



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:33 AM UTC

PDB ID : 2Q9F / pdb_00002q9f
Title : Crystal structure of human cytochrome P450 46A1 in complex with cholesterol-3-sulphate
Authors : White, M.A.; Mast, N.V.; Johnson, E.F.; Stout, C.D.; Pikuleva, I.A.
Deposited on : 2007-06-12
Resolution : 1.90 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

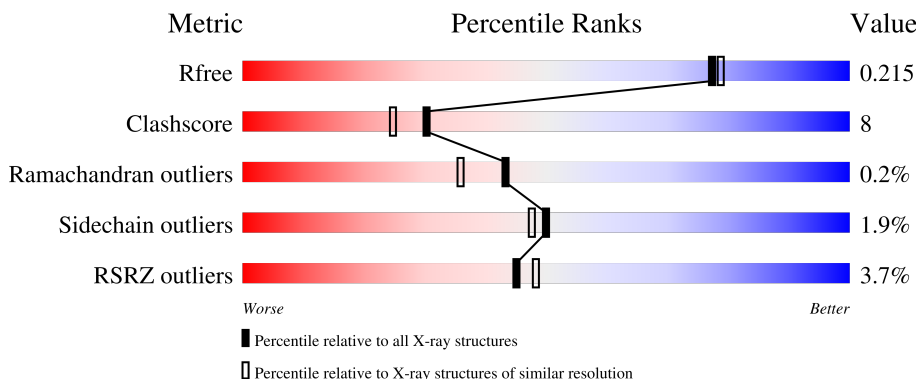
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	456	

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	C3S	A	600	X	-	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3734 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.

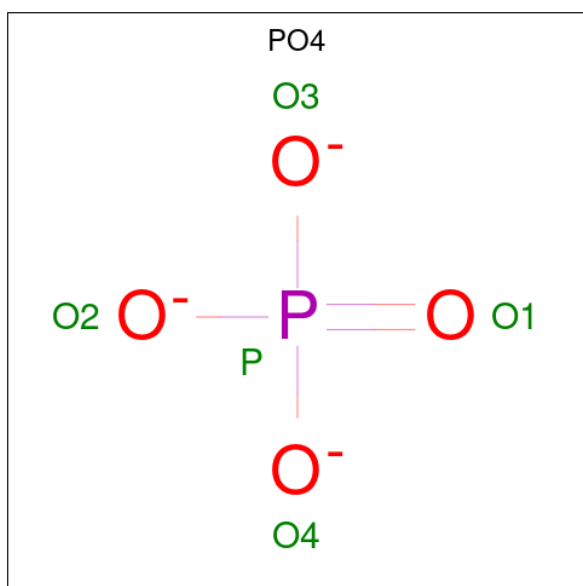
- Molecule 1 is a protein called Cytochrome P450 46A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	433	3488	2231	609	631	17	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

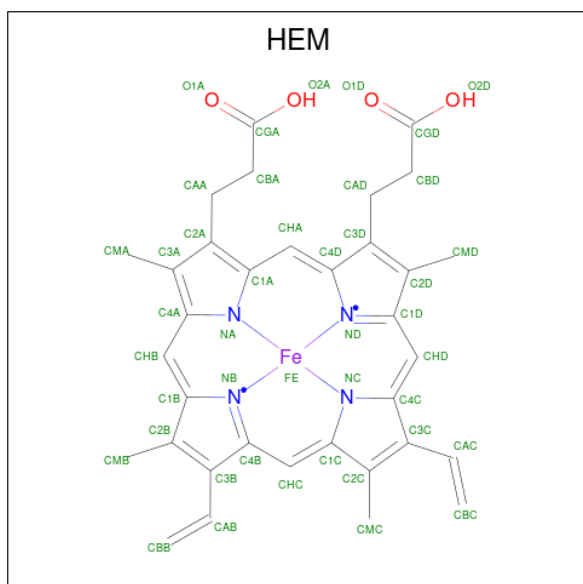
Chain	Residue	Modelled	Actual	Comment	Reference
A	49	MET	-	expression tag	UNP Q9Y6A2
A	50	ALA	-	expression tag	UNP Q9Y6A2
A	501	HIS	-	expression tag	UNP Q9Y6A2
A	502	HIS	-	expression tag	UNP Q9Y6A2
A	503	HIS	-	expression tag	UNP Q9Y6A2
A	504	HIS	-	expression tag	UNP Q9Y6A2

