



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 10:29 am BST

PDB ID : 5VCE  
Title : Crystal structure of the cysteine depleted CYP3A4 bound to ritonavir  
Authors : Sevrioukova, I.  
Deposited on : 2017-03-31  
Resolution : 2.20 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

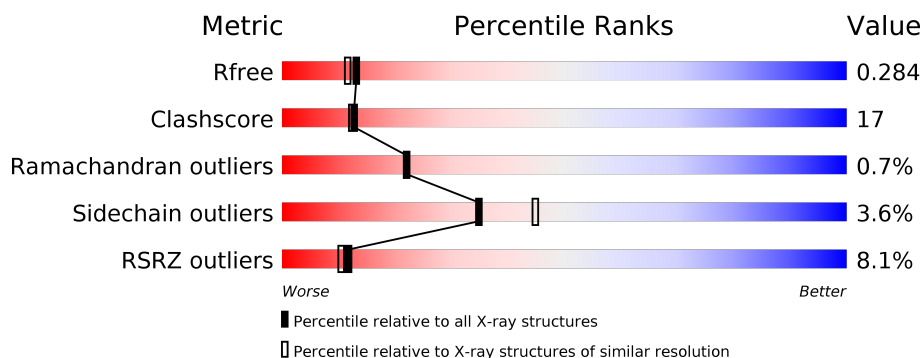
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	<div> <div>8%</div> <div>69%</div> <div>23%</div> <div>6%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3793 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	457	Total	C	N	O	S	0	2	0
			3683	2402	606	656	19			

There are 12 discrepancies between the modelled and reference sequences:

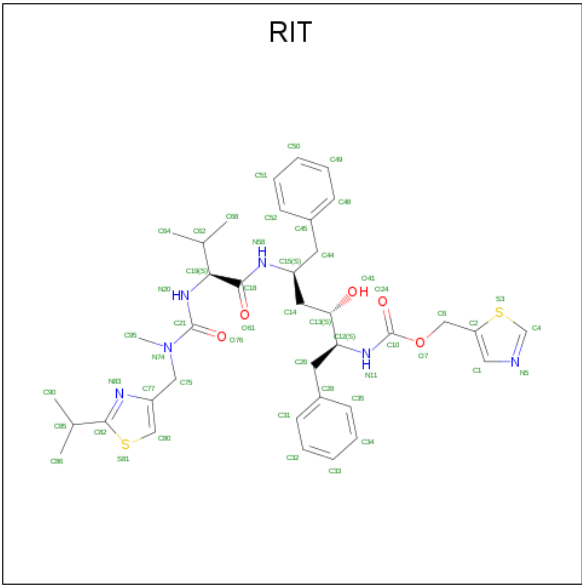
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	see remark 999	UNP P08684
A	2	ALA	-	see remark 999	UNP P08684
A	58	ALA	CYS	engineered mutation	UNP P08684
A	64	MET	CYS	engineered mutation	UNP P08684
A	98	ALA	CYS	engineered mutation	UNP P08684
A	239	THR	CYS	engineered mutation	UNP P08684
A	377	ALA	CYS	engineered mutation	UNP P08684
A	468	SER	CYS	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
A	507	HIS	-	expression tag	UNP P08684

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



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

- Molecule 3 is RITONAVIR (three-letter code: RIT) (formula:  $C_{37}H_{48}N_6O_5S_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			50	37	6	5	2		

