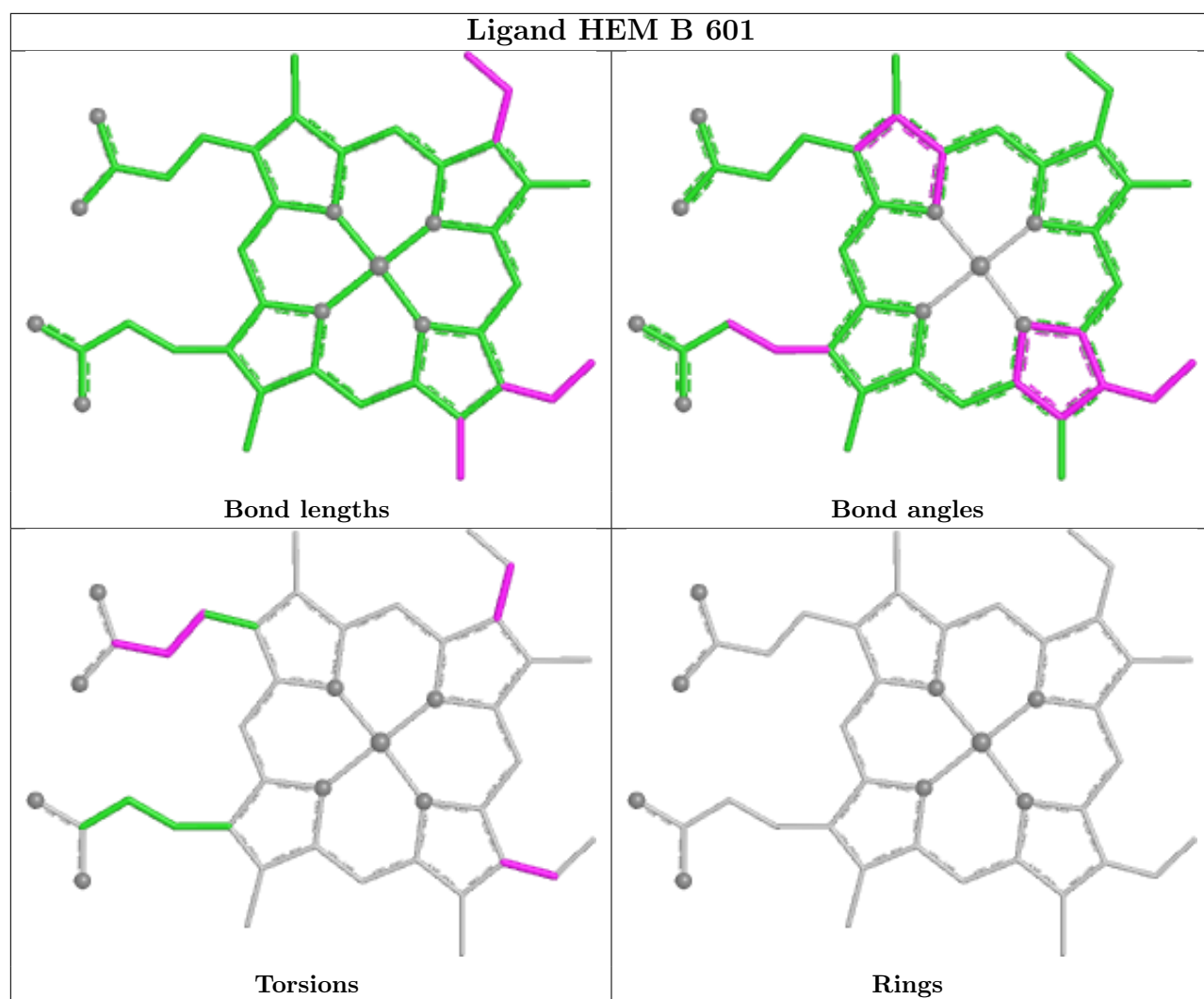
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	2701	FL2	7	0
3	A	1681	NAG	3	0
3	B	1672	NAG	2	0
2	B	2802	BOG	4	0
3	A	1661	NAG	1	0
4	A	601	HEM	5	0
2	A	1802	BOG	8	0
3	A	1671	NAG	7	0
2	B	2801	BOG	2	0
3	B	2671	NAG	5	0
5	A	1701	FL2	4	0
3	B	2681	NAG	2	0
4	B	601	HEM	8	0