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



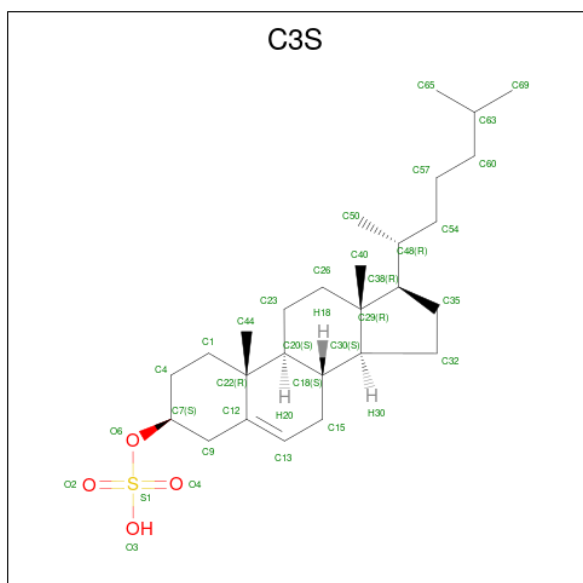
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 4 is CHOLEST-5-EN-3-YL HYDROGEN SULFATE (CCD ID: C3S) (formula: $C_{27}H_{46}O_4S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			32	27	4	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 6 is water.

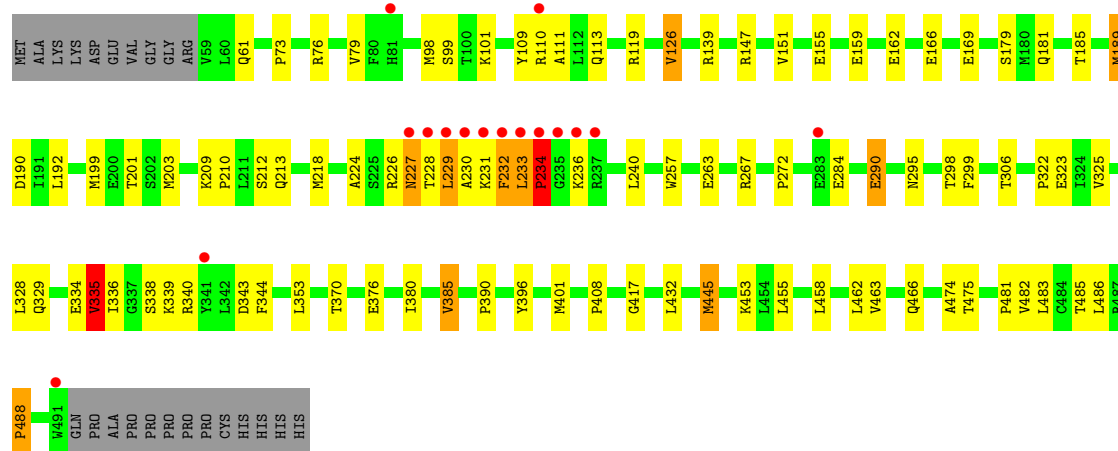
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	130	Total	O	0	0
			130	130		

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 46A1

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.85Å 121.85Å 144.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.08 – 1.90 43.08 – 1.90	Depositor EDS
% Data completeness (in resolution range)	95.4 (43.08-1.90) 96.1 (43.08-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 1.89Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.183 , 0.216 0.184 , 0.215	Depositor DCC
R_{free} test set	2100 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.8	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	0.97	EDS
Total number of atoms	3734	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% 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: GOL, C3S, PO4, 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	1.07	7/3560 (0.2%)	1.24	23/4804 (0.5%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	218	MET	SD-CE	8.10	1.99	1.79
1	A	445	MET	SD-CE	6.59	1.96	1.79
1	A	189	MET	SD-CE	6.18	1.95	1.79
1	A	401	MET	SD-CE	5.68	1.93	1.79
1	A	335	VAL	CA-CB	5.23	1.61	1.54
1	A	98	MET	SD-CE	5.20	1.92	1.79
1	A	482	VAL	CA-CB	5.09	1.61	1.54

