



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 10:17 am GMT

PDB ID : 4D78
Title : Cytochrome P450 3A4 bound to an inhibitor
Authors : Sevrioukova, I.; Poulos, T.
Deposited on : 2014-11-21
Resolution : 2.80 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

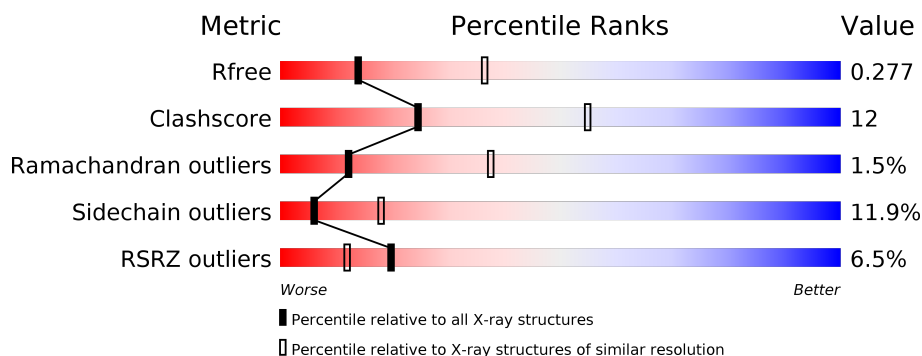
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

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>6%</div> <div>63%</div> <div>27%</div> <div>• • 6%</div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3789 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.

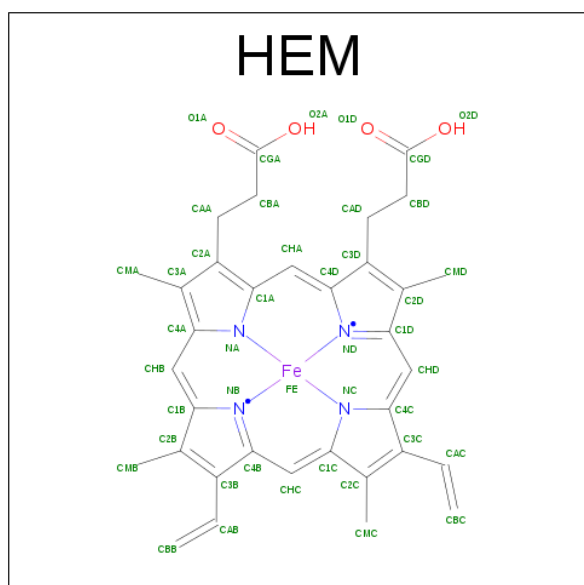
- Molecule 1 is a protein called CYTOCHROME P450 3A4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	460	3698	2406	609	659	24	0	0	1

There are 6 discrepancies between the modelled and reference sequences:

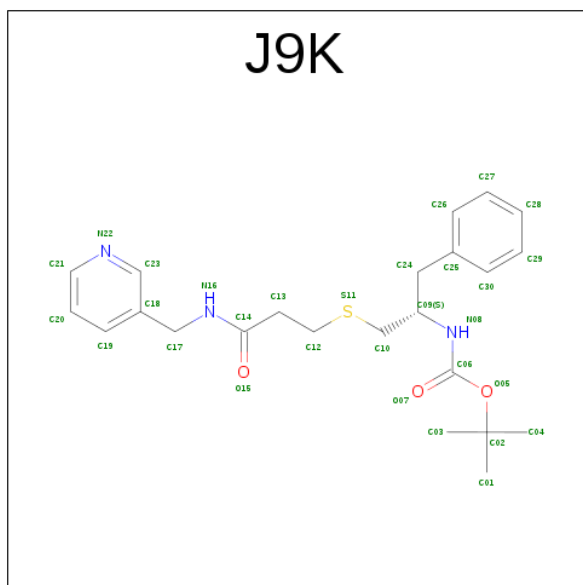
Chain	Residue	Modelled	Actual	Comment	Reference
A	21	MET	-	EXPRESSION TAG	UNP P08684
A	22	ALA	-	EXPRESSION TAG	UNP P08684
A	504	HIS	-	EXPRESSION TAG	UNP P08684
A	505	HIS	-	EXPRESSION TAG	UNP P08684
A	506	HIS	-	EXPRESSION TAG	UNP P08684
A	507	HIS	-	EXPRESSION TAG	UNP P08684

