



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:01 PM GMT

PDB ID : 4K9X
Title : Complex of human CYP3A4 with a desoxyritonavir analog
Authors : Sevrioukova, I.F.; Poulos, T.L.
Deposited on : 2013-04-21
Resolution : 2.76 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

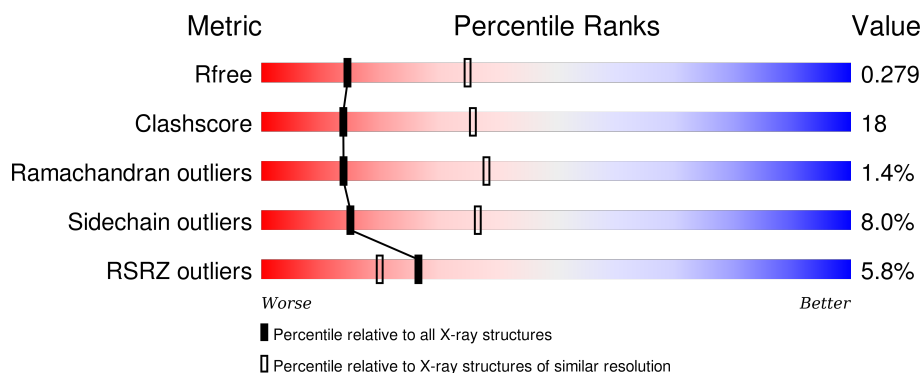
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.76 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}	91344	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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>5%</div> <div>60%</div> <div>28%</div> <div>8%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3683 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.

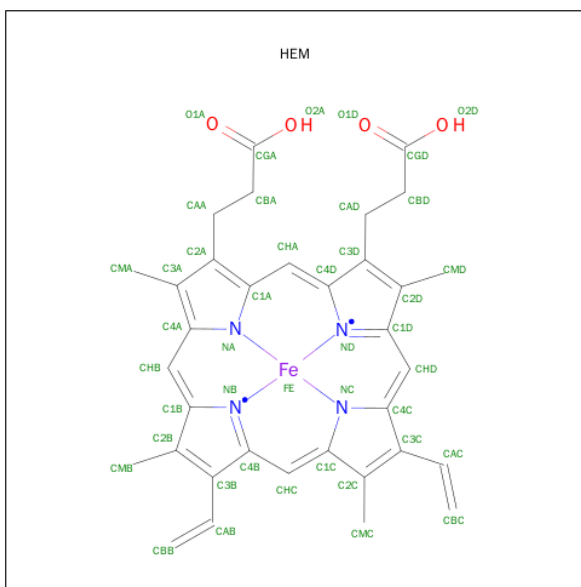
- Molecule 1 is a protein called Cytochrome P450 3A4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	449	Total	C	N	O	S	0	0	0
			3605	2346	591	644	24			

There are 24 discrepancies between the modelled and reference sequences:

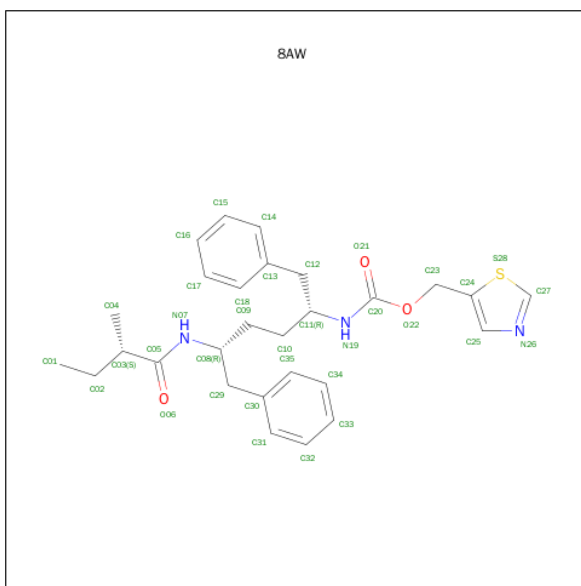
Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	LEU	DELETION	UNP P08684
A	?	-	ILE	DELETION	UNP P08684
A	?	-	PRO	DELETION	UNP P08684
A	?	-	ASP	DELETION	UNP P08684
A	?	-	LEU	DELETION	UNP P08684
A	?	-	ALA	DELETION	UNP P08684
A	?	-	MET	DELETION	UNP P08684
A	?	-	GLU	DELETION	UNP P08684
A	?	-	THR	DELETION	UNP P08684
A	?	-	TRP	DELETION	UNP P08684
A	?	-	LEU	DELETION	UNP P08684
A	?	-	LEU	DELETION	UNP P08684
A	?	-	LEU	DELETION	UNP P08684
A	?	-	ALA	DELETION	UNP P08684
A	?	-	VAL	DELETION	UNP P08684
A	?	-	SER	DELETION	UNP P08684
A	?	-	LEU	DELETION	UNP P08684
A	?	-	VAL	DELETION	UNP P08684
A	?	-	LEU	DELETION	UNP P08684
A	?	-	LEU	DELETION	UNP P08684
A	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 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is 1,3-THIAZOL-5-YLMETHYL [(2R,5R)-5-{|(2S)-2-METHYLBUTANOYL}AMINO}-1,6-DIPHENYLHEXAN-2-YL]CARBAMATE (three-letter code: 8AW) (formula: C₂₈H₃₅N₃O₃S).