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	233	LEU	CA-C-N	8.74	128.06	118.97
1	A	233	LEU	C-N-CA	8.74	128.06	118.97
1	A	232	PHE	N-CA-C	8.31	120.26	108.54
1	A	229	LEU	N-CA-C	-7.24	104.43	113.55
1	A	295	ASN	N-CA-C	7.00	118.91	111.28
1	A	417	GLY	CA-C-N	-6.89	112.89	119.85
1	A	417	GLY	C-N-CA	-6.89	112.89	119.85
1	A	234	PRO	N-CA-C	6.74	123.11	113.47
1	A	390	PRO	CB-CA-C	-6.57	105.42	111.40
1	A	299	PHE	N-CA-C	5.75	117.54	111.28
1	A	126	VAL	CB-CA-C	-5.68	104.62	112.24
1	A	408	PRO	N-CA-C	5.63	121.97	113.81
1	A	213	GLN	N-CA-C	5.48	117.69	111.11
1	A	488	PRO	N-CA-C	-5.47	102.51	111.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	385	VAL	CA-C-N	5.41	125.35	119.78
1	A	385	VAL	C-N-CA	5.41	125.35	119.78
1	A	284	GLU	N-CA-C	5.35	120.15	111.37
1	A	139	ARG	N-CA-C	5.30	117.80	111.71
1	A	353	LEU	N-CA-C	-5.26	105.54	111.28
1	A	290	GLU	N-CA-C	5.14	116.89	111.28
1	A	463	VAL	CA-C-N	-5.10	114.53	119.78
1	A	463	VAL	C-N-CA	-5.10	114.53	119.78
1	A	334	GLU	N-CA-C	5.07	116.50	111.07

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	3488	0	3539	59	0
2	A	5	0	0	0	0
3	A	43	0	30	2	0
4	A	32	0	44	3	0
5	A	36	0	48	3	0
6	A	130	0	0	1	0
All	All	3734	0	3661	61	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 8.