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



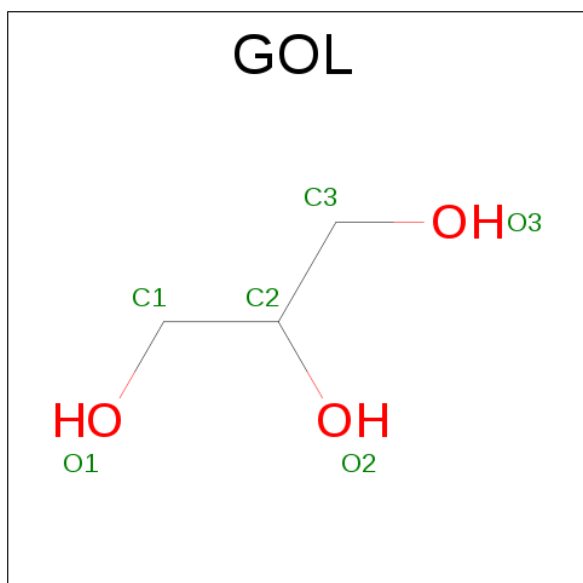
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is TERT-BUTYL [(2S)-1-({3-OXO-3-[(PYRIDIN-3-YLMETHYL)AMINO]PROPYL}SULFANYL)-3-PHENYLPROPAN-2-YL]CARBAMATE (three-letter code: J9K) (formula: $C_{23}H_{31}N_3O_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			30	23	3	3	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is water.

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

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	75.92Å 100.36Å 124.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.85 – 2.80 64.85 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (64.85-2.80) 99.8 (64.85-2.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.193 , 0.272 0.211 , 0.277	Depositor DCC
R_{free} test set	576 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	86.8	Xtriage
Anisotropy	0.305	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 82.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3789	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, J9K, 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	0.47	0/3788	0.75	5/5125 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	242	PRO	N-CA-C	10.91	140.47	112.10
1	A	222	SER	CB-CA-C	-6.05	98.59	110.10
1	A	243	ARG	N-CA-CB	5.54	120.57	110.60
1	A	368	PRO	N-CA-C	-5.35	98.19	112.10
1	A	223	ILE	N-CA-CB	5.05	122.42	110.80

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	3698	0	3779	92	0
2	A	43	0	30	2	0
3	A	30	0	31	3	0
4	A	12	0	16	1	0
5	A	6	0	0	0	0
All	All	3789	0	3856	94	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 12.