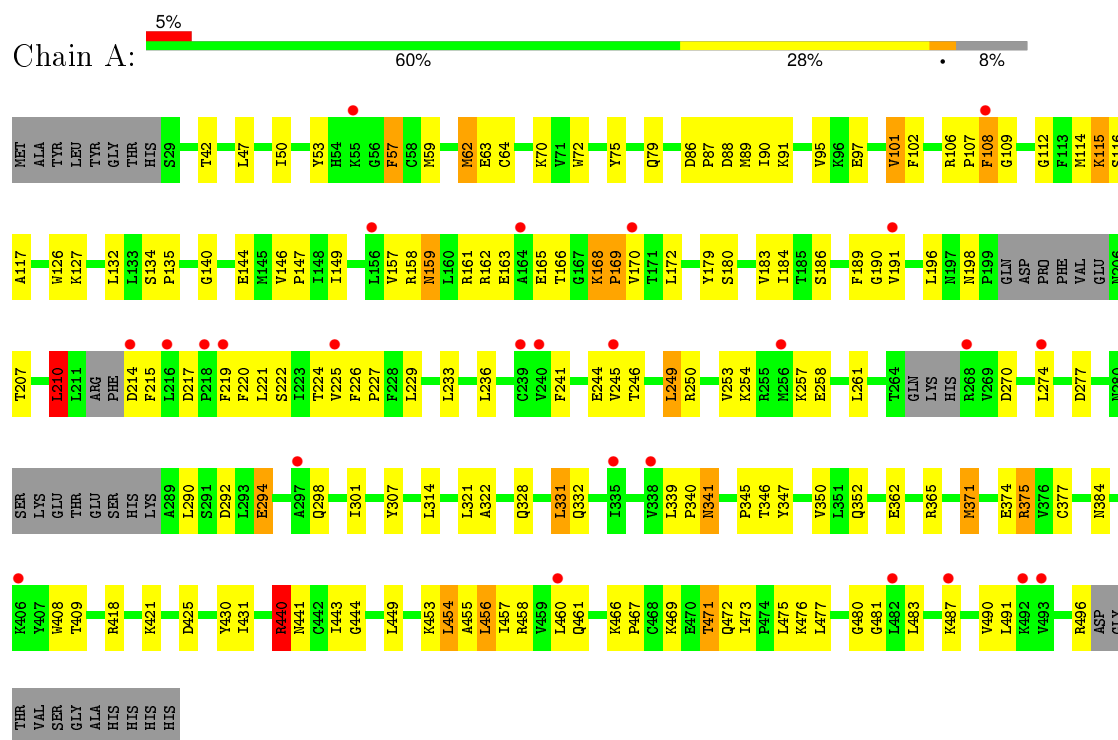


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			35	28	3	3	1		

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 errors 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, α , β , γ	77.27Å 100.33Å 126.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.17 – 2.76 50.17 – 2.76	Depositor EDS
% Data completeness (in resolution range)	100.0 (50.17-2.76) 100.0 (50.17-2.76)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.32 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.211 , 0.287 0.211 , 0.279	Depositor DCC
R_{free} test set	636 reflections (5.13%)	DCC
Wilson B-factor (Å ²)	84.7	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 62.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 13044 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3683	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.59% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, 8AW

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	0.52	1/3688 (0.0%)	0.70	1/4985 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	408	TRP	CD2-CE2	5.36	1.47	1.41

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	440	ARG	NE-CZ-NH1	-5.85	117.38	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3605	0	3697	131	0
2	A	43	0	30	9	0
3	A	35	0	35	8	0
All	All	3683	0	3762	137	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 18.