All (61) 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:A:189:MET:SD	1:A:203:MET:SD	2.81	0.79
1:A:224:ALA:O	1:A:230:ALA:HB2	1.95	0.66
1:A:227:ASN:HD22	1:A:227:ASN:H	1.41	0.66
1:A:155:GLU:O	1:A:159:GLU:HG3	1.97	0.65
1:A:110:ARG:HG3	1:A:111:ALA:N	2.10	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:VAL:O	1:A:329:GLN:HG3	2.00	0.61
1:A:263:GLU:HB3	1:A:267:ARG:HH22	1.66	0.61
1:A:485:THR:C	1:A:486:LEU:HD12	2.26	0.61
1:A:181:GLN:O	1:A:185:THR:HG23	2.00	0.60
4:A:600:C3S:H262	4:A:600:C3S:H501	1.83	0.60
1:A:232:PHE:O	1:A:233:LEU:HD23	2.02	0.60
1:A:306:THR:HG22	1:A:475:THR:HG21	1.82	0.60
1:A:233:LEU:HB3	1:A:234:PRO:HD2	1.83	0.60
1:A:126:VAL:HG11	4:A:600:C3S:H691	1.85	0.59
3:A:602:HEM:HBA2	6:A:715:HOH:O	2.04	0.56
1:A:335:VAL:HG23	1:A:336:ILE:HG23	1.88	0.55
1:A:76:ARG:CZ	1:A:76:ARG:HB2	2.36	0.54
1:A:486:LEU:HD12	1:A:486:LEU:N	2.23	0.54
1:A:370:THR:OG1	3:A:602:HEM:O2A	2.27	0.51
1:A:199:MET:HG3	1:A:257:TRP:CD2	2.45	0.51
1:A:380:ILE:HD12	1:A:385:VAL:HG21	1.91	0.51
1:A:328:LEU:HB3	1:A:455:LEU:HD13	1.93	0.51
1:A:335:VAL:CG2	1:A:336:ILE:HG23	2.41	0.51
1:A:340:ARG:O	1:A:453:LYS:NZ	2.39	0.51
1:A:481:PRO:HG2	1:A:483:LEU:HG	1.93	0.50
1:A:192:LEU:C	1:A:192:LEU:HD13	2.36	0.50
1:A:236:LYS:O	1:A:240:LEU:HD13	2.12	0.50
1:A:339:LYS:NZ	1:A:343:ASP:OD2	2.44	0.50
1:A:169:GLU:HG2	1:A:458:LEU:HD21	1.94	0.49
1:A:110:ARG:NH2	4:A:600:C3S:O4	2.46	0.48
1:A:99:SER:OG	1:A:101:LYS:HG2	2.14	0.48
1:A:226:ARG:NH1	1:A:229:LEU:HD13	2.29	0.48
1:A:147:ARG:O	1:A:151:VAL:HG23	2.14	0.47
1:A:231:LYS:HG3	1:A:232:PHE:CD2	2.48	0.47
1:A:203:MET:HE1	1:A:212:SER:OG	2.16	0.46
1:A:109:TYR:O	1:A:113:GLN:HG3	2.17	0.45
1:A:179:SER:HA	1:A:483:LEU:HD23	1.99	0.45
1:A:432:LEU:HD22	5:A:934:GOL:H31	1.99	0.45
1:A:263:GLU:HB3	1:A:267:ARG:NH2	2.32	0.44
1:A:344:PHE:CD2	1:A:445:MET:HE3	2.52	0.44
1:A:79:VAL:HA	1:A:228:THR:HG21	2.00	0.44
1:A:322:PRO:HG2	1:A:323:GLU:OE1	2.17	0.44
1:A:231:LYS:HA	1:A:231:LYS:HD2	1.74	0.44
1:A:190:ASP:OD1	1:A:201:THR:OG1	2.31	0.43
1:A:227:ASN:H	1:A:227:ASN:ND2	2.10	0.43
1:A:101:LYS:HE3	1:A:101:LYS:HA	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:LYS:HZ1	5:A:935:GOL:C1	2.31	0.43
1:A:119:ARG:NH1	1:A:290:GLU:HB3	2.34	0.43
1:A:209:LYS:NZ	5:A:935:GOL:C1	2.82	0.43
1:A:126:VAL:HG23	1:A:298:THR:HG23	2.00	0.43
1:A:226:ARG:CZ	1:A:229:LEU:HB2	2.48	0.43
1:A:462:LEU:HG	1:A:466:GLN:HB2	2.00	0.43
1:A:226:ARG:NH1	1:A:229:LEU:HD22	2.34	0.43
1:A:344:PHE:HD2	1:A:445:MET:HE3	1.84	0.42
1:A:376:GLU:O	1:A:376:GLU:HG3	2.19	0.42
1:A:162:GLU:OE2	1:A:166:GLU:OE2	2.37	0.42
1:A:226:ARG:HH12	1:A:229:LEU:HD13	1.85	0.42
1:A:483:LEU:HD23	1:A:483:LEU:HA	1.91	0.41
1:A:61:GLN:HB3	1:A:396:TYR:CD1	2.55	0.41
1:A:227:ASN:HD22	1:A:227:ASN:N	2.15	0.41
1:A:338:SER:O	1:A:339:LYS:C	2.62	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	431/456 (94%)	420 (97%)	10 (2%)	1 (0%)	43	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	474	ALA

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	378/397 (95%)	371 (98%)	7 (2%)	50 47

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	PRO
1	A	210	PRO
1	A	227	ASN
1	A	234	PRO
1	A	272	PRO
1	A	335	VAL
1	A	488	PRO

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	213	GLN
1	A	227	ASN
1	A	355	GLN
1	A	444	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

9 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	GOL	A	932	-	5,5,5	0.66	0	5,5,5	0.40	0
4	C3S	A	600	-	35,35,35	2.34	6 (17%)	51,55,55	2.98	16 (31%)
5	GOL	A	936	-	5,5,5	0.45	0	5,5,5	0.50	0
5	GOL	A	931	-	5,5,5	0.21	0	5,5,5	0.42	0
5	GOL	A	933	-	5,5,5	0.28	0	5,5,5	0.62	0
3	HEM	A	602	1	50,50,50	2.57	27 (54%)	67,82,82	1.68	13 (19%)
5	GOL	A	934	-	5,5,5	0.64	0	5,5,5	0.29	0
2	PO4	A	601	-	4,4,4	1.59	1 (25%)	6,6,6	0.44	0
5	GOL	A	935	-	5,5,5	0.24	0	5,5,5	0.45	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
5	GOL	A	932	-	-	0/4/4/4	-
4	C3S	A	600	-	2/2/11/12	2/15/73/73	0/4/4/4
5	GOL	A	936	-	-	0/4/4/4	-
5	GOL	A	931	-	-	0/4/4/4	-
5	GOL	A	933	-	-	0/4/4/4	-
3	HEM	A	602	1	-	4/14/54/54	-
5	GOL	A	934	-	-	0/4/4/4	-
5	GOL	A	935	-	-	0/4/4/4	-