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	551/551 (100%)	0.18	14 (2%) 58 56	22, 37, 53, 64	0
1	B	551/551 (100%)	0.15	14 (2%) 58 56	23, 37, 53, 64	0
All	All	1102/1102 (100%)	0.17	28 (2%) 58 56	22, 37, 53, 64	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	294	LEU	3.6
1	A	33	VAL	3.4
1	A	107	PHE	3.4
1	A	581	HIS	3.3
1	B	288	GLY	3.2
1	B	409	PHE	3.2
1	B	74	ILE	3.0
1	A	358	GLN	2.9
1	B	239	GLU	2.8
1	A	278	GLY	2.7
1	A	409	PHE	2.6
1	B	54	ARG	2.4
1	B	75	TRP	2.4
1	A	485	LYS	2.3
1	B	485	LYS	2.3
1	B	282	GLN	2.3
1	A	294	LEU	2.3
1	A	98	TRP	2.3
1	A	99	LEU	2.2
1	B	408	LEU	2.2
1	B	215	LYS	2.1
1	A	554	VAL	2.1
1	A	102	PHE	2.1
1	B	358	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	33	VAL	2.1
1	B	225	GLY	2.1
1	A	51	GLY	2.0
1	A	215	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.

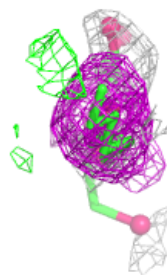
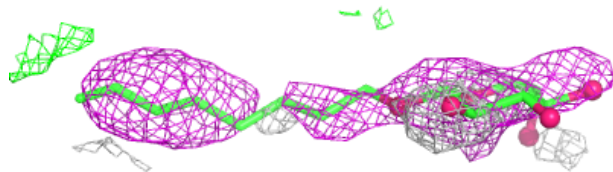
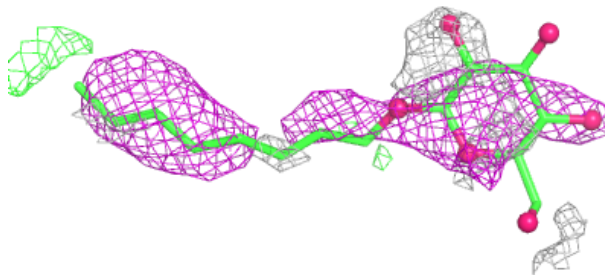
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BOG	B	2801	20/20	0.40	0.31	37,63,68,68	0
3	NAG	B	2662	14/15	0.55	0.28	63,84,96,96	0
3	NAG	A	1662	14/15	0.57	0.27	63,84,96,96	0
3	NAG	A	1681	14/15	0.76	0.19	31,36,42,42	0
3	NAG	A	2672	14/15	0.78	0.20	32,39,65,76	0
3	NAG	B	2681	14/15	0.78	0.18	31,36,42,42	0
2	BOG	A	1802	20/20	0.79	0.19	37,54,64,67	0
2	BOG	B	2802	20/20	0.79	0.19	37,54,64,67	0
3	NAG	B	1672	14/15	0.83	0.18	32,39,65,76	0
3	NAG	A	1661	14/15	0.84	0.17	50,60,64,78	0
3	NAG	B	2661	14/15	0.84	0.17	50,60,64,78	0
3	NAG	A	1671	14/15	0.91	0.10	6,16,23,25	0
5	FL2	A	1701	19/19	0.91	0.16	50,51,56,59	0
5	FL2	B	2701	19/19	0.92	0.14	50,52,56,59	0
3	NAG	B	2671	14/15	0.94	0.08	6,16,23,25	0
4	HEM	A	601	43/43	0.95	0.12	36,41,53,61	0
4	HEM	B	601	43/43	0.96	0.11	36,40,55,65	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.