All (137) 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:50:ILE:HG12	1:A:225:VAL:HG22	1.41	0.99
2:A:601:HEM:C1D	3:A:602:8AW:H34	1.96	0.99
1:A:168:LYS:HB2	1:A:169:PRO:HD2	1.47	0.95
2:A:601:HEM:ND	3:A:602:8AW:H34	1.83	0.93
1:A:214:ASP:HB3	1:A:217:ASP:CB	2.00	0.91
1:A:53:TYR:CD2	1:A:57:PHE:HB3	2.05	0.91
1:A:159:ASN:O	1:A:163:GLU:HG2	1.71	0.91
1:A:161:ARG:HG3	1:A:161:ARG:HH11	1.34	0.90
1:A:169:PRO:O	1:A:170:VAL:HG23	1.74	0.88
1:A:108:PHE:HD2	1:A:109:GLY:H	1.21	0.86
1:A:140:GLY:O	1:A:144:GLU:HG3	1.75	0.86
1:A:170:VAL:O	1:A:490:VAL:HA	1.76	0.86
1:A:214:ASP:HB3	1:A:217:ASP:HB3	1.60	0.83
1:A:168:LYS:HB2	1:A:169:PRO:CD	2.09	0.82
1:A:210:LEU:HD12	1:A:245:VAL:HG11	1.59	0.82
1:A:53:TYR:HD2	1:A:57:PHE:HB3	1.43	0.81
1:A:170:VAL:H	1:A:490:VAL:HG13	1.47	0.79
1:A:186:SER:O	1:A:190:GLY:HA2	1.82	0.79
1:A:106:ARG:HG2	1:A:107:PRO:HD2	1.66	0.78
1:A:162:ARG:O	1:A:165:GLU:HG2	1.84	0.76
1:A:214:ASP:HB3	1:A:217:ASP:HB2	1.64	0.76
1:A:149:ILE:HG12	1:A:183:VAL:HG13	1.70	0.73
1:A:108:PHE:CD2	1:A:109:GLY:N	2.50	0.73
1:A:97:GLU:HB3	1:A:101:VAL:HG13	1.72	0.71
1:A:371:MET:HE1	1:A:483:LEU:HD13	1.73	0.70
1:A:117:ALA:HB1	1:A:301:ILE:HG13	1.73	0.70
1:A:170:VAL:O	1:A:490:VAL:HG22	1.90	0.70
1:A:112:GLY:O	1:A:115:LYS:HG2	1.93	0.68
1:A:87:PRO:HG3	1:A:431:ILE:HD11	1.76	0.68
1:A:159:ASN:ND2	1:A:162:ARG:HH21	1.92	0.68
1:A:172:LEU:HD11	1:A:491:LEU:HD12	1.75	0.67
1:A:157:VAL:HG23	1:A:158:ARG:N	2.11	0.66
1:A:159:ASN:HD22	1:A:162:ARG:HH21	1.43	0.64
1:A:375:ARG:NH2	2:A:601:HEM:CGA	2.60	0.64
1:A:47:LEU:O	1:A:50:ILE:HB	1.97	0.64
1:A:126:TRP:CZ2	1:A:440:ARG:HG2	2.33	0.63
1:A:371:MET:CE	1:A:483:LEU:HD13	2.28	0.62
1:A:249:LEU:O	1:A:253:VAL:HG23	1.98	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:ARG:NH2	2:A:601:HEM:O2A	2.32	0.62
1:A:166:THR:HB	1:A:168:LYS:HG2	1.84	0.59
1:A:375:ARG:HH22	2:A:601:HEM:CGA	2.16	0.58
1:A:458:ARG:HG3	1:A:458:ARG:HH21	1.69	0.58
1:A:294:GLU:O	1:A:298:GLN:HG2	2.05	0.57
1:A:169:PRO:O	1:A:170:VAL:CG2	2.49	0.57