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	600	C3S	C15-C18	-8.35	1.39	1.53
4	A	600	C3S	C18-C20	-7.27	1.40	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	HEM	O2A-CGA	-6.28	1.10	1.30
3	A	602	HEM	CBA-CAA	-5.22	1.33	1.51
4	A	600	C3S	C15-C13	-5.07	1.39	1.50
4	A	600	C3S	O6-S1	-4.36	1.44	1.57
3	A	602	HEM	FE-NC	4.34	2.09	1.95
3	A	602	HEM	C1C-NC	-4.28	1.31	1.39
3	A	602	HEM	FE-ND	3.90	2.06	1.94
3	A	602	HEM	FE-NA	3.89	2.08	1.95
3	A	602	HEM	CMA-C3A	3.87	1.58	1.50
3	A	602	HEM	C3C-C2C	3.38	1.44	1.37
3	A	602	HEM	O1A-CGA	3.33	1.33	1.22
3	A	602	HEM	CMD-C2D	3.23	1.57	1.50
3	A	602	HEM	CMB-C2B	3.15	1.57	1.50
4	A	600	C3S	C13-C12	3.13	1.39	1.33
3	A	602	HEM	C4B-NB	-3.10	1.32	1.38
3	A	602	HEM	C4A-NA	-3.09	1.33	1.39
3	A	602	HEM	CBA-CGA	2.95	1.57	1.50
3	A	602	HEM	CBB-CAB	2.93	1.44	1.30
3	A	602	HEM	CHA-C1A	2.78	1.45	1.39
3	A	602	HEM	C3B-C2B	2.70	1.42	1.37
3	A	602	HEM	FE-NB	2.61	2.02	1.94
3	A	602	HEM	C4D-ND	-2.56	1.35	1.40
3	A	602	HEM	C1A-C2A	-2.55	1.39	1.44
3	A	602	HEM	CBC-CAC	2.49	1.42	1.30
3	A	602	HEM	CAB-C3B	-2.47	1.40	1.47
3	A	602	HEM	CAA-C2A	2.26	1.57	1.51
3	A	602	HEM	CHA-C4D	2.17	1.42	1.38
3	A	602	HEM	CBD-CAD	2.11	1.59	1.51
3	A	602	HEM	CMC-C2C	2.05	1.55	1.50
3	A	602	HEM	C1C-C2C	-2.04	1.41	1.45
2	A	601	PO4	P-O4	-2.04	1.48	1.54
4	A	600	C3S	O6-C7	-2.02	1.44	1.48

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	600	C3S	C15-C18-C20	9.74	120.98	109.72
4	A	600	C3S	C23-C20-C18	8.22	123.24	111.78
3	A	602	HEM	CBA-CAA-C2A	7.11	132.20	112.53
4	A	600	C3S	C18-C15-C13	7.04	122.52	112.76
4	A	600	C3S	C20-C18-C30	7.02	118.26	109.09
4	A	600	C3S	C44-C22-C20	-6.97	103.84	111.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	600	C3S	C15-C18-C30	6.43	120.04	110.93
4	A	600	C3S	C22-C20-C18	4.20	118.85	112.71
3	A	602	HEM	C4C-C3C-C2C	-3.55	103.73	106.81
3	A	602	HEM	CAA-CBA-CGA	-3.26	105.03	113.67
4	A	600	C3S	C23-C20-C22	3.22	117.06	113.08
3	A	602	HEM	CHD-C4C-NC	-2.91	121.28	124.45
3	A	602	HEM	C3B-C2B-C1B	-2.91	104.23	106.41
3	A	602	HEM	CBD-CAD-C3D	-2.81	104.75	112.53
4	A	600	C3S	C35-C32-C30	-2.72	99.82	105.14
4	A	600	C3S	C30-C29-C38	-2.71	96.99	100.10
4	A	600	C3S	C20-C22-C12	2.65	113.54	109.65
3	A	602	HEM	O1A-CGA-CBA	2.56	131.20	123.09
4	A	600	C3S	C9-C12-C22	2.49	119.61	116.42
3	A	602	HEM	C4D-ND-C1D	2.44	108.10	105.21
4	A	600	C3S	C15-C13-C12	-2.42	120.93	125.02
4	A	600	C3S	C1-C4-C7	2.36	114.18	110.33
3	A	602	HEM	C1C-CHC-C4B	2.30	130.90	126.02
4	A	600	C3S	C29-C38-C48	-2.24	116.04	119.50
3	A	602	HEM	C1B-NB-C4B	2.22	107.83	105.21
4	A	600	C3S	C1-C22-C20	2.19	111.64	108.74
3	A	602	HEM	CHA-C1A-NA	-2.15	119.97	123.86
3	A	602	HEM	O2A-CGA-O1A	-2.06	118.04	123.33
3	A	602	HEM	CHB-C4A-NA	-2.03	120.17	123.86