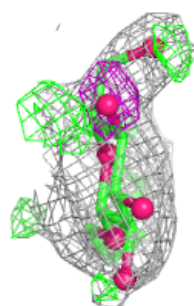
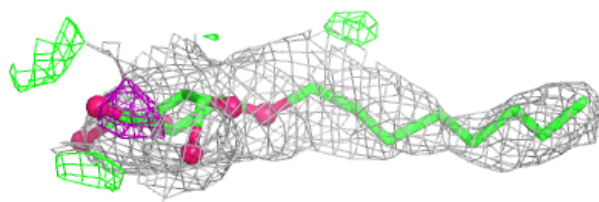
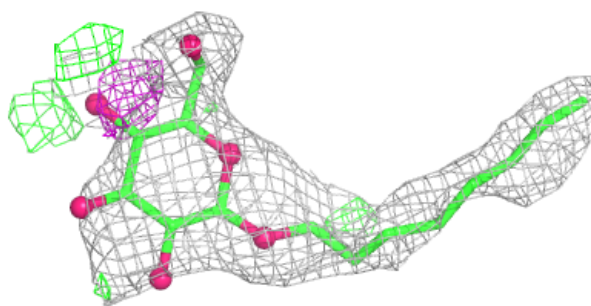
Electron density around BOG B 2801:

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

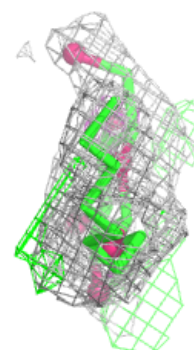
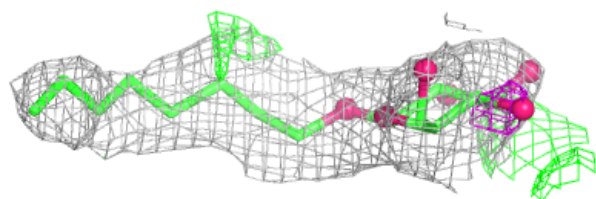
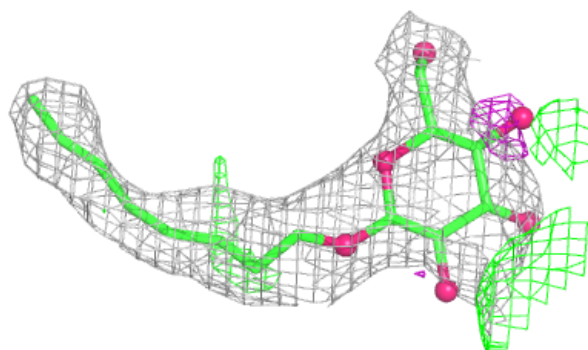


Electron density around BOG A 1802:

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

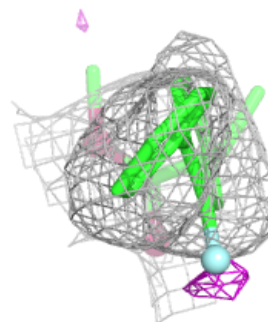
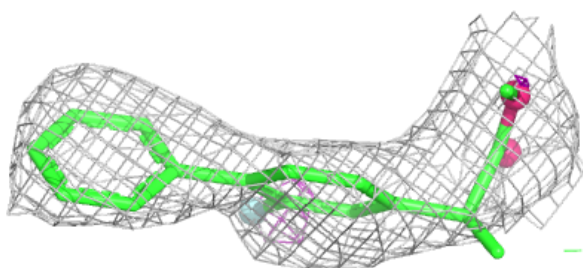
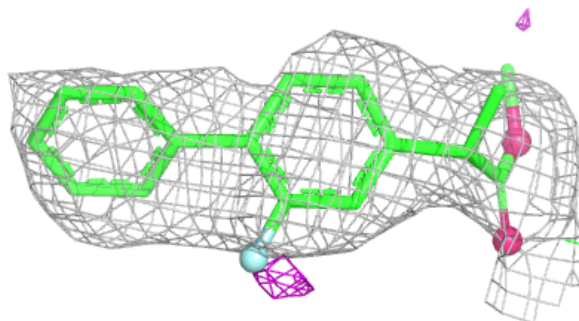
**Electron density around BOG B 2802:**