All (94) 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:58:CYS:SG	1:A:371:MET:HG2	1.78	1.22
1:A:226:PHE:HB3	1:A:229:LEU:HD12	1.28	1.09
1:A:58:CYS:SG	1:A:371:MET:CG	2.60	0.89
1:A:226:PHE:CB	1:A:229:LEU:HD12	2.04	0.86
1:A:58:CYS:HG	1:A:371:MET:HG2	1.39	0.85
1:A:210:LEU:O	1:A:211:LEU:HD23	1.82	0.79
3:A:1600:J9K:H19	3:A:1600:J9K:H13A	1.67	0.76
1:A:232:ILE:HG22	1:A:233:LEU:HD13	1.68	0.76
1:A:414:PHE:O	1:A:415:LEU:HD23	1.91	0.70
1:A:146:VAL:HG21	1:A:347:TYR:HB2	1.74	0.68
1:A:232:ILE:HG22	1:A:233:LEU:CD1	2.23	0.68
1:A:350:VAL:HG21	1:A:454:LEU:CD1	2.31	0.60
1:A:126:TRP:CZ2	1:A:440:ARG:HG2	2.37	0.59
1:A:301:ILE:HG22	1:A:302:PHE:CD1	2.38	0.58
1:A:230:ILE:HB	1:A:231:PRO:HD3	1.84	0.58
1:A:27:THR:OG1	1:A:28:HIS:N	2.37	0.58
1:A:218:PRO:O	1:A:222:SER:N	2.34	0.57
1:A:213:PHE:O	1:A:214:ASP:OD1	2.22	0.57
1:A:28:HIS:O	1:A:29:SER:HB3	2.04	0.57
1:A:183:VAL:HG21	1:A:451:ASN:HD21	1.69	0.57
1:A:270:ASP:H	1:A:273:GLN:HB3	1.70	0.57
1:A:168:LYS:HB2	1:A:169:PRO:HD2	1.87	0.56
1:A:53:TYR:CD1	1:A:57:PHE:HB3	2.40	0.56
1:A:191:VAL:HG12	1:A:193:ILE:HG23	1.86	0.56
1:A:301:ILE:HG22	1:A:302:PHE:HD1	1.69	0.56
1:A:149:ILE:C	1:A:151:GLN:H	2.09	0.56
1:A:58:CYS:HB3	1:A:399:TYR:CD1	2.42	0.55
1:A:94:LEU:HD21	1:A:396:ILE:HD13	1.89	0.55
1:A:322:ALA:HB2	1:A:491:LEU:HD13	1.88	0.54
1:A:350:VAL:HG21	1:A:454:LEU:HD13	1.90	0.54
1:A:247:ASN:HA	1:A:250:ARG:HB2	1.90	0.53
1:A:207:THR:O	1:A:210:LEU:HB2	2.08	0.53
1:A:223:ILE:C	1:A:225:VAL:H	2.11	0.53
1:A:409:THR:O	1:A:418:ARG:NH2	2.41	0.53
1:A:442:CYS:HB2	2:A:1500:HEM:NA	2.24	0.53
1:A:210:LEU:HG	1:A:245:VAL:HG11	1.91	0.53
1:A:249:LEU:HA	1:A:252:SER:HB2	1.92	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:HIS:CG	1:A:29:SER:N	2.77	0.51
1:A:229:LEU:HD23	1:A:232:ILE:HD12	1.93	0.50
1:A:246:THR:O	1:A:250:ARG:HG3	2.12	0.50
1:A:471:THR:HG22	1:A:473:ILE:HG13	1.93	0.50
1:A:368:PRO:HG2	1:A:368:PRO:O	2.12	0.50
1:A:240:VAL:HG23	1:A:241:PHE:CD2	2.48	0.49
1:A:312:SER:HB3	1:A:369:ILE:HD11	1.93	0.48
1:A:322:ALA:HB1	1:A:467:PRO:HD3	1.95	0.48
1:A:370:ALA:O	1:A:371:MET:HB2	2.13	0.48
1:A:29:SER:OG	1:A:78:GLN:HG3	2.13	0.48
1:A:158:ARG:NH2	1:A:197:ASN:OD1	2.46	0.48
1:A:241:PHE:HD1	1:A:241:PHE:O	1.97	0.47
3:A:1600:J9K:H24A	3:A:1600:J9K:O07	2.13	0.47
1:A:206:ASN:HB3	1:A:245:VAL:HG13	1.96	0.47
1:A:316:PHE:HB3	1:A:367:PHE:CD2	2.51	0.46
1:A:301:ILE:HG22	1:A:302:PHE:N	2.30	0.46
1:A:146:VAL:HB	1:A:147:PRO:HD3	1.96	0.46
1:A:414:PHE:C	1:A:415:LEU:HD23	2.36	0.45
1:A:168:LYS:CB	1:A:169:PRO:HD2	2.46	0.45
1:A:134:SER:N	1:A:135:PRO:CD	2.80	0.45
1:A:149:ILE:C	1:A:151:GLN:N	2.70	0.44
1:A:91:LYS:HG3	1:A:430:TYR:CZ	2.52	0.44
1:A:461:GLN:HA	1:A:496:ARG:HH21	1.83	0.44
1:A:374:GLU:C	1:A:375:ARG:HG2	2.38	0.43
1:A:415:LEU:HB2	1:A:418:ARG:HD2	1.99	0.43
1:A:149:ILE:O	1:A:151:GLN:N	2.52	0.43
1:A:270:ASP:N	1:A:270:ASP:OD1	2.51	0.43
1:A:403:ARG:O	1:A:405:PRO:HD3	2.19	0.43
1:A:339:LEU:HB3	1:A:343:ALA:HB3	1.99	0.42
1:A:219:PHE:HE2	1:A:240:VAL:HG12	1.84	0.42
1:A:191:VAL:HG21	1:A:256:MET:HG2	2.01	0.42
1:A:114:MET:HG3	1:A:241:PHE:CZ	2.54	0.42
1:A:168:LYS:HB2	1:A:168:LYS:HE2	1.32	0.42
1:A:155:VAL:HB	1:A:196:LEU:HD23	2.01	0.42
1:A:185:THR:O	1:A:189:PHE:HB2	2.20	0.42
1:A:260:ARG:C	1:A:262:GLU:H	2.23	0.42
1:A:184:ILE:HA	1:A:184:ILE:HD12	1.96	0.42
1:A:132:LEU:HD21	1:A:290:LEU:HD22	2.02	0.42
1:A:57:PHE:N	4:A:1602:GOL:H12	2.34	0.42
1:A:94:LEU:HD21	1:A:396:ILE:CD1	2.50	0.41
1:A:294:GLU:O	1:A:298:GLN:HG2	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:466:LYS:O	1:A:492:LYS:N	2.47	0.41
1:A:356:LEU:HD13	1:A:453:LYS:HG2	2.01	0.41
1:A:317:ILE:HD13	1:A:456:LEU:HD11	2.01	0.41
1:A:342:LYS:HB2	1:A:342:LYS:HE2	1.93	0.41
1:A:241:PHE:CD1	1:A:241:PHE:O	2.72	0.41
1:A:301:ILE:HD13	3:A:1600:J9K:H03B	2.03	0.41
1:A:237:ASN:OD1	1:A:243:ARG:HD2	2.21	0.41
1:A:316:PHE:CG	1:A:367:PHE:CD2	3.09	0.41
1:A:420:SER:O	1:A:424:LYS:N	2.47	0.41
1:A:428:ASP:HA	1:A:429:PRO:HD2	1.86	0.41
1:A:92:THR:HA	1:A:96:LYS:HB2	2.02	0.40
1:A:442:CYS:HB2	2:A:1500:HEM:C1A	2.56	0.40
1:A:97:GLU:HB3	1:A:101:VAL:HG13	2.03	0.40
1:A:344:PRO:HA	1:A:345:PRO:HD2	1.97	0.40
1:A:219:PHE:HD2	1:A:220:PHE:CD1	2.39	0.40
1:A:133:LEU:HD22	1:A:271:PHE:HE1	1.87	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	454/487 (93%)	405 (89%)	42 (9%)	7 (2%)	12	37