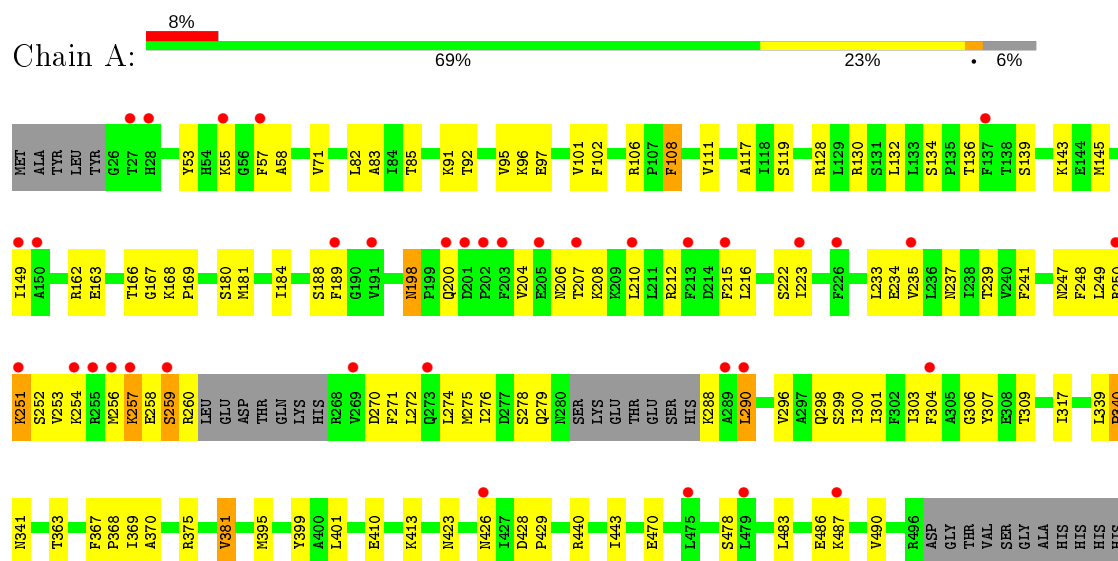
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	17	Total	O	0	0
			17	17		

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.37Å 100.71Å 129.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	79.43 – 2.20 50.35 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (79.43-2.20) 99.2 (50.35-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.29 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.240 , 0.275 0.249 , 0.284	Depositor DCC
$R_{free}$ test set	1318 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	59.3	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 61.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3793	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, RIT

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	0.46	0/3778	0.64	0/5110

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	3683	0	3786	124	0
2	A	43	0	30	3	0
3	A	50	0	47	11	0
4	A	17	0	0	1	0
All	All	3793	0	3863	127	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 17.