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	600	C3S	C18
4	A	600	C3S	C20

All (6) torsion outliers are listed below:

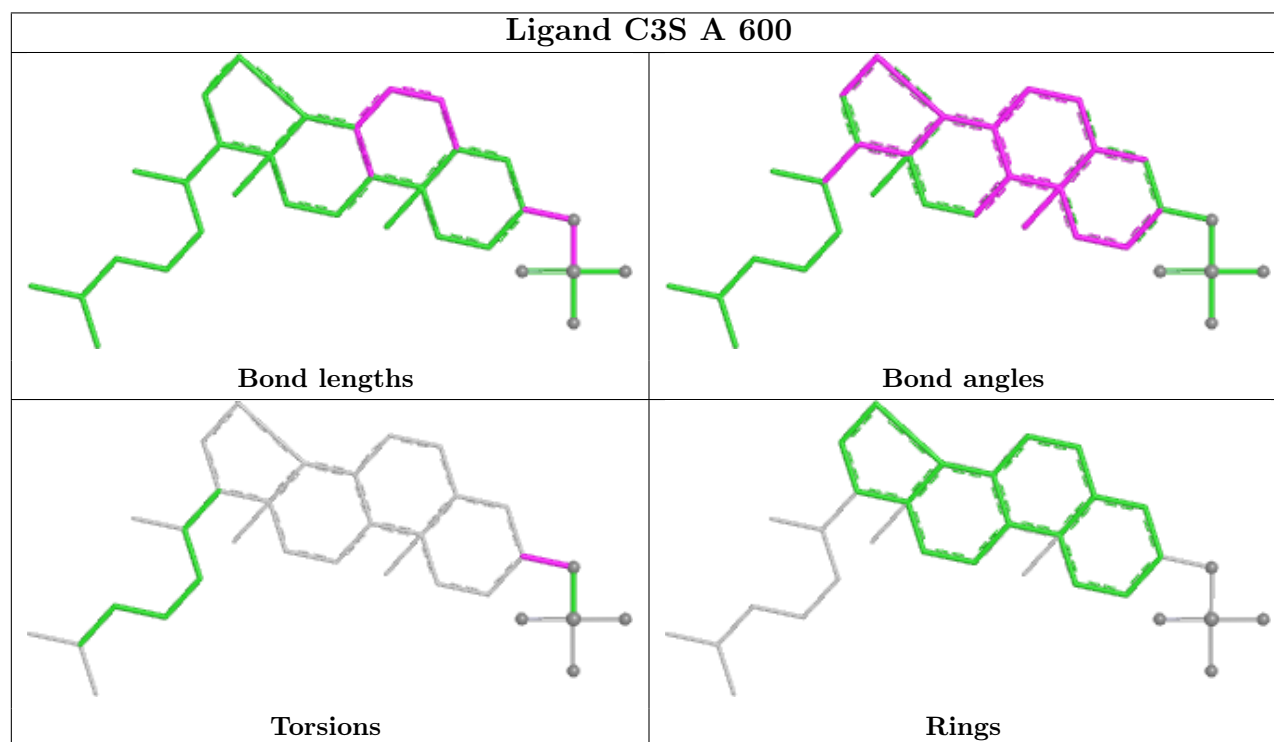
Mol	Chain	Res	Type	Atoms
4	A	600	C3S	C9-C7-O6-S1
4	A	600	C3S	C4-C7-O6-S1
3	A	602	HEM	CAA-CBA-CGA-O2A
3	A	602	HEM	CAA-CBA-CGA-O1A
3	A	602	HEM	CAD-CBD-CGD-O1D
3	A	602	HEM	CAD-CBD-CGD-O2D