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	242	PRO
1	A	150	ALA
1	A	224	THR
1	A	261	LEU
1	A	28	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	219	PHE
1	A	169	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	420/443 (95%)	370 (88%)	50 (12%)	6 18

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	HIS
1	A	42	THR
1	A	50	ILE
1	A	59	MET
1	A	62	MET
1	A	70	LYS
1	A	88	ASP
1	A	93	VAL
1	A	101	VAL
1	A	116	SER
1	A	155	VAL
1	A	157	VAL
1	A	168	LYS
1	A	175	VAL
1	A	196	LEU
1	A	198	ASN
1	A	210	LEU
1	A	211	LEU
1	A	213	PHE
1	A	216	LEU
1	A	217	ASP
1	A	222	SER
1	A	223	ILE
1	A	233	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	236	LEU
1	A	246	THR
1	A	263	ASP
1	A	268	ARG
1	A	272	LEU
1	A	274	LEU
1	A	301	ILE
1	A	310	THR
1	A	312	SER
1	A	331	LEU
1	A	342	LYS
1	A	356	LEU
1	A	369	ILE
1	A	371	MET
1	A	390	LYS
1	A	425	ASP
1	A	426	ASN
1	A	437	SER
1	A	440	ARG
1	A	458	ARG
1	A	459	VAL
1	A	469	LYS
1	A	478	SER
1	A	482	LEU
1	A	490	VAL
1	A	495	SER

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	28	HIS
1	A	30	HIS
1	A	78	GLN
1	A	206	ASN
1	A	384	ASN
1	A	426	ASN
1	A	451	ASN
1	A	461	GLN
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 ⓘ