All (127) 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:253:VAL:HG12	1:A:257:LYS:HG2	1.32	1.10
1:A:251:LYS:HD3	1:A:254:LYS:HB2	1.09	1.08
1:A:92:THR:HA	1:A:96:LYS:HG2	1.40	1.03
1:A:215:PHE:HB2	3:A:602:RIT:H953	1.39	1.03
1:A:251:LYS:CD	1:A:254:LYS:HB2	1.95	0.96
1:A:248:PHE:O	1:A:252:SER:OG	1.85	0.95
1:A:253:VAL:HG12	1:A:257:LYS:CG	1.97	0.94
1:A:251:LYS:HA	1:A:251:LYS:HE2	1.52	0.92
1:A:136:THR:HG21	1:A:271:PHE:CD1	2.10	0.86
1:A:117:ALA:HB1	1:A:301:ILE:HG13	1.56	0.86
1:A:239:THR:HG22	1:A:241:PHE:H	1.40	0.85
1:A:212:ARG:HG2	1:A:304:PHE:HZ	1.38	0.85
1:A:253:VAL:CG1	1:A:257:LYS:HG2	2.06	0.84
1:A:184:ILE:HG21	1:A:306:GLY:HA3	1.62	0.82
1:A:251:LYS:HD3	1:A:254:LYS:CB	2.04	0.79
1:A:180:SER:OG	1:A:307:TYR:HA	1.86	0.76
1:A:423:ASN:O	1:A:426:ASN:HB2	1.88	0.72
1:A:136:THR:CG2	1:A:271:PHE:CD1	2.72	0.72
3:A:602:RIT:H48	3:A:602:RIT:H643	1.70	0.72
1:A:257:LYS:HA	1:A:258:GLU:HB2	1.72	0.71
2:A:601:HEM:HBC2	2:A:601:HEM:HHD	1.71	0.71
1:A:251:LYS:HD2	1:A:254:LYS:HD2	1.73	0.71
1:A:136:THR:CG2	1:A:271:PHE:CE1	2.74	0.70
1:A:258:GLU:CG	1:A:259:SER:H	2.05	0.70
1:A:55:LYS:O	1:A:55:LYS:HG2	1.92	0.69
1:A:145:MET:O	1:A:149:ILE:HG13	1.92	0.69
1:A:239:THR:HG22	1:A:241:PHE:N	2.06	0.69
1:A:258:GLU:HG2	1:A:259:SER:H	1.58	0.69
1:A:136:THR:HG21	1:A:271:PHE:HD1	1.57	0.69
1:A:259:SER:O	1:A:260:ARG:HG2	1.94	0.67
1:A:136:THR:HG22	1:A:271:PHE:CE1	2.30	0.67
1:A:204:VAL:O	1:A:208:LYS:HG2	1.94	0.67
1:A:206:ASN:ND2	1:A:248:PHE:HB3	2.11	0.65
1:A:117:ALA:HB1	1:A:301:ILE:CG1	2.24	0.65
1:A:128:ARG:HG3	1:A:288:LYS:HE2	1.77	0.65
1:A:85:THR:HB	1:A:401:LEU:HD21	1.78	0.65
1:A:108:PHE:HB2	3:A:602:RIT:N83	2.13	0.64
1:A:270:ASP:O	1:A:274:LEU:HG	1.98	0.63
1:A:57:PHE:CZ	1:A:215:PHE:HD2	2.16	0.63
1:A:367:PHE:HB3	1:A:483:LEU:HD21	1.81	0.62
2:A:601:HEM:C1D	3:A:602:RIT:H1	2.34	0.62
1:A:257:LYS:HB3	1:A:272:LEU:HD21	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:VAL:HG12	1:A:257:LYS:CB	2.31	0.60
1:A:139:SER:O	1:A:143:LYS:HB2	2.02	0.60
1:A:256:MET:C	1:A:258:GLU:HA	2.24	0.58
1:A:92:THR:CA	1:A:96:LYS:HG2	2.27	0.58
1:A:184:ILE:CG2	1:A:306:GLY:HA3	2.32	0.58
1:A:188:SER:HA	1:A:271:PHE:HB2	1.85	0.57
1:A:223:ILE:HD11	1:A:233:LEU:HD12	1.86	0.57
1:A:95:VAL:HG12	1:A:96:LYS:N	2.20	0.57
1:A:189:PHE:HZ	1:A:299:SER:HB3	1.69	0.57
1:A:57:PHE:CZ	1:A:215:PHE:CD2	2.93	0.56
1:A:189:PHE:CZ	1:A:299:SER:HB3	2.40	0.56
1:A:57:PHE:HZ	1:A:215:PHE:HB3	1.70	0.56
1:A:276:ILE:HA	1:A:279:GLN:OE1	2.07	0.55
1:A:101:VAL:HG21	1:A:381:VAL:HG11	1.88	0.55
1:A:470:GLU:HG2	1:A:490:VAL:HG21	1.88	0.55
1:A:96:LYS:HG3	1:A:97:GLU:HG2	1.88	0.55
1:A:71[A]:VAL:HG13	1:A:82:LEU:HD11	1.90	0.54
1:A:136:THR:HG21	1:A:271:PHE:CE1	2.39	0.54
1:A:212:ARG:NH1	3:A:602:RIT:O61	2.41	0.54
1:A:247:ASN:O	1:A:250:ARG:HB2	2.08	0.53
1:A:251:LYS:CE	1:A:251:LYS:HA	2.21	0.53
1:A:251:LYS:CD	1:A:254:LYS:HD2	2.39	0.53
1:A:253:VAL:HG12	1:A:257:LYS:HB3	1.91	0.53
1:A:370:ALA:HA	3:A:602:RIT:H50	1.90	0.53