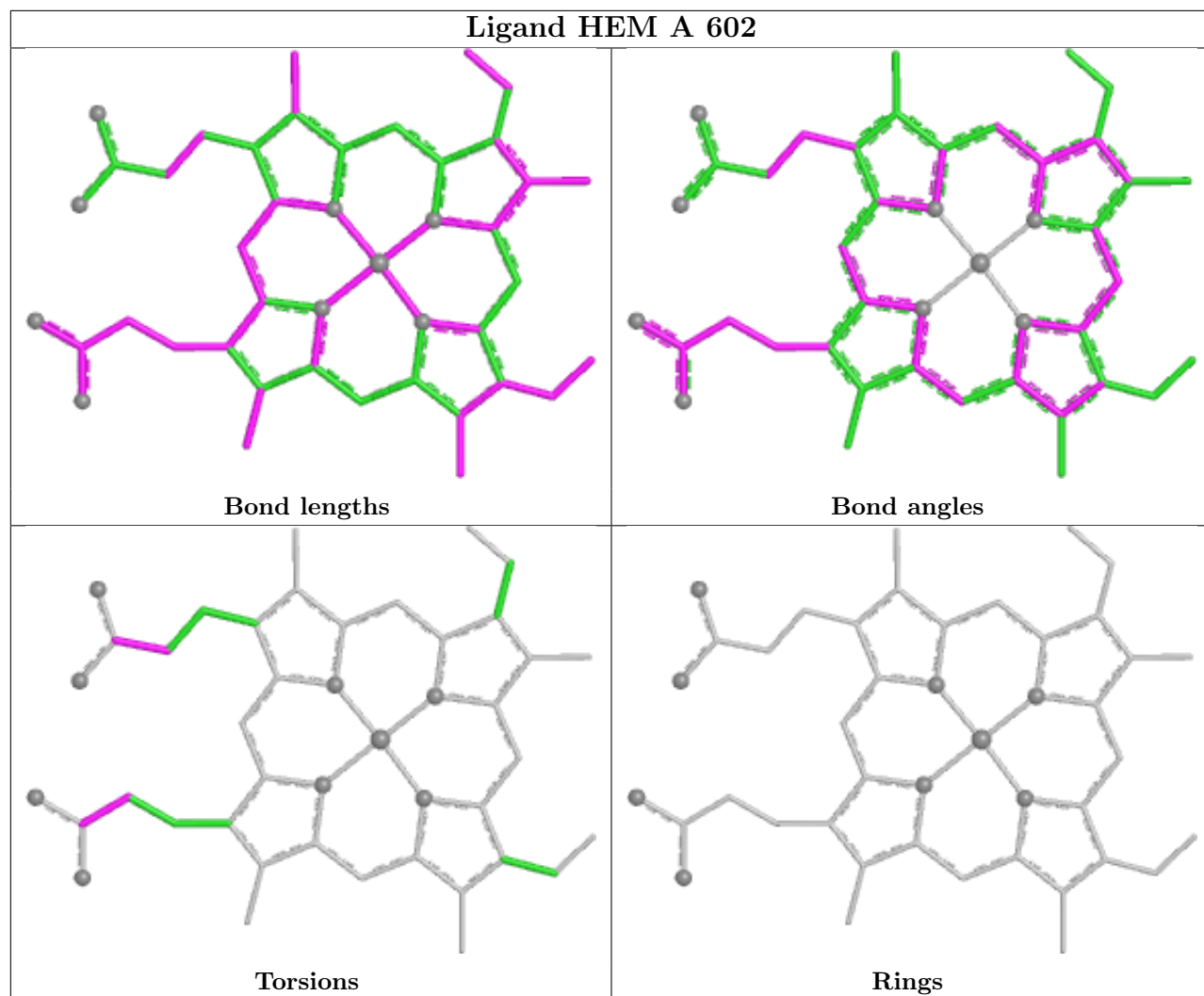
There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	600	C3S	3	0
3	A	602	HEM	2	0
5	A	934	GOL	1	0
5	A	935	GOL	2	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	433/456 (94%)	0.17	16 (3%) 45 48	27, 42, 74, 111	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	491	TRP	5.6
1	A	234	PRO	4.7
1	A	232	PHE	4.0
1	A	341	TYR	3.8
1	A	230	ALA	3.8
1	A	233	LEU	3.4
1	A	236	LYS	3.3
1	A	235	GLY	3.2
1	A	81	HIS	3.1
1	A	283	GLU	3.1
1	A	231	LYS	2.7
1	A	229	LEU	2.5
1	A	227	ASN	2.4
1	A	237	ARG	2.2
1	A	110	ARG	2.1
1	A	228	THR	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](#)