1:A:161:ARG:NH1	1:A:161:ARG:HG3	2.05	0.57
1:A:179:TYR:CE1	1:A:455:ALA:HB2	2.39	0.57
1:A:472:GLN:NE2	1:A:476:LYS:H	2.01	0.57
1:A:89:MET:HA	1:A:384:ASN:ND2	2.21	0.56
1:A:168:LYS:CB	1:A:169:PRO:CD	2.80	0.56
1:A:471:THR:CG2	1:A:473:ILE:HG13	2.36	0.55
1:A:409:THR:O	1:A:418:ARG:NH2	2.38	0.55
1:A:53:TYR:CD2	1:A:57:PHE:CB	2.85	0.55
1:A:157:VAL:CG2	1:A:158:ARG:N	2.70	0.55
1:A:170:VAL:HB	1:A:491:LEU:O	2.05	0.55
1:A:91:LYS:HG3	1:A:430:TYR:CZ	2.42	0.55
1:A:157:VAL:CG2	1:A:158:ARG:H	2.20	0.55
1:A:114:MET:HG3	1:A:241:PHE:CE1	2.42	0.54
1:A:371:MET:CE	1:A:483:LEU:HD22	2.37	0.54
1:A:340:PRO:O	1:A:341:ASN:HB2	2.06	0.54
1:A:189:PHE:HB2	1:A:191:VAL:HG23	1.89	0.54
1:A:332:GLN:NE2	1:A:496:ARG:HD3	2.23	0.54
1:A:254:LYS:O	1:A:258:GLU:HG3	2.08	0.53
1:A:146:VAL:HG21	1:A:347:TYR:HB2	1.90	0.53
1:A:134:SER:OG	1:A:135:PRO:HD3	2.09	0.53
1:A:106:ARG:HG2	1:A:107:PRO:CD	2.38	0.53
1:A:472:GLN:HE22	1:A:476:LYS:H	1.57	0.53
1:A:222:SER:O	1:A:226:PHE:HB2	2.09	0.52
3:A:602:8AW:H22	3:A:602:8AW:C31	2.40	0.51
1:A:362:GLU:OE2	1:A:418:ARG:HD3	2.10	0.51
1:A:50:ILE:HG12	1:A:225:VAL:CG2	2.28	0.51
1:A:440:ARG:HD3	2:A:601:HEM:O1D	2.09	0.51
1:A:108:PHE:HD2	1:A:109:GLY:N	2.00	0.51
1:A:458:ARG:HA	1:A:461:GLN:HG2	1.91	0.51
1:A:157:VAL:HG23	1:A:158:ARG:H	1.74	0.51
1:A:346:THR:O	1:A:350:VAL:HG23	2.11	0.51
1:A:471:THR:HG23	1:A:473:ILE:HG13	1.94	0.50
1:A:86:ASP:O	1:A:90:ILE:HG13	2.11	0.50
1:A:374:GLU:C	1:A:375:ARG:HG2	2.32	0.49
1:A:116:SER:HB3	1:A:294:GLU:HG3	1.94	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:472:GLN:HE21	1:A:475:LEU:HA	1.78	0.49
1:A:75:TYR:HA	1:A:79:GLN:O	2.12	0.49
1:A:274:LEU:O	1:A:277:ASP:HB2	2.12	0.49
1:A:180:SER:HB3	1:A:307:TYR:HA	1.94	0.49
1:A:107:PRO:O	1:A:108:PHE:O	2.30	0.49
1:A:339:LEU:HD11	1:A:345:PRO:HA	1.94	0.49
1:A:146:VAL:HB	1:A:147:PRO:HD3	1.95	0.48
1:A:168:LYS:O	1:A:169:PRO:O	2.30	0.48
1:A:53:TYR:CE2	1:A:57:PHE:CB	2.96	0.48
1:A:134:SER:HA	1:A:443:ILE:HD11	1.95	0.48
1:A:371:MET:HE2	1:A:371:MET:HA	1.95	0.48
1:A:219:PHE:HD1	1:A:220:PHE:CD2	2.31	0.48