1:A:253:VAL:O	1:A:257:LYS:HG3	2.08	0.53
1:A:212:ARG:HG2	1:A:304:PHE:CZ	2.30	0.52
1:A:258:GLU:CG	1:A:259:SER:N	2.73	0.51
1:A:317:ILE:HG13	1:A:363:THR:HG21	1.93	0.50
1:A:210:LEU:HD11	1:A:300:ILE:HG23	1.93	0.50
1:A:258:GLU:C	1:A:260:ARG:H	2.14	0.50
1:A:162:ARG:CZ	1:A:163:GLU:OE2	2.59	0.50
1:A:250:ARG:HA	1:A:296:VAL:HG11	1.94	0.49
1:A:257:LYS:HD2	1:A:257:LYS:O	2.12	0.49
1:A:259:SER:O	1:A:260:ARG:CG	2.59	0.49
1:A:256:MET:O	1:A:258:GLU:HG3	2.12	0.49
1:A:248:PHE:C	1:A:252:SER:OG	2.50	0.49
1:A:304:PHE:HD2	3:A:602:RIT:H62A	1.75	0.49
1:A:206:ASN:ND2	1:A:248:PHE:CB	2.76	0.48
1:A:249:LEU:HA	1:A:252:SER:HB2	1.94	0.48
1:A:189:PHE:HB3	1:A:256:MET:SD	2.53	0.48
1:A:111:VAL:O	1:A:234:GLU:OE2	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:PHE:HZ	1:A:215:PHE:CD2	2.32	0.47
1:A:271:PHE:HA	1:A:274:LEU:HD12	1.95	0.47
1:A:253:VAL:HA	1:A:256:MET:HB3	1.97	0.47
1:A:58:ALA:HB1	1:A:399:TYR:CD2	2.49	0.47
1:A:239:THR:CG2	1:A:241:PHE:H	2.21	0.46
1:A:132:LEU:HD11	1:A:288:LYS:O	2.15	0.46
1:A:275:MET:HE1	1:A:298:GLN:HB2	1.98	0.46
1:A:166:THR:O	1:A:168:LYS:N	2.47	0.46
1:A:128:ARG:O	1:A:132:LEU:HG	2.17	0.45
1:A:184:ILE:HD13	1:A:303:ILE:HA	1.97	0.45
1:A:162:ARG:NH2	1:A:163:GLU:OE2	2.50	0.45
1:A:130:ARG:O	1:A:134:SER:HB2	2.16	0.45
1:A:235:VAL:HG23	4:A:702:HOH:O	2.16	0.45
1:A:257:LYS:HB3	1:A:272:LEU:HD11	1.98	0.44
1:A:180:SER:OG	1:A:307:TYR:O	2.21	0.44
1:A:257:LYS:HB2	1:A:276:ILE:HD11	2.00	0.44
1:A:180:SER:OG	1:A:307:TYR:CA	2.60	0.44
1:A:91:LYS:O	1:A:95:VAL:HB	2.18	0.43
1:A:181:MET:SD	1:A:184:ILE:HD11	2.58	0.43
1:A:428:ASP:HA	1:A:429:PRO:HD2	1.83	0.43
1:A:181:MET:HG2	1:A:208:LYS:NZ	2.33	0.43
1:A:132:LEU:HD13	1:A:290:LEU:HG	2.01	0.43
1:A:215:PHE:CB	3:A:602:RIT:H953	2.29	0.43
1:A:368:PRO:O	1:A:483:LEU:HD22	2.19	0.43
1:A:101:VAL:HG21	1:A:381:VAL:CG1	2.48	0.43
1:A:53:TYR:HB3	1:A:216:LEU:HD22	2.01	0.43
1:A:212:ARG:HH12	3:A:602:RIT:H12	1.84	0.43
1:A:181:MET:CE	1:A:207:THR:HB	2.49	0.42
1:A:470:GLU:CG	1:A:490:VAL:HG21	2.48	0.42
1:A:253:VAL:CG1	1:A:257:LYS:HB3	2.48	0.42
1:A:486:GLU:HG3	1:A:487:LYS:N	2.34	0.42
1:A:259:SER:C	1:A:260:ARG:HG2	2.38	0.42
1:A:256:MET:O	1:A:258:GLU:HA	2.19	0.42
1:A:486:GLU:HG3	1:A:487:LYS:H	1.84	0.42
1:A:119:SER:HB2	3:A:602:RIT:H141	2.01	0.42
1:A:198:ASN:C	1:A:200:GLN:H	2.23	0.41
1:A:443:ILE:CG2	2:A:601:HEM:HMD3	2.50	0.41
1:A:83:ALA:HA	1:A:395:MET:O	2.20	0.41
1:A:258:GLU:C	1:A:260:ARG:N	2.73	0.41
1:A:339:LEU:HA	1:A:340:PRO:HD2	1.78	0.41
1:A:410:GLU:OE1	1:A:413:LYS:HD3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:PHE:HB3	1:A:375:ARG:CB	2.51	0.41
1:A:253:VAL:O	1:A:257:LYS:N	2.41	0.40
1:A:369:ILE:HD12	3:A:602:RIT:H50	2.02	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	453/487 (93%)	417 (92%)	33 (7%)	3 (1%)	22	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	340	PRO
1	A	167	GLY
1	A	169	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	416/440 (94%)	401 (96%)	15 (4%)	35	45