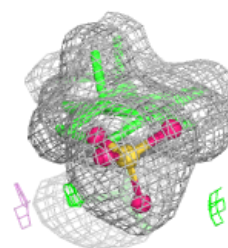
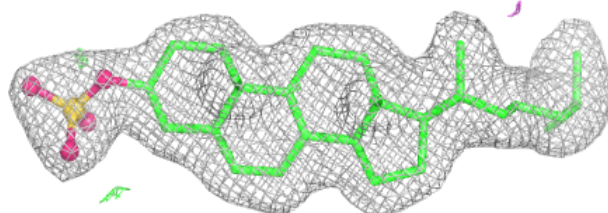
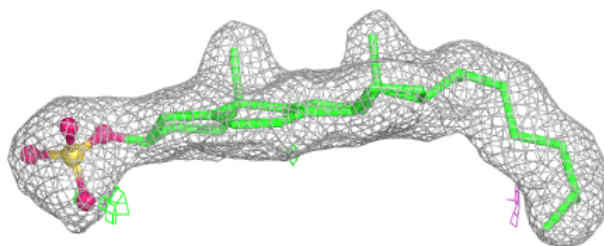
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GOL	A	933	6/6	0.79	0.23	57,70,82,85	0
5	GOL	A	935	6/6	0.79	0.13	65,73,83,88	0
5	GOL	A	936	6/6	0.83	0.16	61,88,97,102	0
5	GOL	A	934	6/6	0.84	0.12	46,71,77,79	0
5	GOL	A	931	6/6	0.86	0.14	66,78,84,93	0
2	PO4	A	601	5/5	0.88	0.09	87,91,114,119	0
5	GOL	A	932	6/6	0.89	0.10	47,79,80,87	0
4	C3S	A	600	32/32	0.96	0.07	31,36,52,58	0
3	HEM	A	602	43/43	0.98	0.05	26,28,36,46	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.

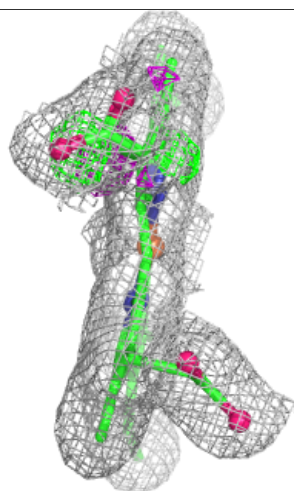
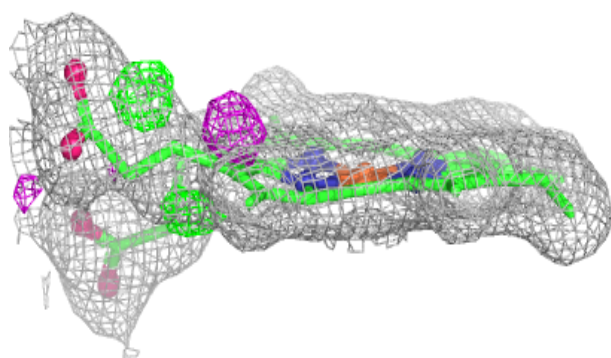
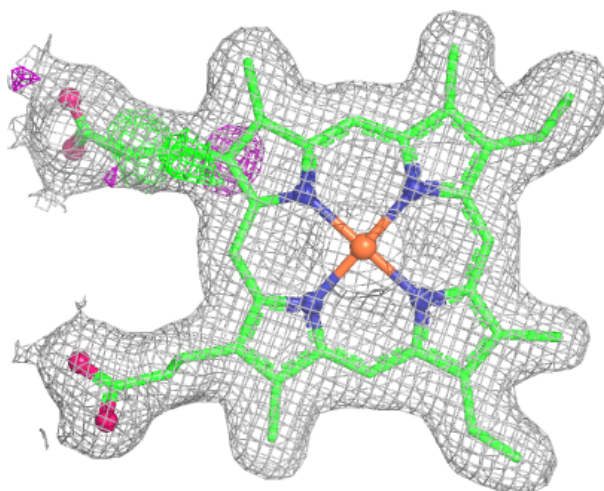
Electron density around C3S A 600:

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 602:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.