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

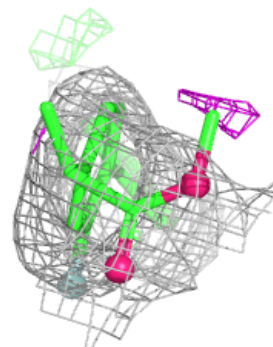
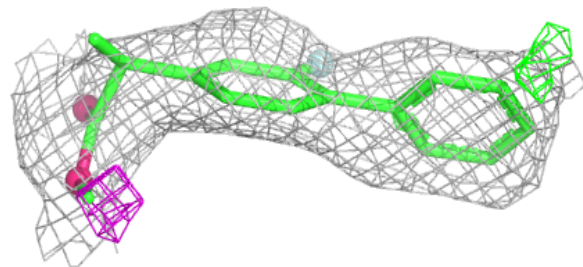
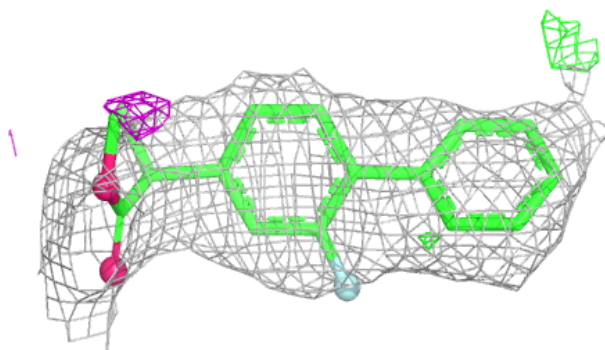


Electron density around FL2 A 1701:

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

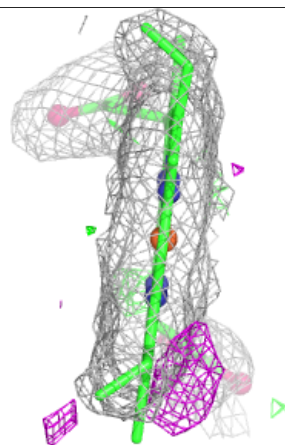
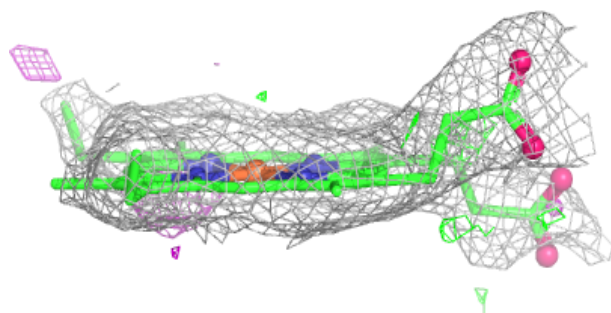
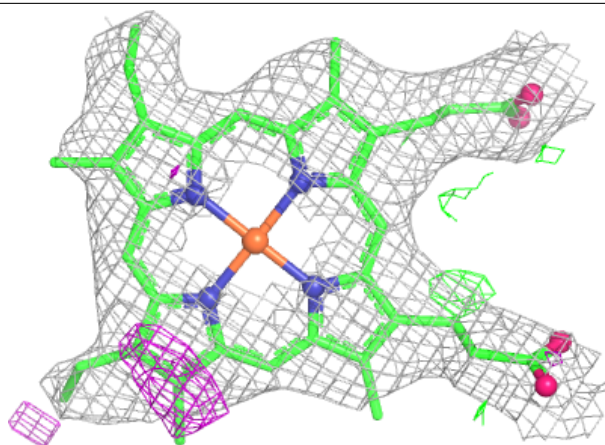
**Electron density around FL2 B 2701:**