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	ARG
1	A	108	PHE
1	A	198	ASN
1	A	222	SER
1	A	237	ASN
1	A	251	LYS
1	A	257	LYS
1	A	259	SER
1	A	278	SER
1	A	290	LEU
1	A	309	THR
1	A	341	ASN
1	A	381	VAL
1	A	440	ARG
1	A	478	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	206	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	RIT	A	602	2	48,53,53	2.56	10 (20%)	55,71,71	1.82	9 (16%)
2	HEM	A	601	1,3	27,50,50	0.79	1 (3%)	17,82,82	1.40	4 (23%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	RIT	A	602	2	-	9/49/53/53	0/4/4/4
2	HEM	A	601	1,3	-	0/6/54/54	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	RIT	O41-C13	-10.51	1.21	1.43
3	A	602	RIT	C18-N58	6.47	1.48	1.34
3	A	602	RIT	C21-N20	5.99	1.48	1.35
3	A	602	RIT	C21-N74	5.52	1.48	1.37
3	A	602	RIT	C10-N11	4.85	1.46	1.34
3	A	602	RIT	O7-C10	4.37	1.43	1.35
3	A	602	RIT	C1-N5	-3.19	1.25	1.35
3	A	602	RIT	C6-C2	2.53	1.53	1.50
2	A	601	HEM	C4D-C3D	2.34	1.47	1.42
3	A	602	RIT	C75-C77	2.16	1.54	1.51
3	A	602	RIT	C82-C85	2.01	1.54	1.50