1:A:226:PHE:HA	1:A:227:PRO:HD3	1.83	0.48
1:A:221:LEU:O	1:A:225:VAL:HG23	2.14	0.47
1:A:89:MET:HA	1:A:384:ASN:HD22	1.79	0.47
1:A:210:LEU:CD2	3:A:602:8AW:H28	2.44	0.47
1:A:371:MET:HE3	1:A:483:LEU:HD22	1.96	0.47
1:A:257:LYS:O	1:A:261:LEU:HG	2.15	0.47
1:A:449:LEU:O	1:A:453:LYS:HG3	2.14	0.47
1:A:107:PRO:O	1:A:108:PHE:C	2.54	0.47
1:A:170:VAL:O	1:A:490:VAL:CA	2.55	0.46
1:A:220:PHE:O	1:A:224:THR:N	2.43	0.46
1:A:322:ALA:HB1	1:A:467:PRO:HD3	1.98	0.46
1:A:246:THR:O	1:A:250:ARG:CG	2.63	0.46
1:A:221:LEU:O	1:A:225:VAL:CG2	2.64	0.46
2:A:601:HEM:CHD	3:A:602:8AW:H34	2.44	0.45
1:A:102:PHE:HE2	1:A:377:CYS:HB2	1.81	0.45
1:A:64:CYS:HB3	1:A:72:TRP:CE2	2.52	0.45
1:A:314:LEU:HD22	1:A:456:LEU:HD13	1.99	0.45
1:A:219:PHE:HD1	1:A:220:PHE:CE2	2.35	0.44
1:A:331:LEU:HD13	1:A:460:LEU:HD12	1.98	0.44
1:A:233:LEU:HA	1:A:236:LEU:HB2	1.98	0.44
1:A:184:ILE:HA	1:A:184:ILE:HD12	1.90	0.44
1:A:444:GLY:HA3	2:A:601:HEM:C3C	2.53	0.44
1:A:89:MET:HB3	1:A:89:MET:HE2	1.90	0.44
1:A:210:LEU:HD12	1:A:245:VAL:CG1	2.40	0.44
1:A:321:LEU:O	1:A:328:GLN:HB2	2.18	0.43
1:A:169:PRO:HB3	1:A:490:VAL:CG1	2.49	0.43
3:A:602:8AW:H22	3:A:602:8AW:C30	2.49	0.42
1:A:168:LYS:HE2	1:A:168:LYS:HB2	1.92	0.42
1:A:91:LYS:O	1:A:95:VAL:HB	2.20	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:487:LYS:HA	1:A:487:LYS:HD3	1.71	0.42
1:A:207:THR:O	1:A:210:LEU:HB2	2.20	0.41
1:A:159:ASN:HD22	1:A:162:ARG:NH2	2.15	0.41
1:A:467:PRO:HB3	1:A:471:THR:HG21	2.01	0.41
1:A:375:ARG:NH2	2:A:601:HEM:O1A	2.52	0.41
1:A:270:ASP:O	1:A:274:LEU:HB2	2.20	0.41
1:A:132:LEU:HD21	1:A:290:LEU:HD22	2.02	0.41
3:A:602:8AW:C10	3:A:602:8AW:C30	2.98	0.41
1:A:371:MET:SD	1:A:483:LEU:HD13	2.60	0.41
1:A:477:LEU:HD22	1:A:483:LEU:HG	2.02	0.41
1:A:217:ASP:OD1	1:A:219:PHE:HB2	2.21	0.41
1:A:345:PRO:HB3	1:A:457:ILE:HG21	2.01	0.41
1:A:134:SER:N	1:A:135:PRO:CD	2.84	0.41
1:A:62:MET:O	1:A:63:GLU:C	2.59	0.40
1:A:441:ASN:C	1:A:441:ASN:OD1	2.59	0.40
1:A:210:LEU:HD23	3:A:602:8AW:H28	2.04	0.40
1:A:350:VAL:HG21	1:A:454:LEU:HD13	2.02	0.40