4 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	1500	1,3	28,50,50	0.93	1 (3%)	17,82,82	1.52	4 (23%)
3	J9K	A	1600	2	30,31,31	1.64	3 (10%)	38,40,40	2.51	13 (34%)
4	GOL	A	1601	-	5,5,5	0.35	0	5,5,5	0.35	0
4	GOL	A	1602	-	5,5,5	0.26	0	5,5,5	0.65	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
2	HEM	A	1500	1,3	-	0/6/54/54	0/0/8/8
3	J9K	A	1600	2	-	0/25/25/25	0/2/2/2
4	GOL	A	1601	-	-	0/4/4/4	0/0/0/0
4	GOL	A	1602	-	-	0/4/4/4	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1500	HEM	C1B-NB	-3.52	1.32	1.36
3	A	1600	J9K	O05-C06	3.67	1.42	1.34
3	A	1600	J9K	C14-N16	5.11	1.45	1.33
3	A	1600	J9K	C06-N08	5.24	1.48	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1600	J9K	C09-N08-C06	-4.75	114.89	122.22
3	A	1600	J9K	O15-C14-N16	-2.94	117.36	122.97
3	A	1600	J9K	O07-C06-N08	-2.77	120.11	124.87
2	A	1500	HEM	CBA-CAA-C2A	-2.61	107.50	112.48
2	A	1500	HEM	CMA-C3A-C4A	-2.51	124.60	128.46
2	A	1500	HEM	CBD-CAD-C3D	-2.35	107.98	112.47
3	A	1600	J9K	O05-C06-O07	-2.29	121.10	125.56
3	A	1600	J9K	C21-N22-C23	2.18	120.64	116.83
3	A	1600	J9K	C10-S11-C12	2.40	109.56	102.29
3	A	1600	J9K	C25-C24-C09	2.46	119.43	113.99
2	A	1500	HEM	CMB-C2B-C3B	2.63	129.78	124.89
3	A	1600	J9K	C24-C09-N08	2.87	116.09	110.47
3	A	1600	J9K	C02-O05-C06	3.13	126.11	121.04
3	A	1600	J9K	C13-C14-N16	4.30	123.91	116.49
3	A	1600	J9K	O05-C06-N08	4.40	117.95	110.06
3	A	1600	J9K	C17-N16-C14	5.55	131.48	122.35
3	A	1600	J9K	C18-C17-N16	8.24	131.13	112.96

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1500	HEM	2	0
3	A	1600	J9K	3	0
4	A	1602	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	460/487 (94%)	0.26	30 (6%) 20 12	49, 97, 163, 201	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	240	VAL	7.5
1	A	210	LEU	5.8
1	A	216	LEU	5.4
1	A	491	LEU	5.1
1	A	201	ASP	5.0
1	A	204	VAL	4.9
1	A	242	PRO	4.4
1	A	260	ARG	4.3
1	A	245	VAL	4.2
1	A	274	LEU	3.9
1	A	259	SER	3.7
1	A	167	GLY	3.5
1	A	241	PHE	3.5
1	A	269	VAL	3.4
1	A	196	LEU	3.1
1	A	213	PHE	2.7
1	A	239	CYS	2.7
1	A	207	THR	2.6
1	A	197	ASN	2.6
1	A	208	LYS	2.5
1	A	479	LEU	2.4
1	A	459	VAL	2.4
1	A	200	GLN	2.3
1	A	205	GLU	2.3
1	A	170	VAL	2.2
1	A	383	ILE	2.1
1	A	290	LEU	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	154	ASP	2.1
1	A	271	PHE	2.0
1	A	215	PHE	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(Å ²)	Q<0.9
4	GOL	A	1601	6/6	0.74	0.21	0.50	98,106,111,112	0
2	HEM	A	1500	43/43	0.99	0.21	0.44	44,60,66,70	0
3	J9K	A	1600	30/30	0.95	0.29	0.25	51,111,166,170	0
4	GOL	A	1602	6/6	0.94	0.25	0.24	80,85,97,99	0

6.5 Other polymers [i](#)

There are no such residues in this entry.