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	RIT	C1-N5-C4	7.66	117.73	105.78
3	A	602	RIT	C75-C77-C80	-4.89	122.26	129.62
3	A	602	RIT	O7-C10-N11	3.71	118.05	110.50
3	A	602	RIT	C45-C44-C15	-3.68	105.43	113.78
3	A	602	RIT	C19-C18-N58	3.65	124.36	116.48
3	A	602	RIT	O24-C10-N11	-3.07	119.82	124.85
3	A	602	RIT	O61-C18-N58	-2.86	117.63	122.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	HEM	CMB-C2B-C3B	2.31	129.00	124.68
2	A	601	HEM	CMA-C3A-C4A	-2.31	124.92	128.46
2	A	601	HEM	C3C-C4C-NC	-2.17	106.84	110.94
2	A	601	HEM	CBA-CAA-C2A	-2.16	108.50	112.49
3	A	602	RIT	C15-N58-C18	-2.11	119.84	123.20
3	A	602	RIT	C26-C12-C13	-2.06	108.20	111.65

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	RIT	O7-C10-N11-C12
3	A	602	RIT	C12-C13-C14-C15
3	A	602	RIT	N58-C18-C19-C62
3	A	602	RIT	O61-C18-C19-C62
3	A	602	RIT	O24-C10-N11-C12
3	A	602	RIT	N11-C10-O7-C6
3	A	602	RIT	O24-C10-O7-C6
3	A	602	RIT	N11-C12-C26-C28
3	A	602	RIT	N20-C19-C62-C68

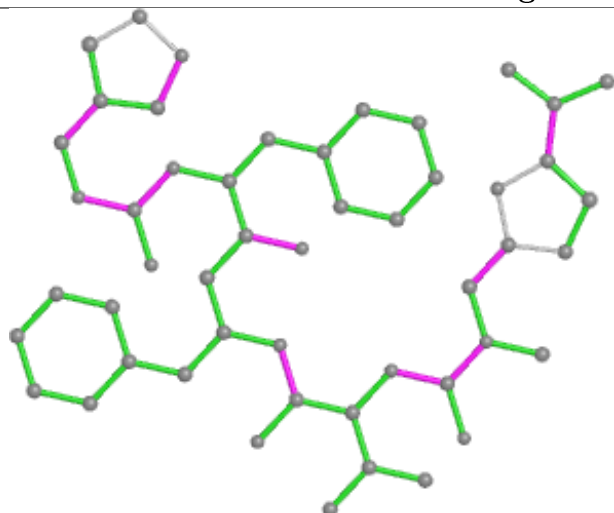
There are no ring outliers.