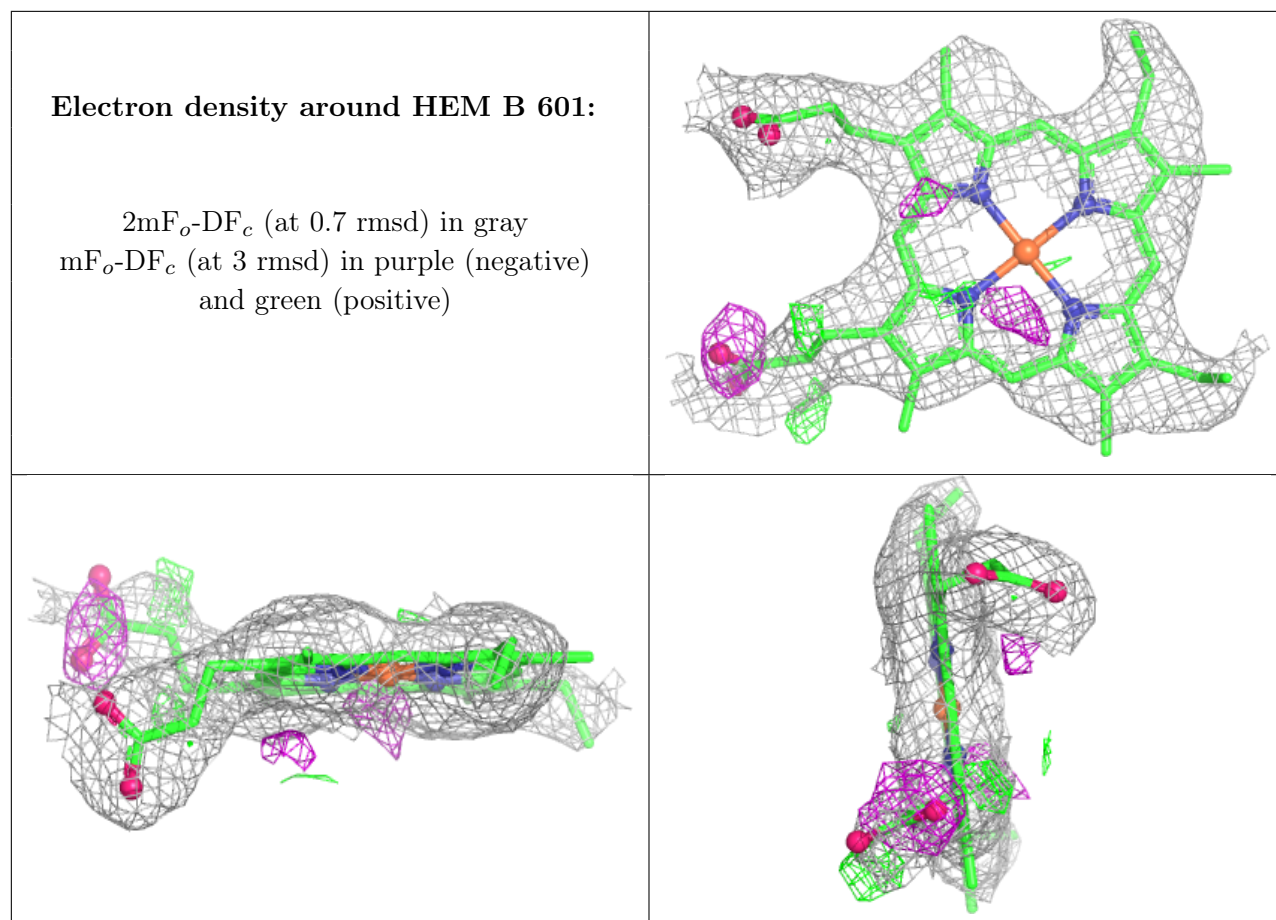
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 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.