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	439/487 (90%)	387 (88%)	46 (10%)	6 (1%)	14	38

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	108	PHE
1	A	169	PRO
1	A	210	LEU
1	A	341	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	481	GLY
1	A	480	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	410/443 (93%)	377 (92%)	33 (8%)	15	37

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	THR
1	A	57	PHE
1	A	59	MET
1	A	62	MET
1	A	70	LYS
1	A	88	ASP
1	A	101	VAL
1	A	115	LYS
1	A	127	LYS
1	A	159	ASN
1	A	168	LYS
1	A	196	LEU
1	A	198	ASN
1	A	210	LEU
1	A	215	PHE
1	A	229	LEU
1	A	244	GLU
1	A	249	LEU
1	A	292	ASP
1	A	294	GLU
1	A	331	LEU
1	A	352	GLN
1	A	365	ARG
1	A	371	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	375	ARG
1	A	421	LYS
1	A	425	ASP
1	A	440	ARG
1	A	454	LEU
1	A	456	LEU
1	A	466	LYS
1	A	469	LYS
1	A	471	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	159	ASN
1	A	192	ASN
1	A	198	ASN
1	A	237	ASN
1	A	279	GLN
1	A	384	ASN
1	A	451	ASN
1	A	472	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	$\# Z > 2$
2	HEM	A	601	1,3	30,50,50	2.99	14 (46%)	24,82,82	3.20	13 (54%)
3	8AW	A	602	2	35,37,37	1.11	1 (2%)	39,47,47	1.67	8 (20%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	601	1,3	-	0/10/54/54	0/0/8/8
3	8AW	A	602	2	-	0/30/32/32	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	HEM	C3B-C4B	-4.09	1.48	1.51
2	A	601	HEM	C2D-C3D	-2.76	1.46	1.54
2	A	601	HEM	C3D-C4D	-2.65	1.48	1.51
2	A	601	HEM	C3C-CAC	-2.19	1.47	1.51
2	A	601	HEM	C2C-C1C	-2.11	1.48	1.52
2	A	601	HEM	FE-ND	2.62	2.11	1.97
2	A	601	HEM	CHD-C1D	2.72	1.46	1.38
2	A	601	HEM	C2A-C3A	2.77	1.45	1.37
2	A	601	HEM	CHC-C4B	2.83	1.46	1.38
2	A	601	HEM	FE-NB	2.89	2.12	1.97
2	A	601	HEM	CHD-C4C	3.61	1.44	1.36
2	A	601	HEM	CHC-C1C	3.98	1.45	1.36
3	A	602	8AW	O22-C20	5.01	1.45	1.35
2	A	601	HEM	C4C-NC	8.41	1.46	1.36
2	A	601	HEM	C1C-NC	8.44	1.46	1.36

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	HEM	CAA-CBA-CGA	-5.96	101.83	112.75
2	A	601	HEM	C3C-CAC-CBC	-4.80	117.09	124.46
2	A	601	HEM	C3B-CAB-CBB	-3.34	119.33	124.46
2	A	601	HEM	CBD-CAD-C3D	-3.15	104.38	113.55
3	A	602	8AW	O22-C20-O21	-2.86	118.34	124.22
2	A	601	HEM	C1D-CHD-C4C	-2.49	121.66	125.82
2	A	601	HEM	C2C-C1C-CHC	-2.33	120.13	123.68
3	A	602	8AW	C10-C11-C12	-2.19	108.09	112.05
2	A	601	HEM	C3B-C4B-NB	2.03	115.51	111.63
3	A	602	8AW	O22-C20-N19	2.11	115.22	110.54
3	A	602	8AW	C23-O22-C20	2.25	121.28	115.91
3	A	602	8AW	C23-C24-C25	3.32	135.24	128.26
2	A	601	HEM	CMD-C2D-C3D	3.37	129.26	114.35
2	A	601	HEM	C2C-C1C-NC	3.62	116.31	110.21
3	A	602	8AW	C08-N07-C05	3.75	129.00	123.18
3	A	602	8AW	C11-N19-C20	4.37	128.39	122.08
2	A	601	HEM	CAD-C3D-C4D	4.41	128.02	112.47
3	A	602	8AW	C25-N26-C27	4.47	112.76	105.71
2	A	601	HEM	CMC-C2C-C3C	5.04	129.11	116.53
2	A	601	HEM	CAD-C3D-C2D	5.89	130.16	113.22
2	A	601	HEM	CMB-C2B-C3B	6.25	132.14	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	HEM	9	0
3	A	602	8AW	8	0

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	449/487 (92%)	0.42	26 (5%)	26 20	43, 84, 137, 177	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	216	LEU	5.7
1	A	170	VAL	3.9
1	A	482	LEU	3.8
1	A	239	CYS	3.8
1	A	240	VAL	3.4
1	A	219	PHE	3.3
1	A	493	VAL	3.3
1	A	225	VAL	3.2
1	A	108	PHE	3.2
1	A	55	LYS	3.2
1	A	256	MET	3.0
1	A	492	LYS	2.7
1	A	245	VAL	2.6
1	A	218	PRO	2.5
1	A	406	LYS	2.3
1	A	191	VAL	2.3
1	A	156	LEU	2.3
1	A	338	VAL	2.2
1	A	335	ILE	2.2
1	A	164	ALA	2.2
1	A	487	LYS	2.1
1	A	297	ALA	2.1
1	A	460	LEU	2.1
1	A	268	ARG	2.1
1	A	214	ASP	2.1
1	A	274	LEU	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	HEM	A	601	43/43	0.98	0.25	1.15	51,55,60,62	0
3	8AW	A	602	35/35	0.95	0.30	0.92	49,106,133,133	0

6.5 Other polymers [i](#)

There are no such residues in this entry.