2 monomers are involved in 13 short contacts:

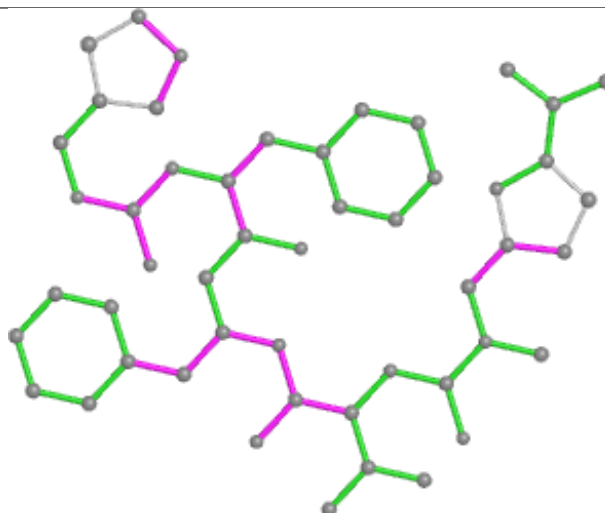
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	RIT	11	0
2	A	601	HEM	3	0

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.

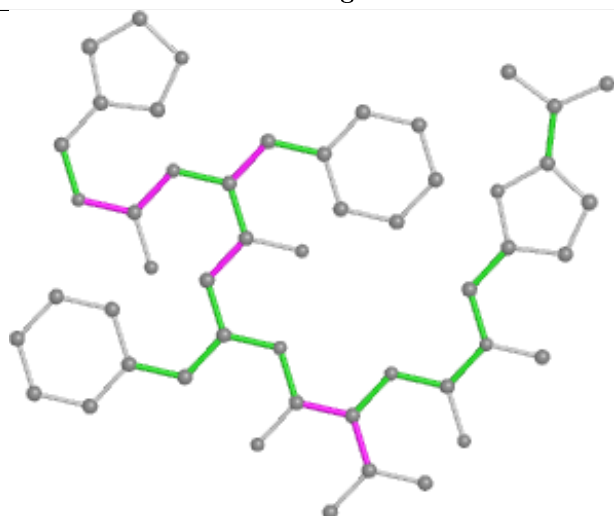
## Ligand RIT A 602



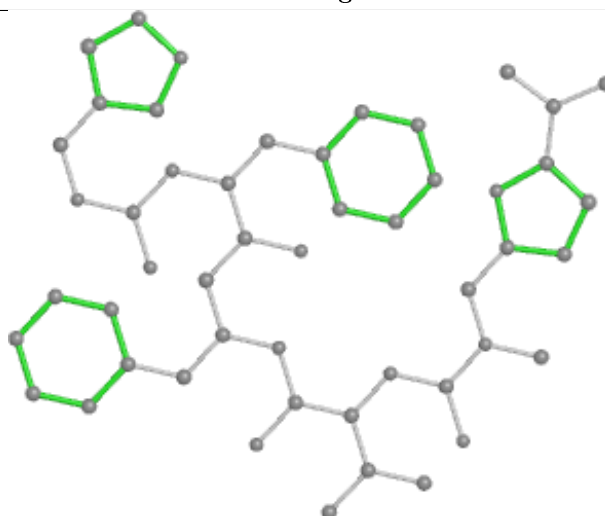
Bond lengths



Bond angles

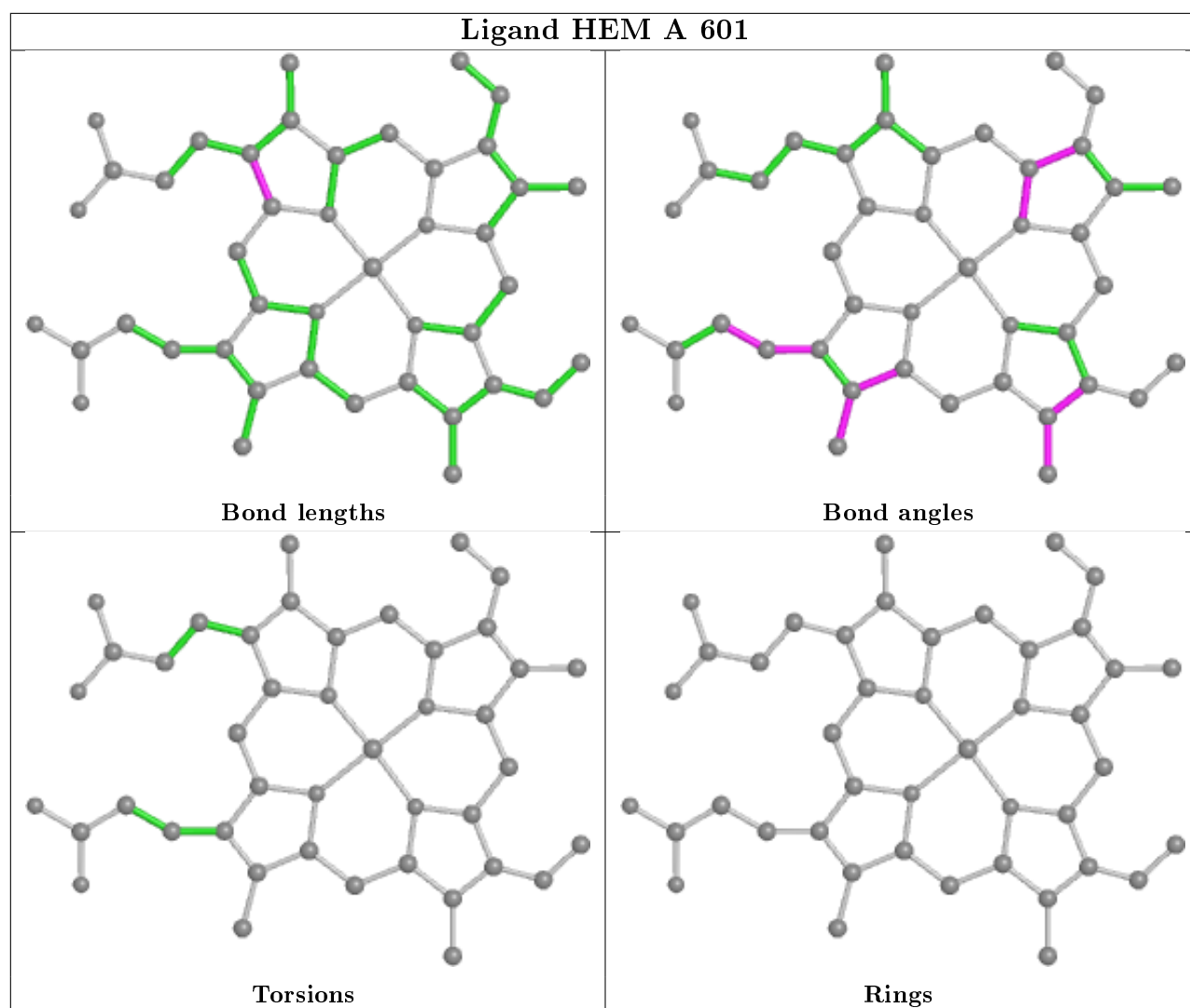


Torsions



Rings





## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	457/487 (93%)	0.59	37 (8%) 12 10	40, 68, 113, 173	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	202	PRO	9.7
1	A	251	LYS	7.9
1	A	203	PHE	7.4
1	A	215	PHE	6.4
1	A	269	VAL	6.3
1	A	250	ARG	6.1
1	A	254	LYS	6.0
1	A	289	ALA	5.5
1	A	255	ARG	5.1
1	A	189	PHE	4.9
1	A	257	LYS	4.5
1	A	55	LYS	3.6
1	A	210	LEU	3.6
1	A	28	HIS	3.5
1	A	226	PHE	3.4
1	A	207	THR	3.2
1	A	475	LEU	3.1
1	A	426	ASN	3.0
1	A	27	THR	3.0
1	A	290	LEU	3.0
1	A	57	PHE	2.7
1	A	205	GLU	2.6
1	A	150	ALA	2.6
1	A	137	PHE	2.6
1	A	304	PHE	2.6
1	A	200	GLN	2.4
1	A	256	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	191	VAL	2.4
1	A	213	PHE	2.3
1	A	149	ILE	2.3
1	A	259	SER	2.3
1	A	479	LEU	2.2
1	A	273	GLN	2.2
1	A	235	VAL	2.2
1	A	223	ILE	2.2
1	A	201	ASP	2.2
1	A	487	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

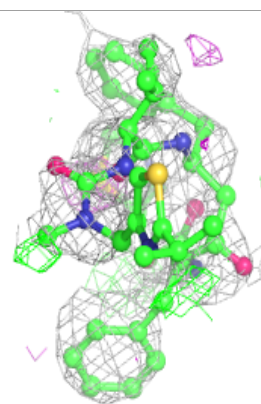
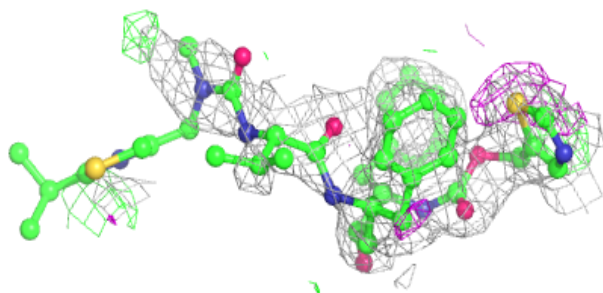
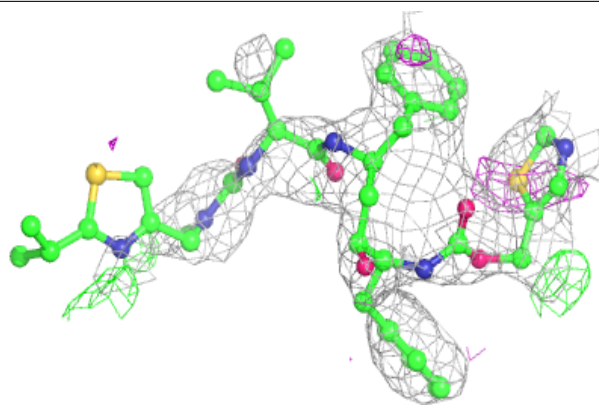
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	RIT	A	602	50/50	0.87	0.28	62,96,178,182	0
2	HEM	A	601	43/43	0.98	0.15	43,47,51,54	0

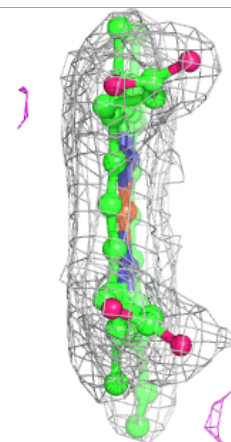
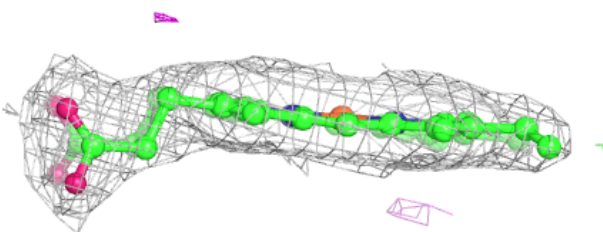
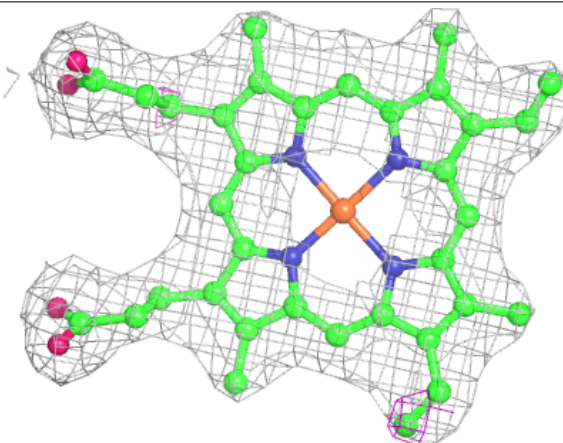
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around RIT A 602:**

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

**Electron density around 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.