



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:47 AM GMT

PDB ID : 3C6G
Title : Crystal structure of CYP2R1 in complex with vitamin D3
Authors : Strushkevich, N.V.; Min, J.; Loppnau, P.; Tempel, W.; Arrowsmith, C.H.;
Edwards, A.M.; Sundstrom, M.; Weigelt, J.; Bochkarev, A.; Plotnikov, A.N.;
Park, H.; Structural Genomics Consortium (SGC)
Deposited on : 2008-02-04
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 (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

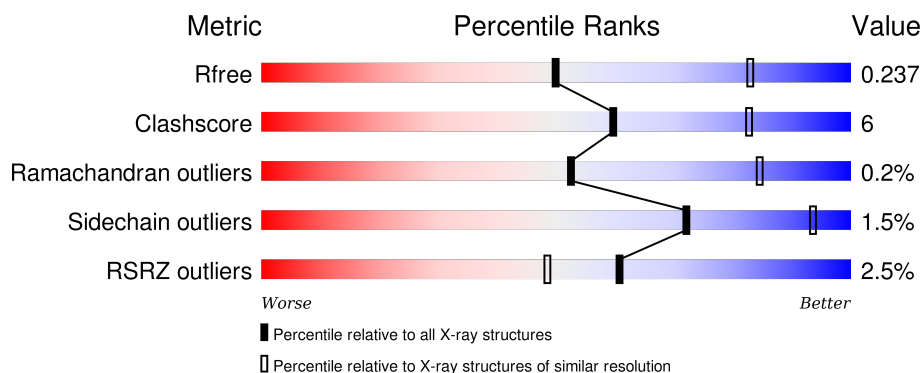
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (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	479	<div> <div>2%</div> <div>87%</div> <div>9%</div> <div>.</div> </div>
1	B	479	<div> <div>3%</div> <div>87%</div> <div>10%</div> <div>..</div> </div>

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
2	BCD	B	800	-	-	-	X
4	VD3	A	701	-	-	-	X
5	UNX	B	4	-	-	-	X
5	UNX	B	6	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7955 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.

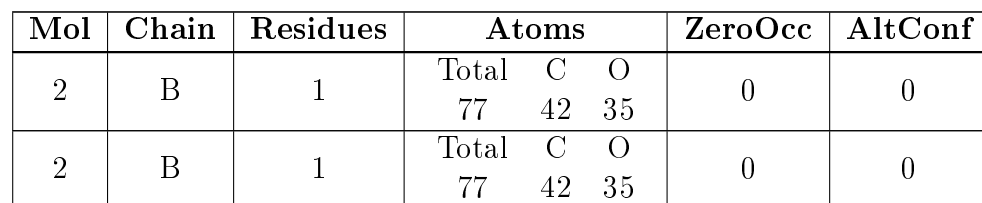
- Molecule 1 is a protein called Cytochrome P450 2R1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3763	2444	632	669	18			
1	B	467	Total	C	N	O	S	0	0	0
			3793	2462	641	672	18			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	27	MET	-	EXPRESSION TAG	UNP Q6V VX0
A	28	ALA	-	EXPRESSION TAG	UNP Q6V VX0
A	29	LYS	-	EXPRESSION TAG	UNP Q6V VX0
A	30	LYS	-	EXPRESSION TAG	UNP Q6V VX0
A	31	THR	-	EXPRESSION TAG	UNP Q6V VX0
A	502	HIS	-	EXPRESSION TAG	UNP Q6V VX0
A	503	HIS	-	EXPRESSION TAG	UNP Q6V VX0
A	504	HIS	-	EXPRESSION TAG	UNP Q6V VX0
A	505	HIS	-	EXPRESSION TAG	UNP Q6V VX0
B	27	MET	-	EXPRESSION TAG	UNP Q6V VX0
B	28	ALA	-	EXPRESSION TAG	UNP Q6V VX0
B	29	LYS	-	EXPRESSION TAG	UNP Q6V VX0
B	30	LYS	-	EXPRESSION TAG	UNP Q6V VX0
B	31	THR	-	EXPRESSION TAG	UNP Q6V VX0
B	502	HIS	-	EXPRESSION TAG	UNP Q6V VX0
B	503	HIS	-	EXPRESSION TAG	UNP Q6V VX0
B	504	HIS	-	EXPRESSION TAG	UNP Q6V VX0
B	505	HIS	-	EXPRESSION TAG	UNP Q6V VX0

- Molecule 2 is SUGAR (BETA-CYCLODEXTRIN) (three-letter code: BCD) (formula: C₄₂H₇₀O₃₅).



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- Chemical structure of HEM (heme) showing a central iron atom coordinated by four nitrogen atoms in a porphyrin-like ring, with various side chains and a central heme group.

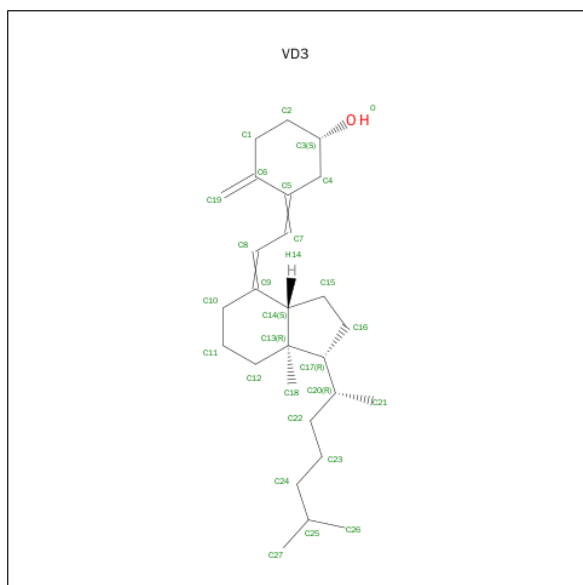
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

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

- Molecule 4 is (1S,3Z)-3-[(2E)-2-[(1R,3AR,7AS)-7A-METHYL-1-[(2R)-6-METHYLHEPTA N-2-YL]-2,3,3A,5,6,7-HEXAHYDRO-1H-INDEN-4-YLIDENE]ETHYLIDENE]-4-METHYL IDENE-CYCLOHEXAN-1-OL (three-letter code: VD3) (formula: C₂₇H₄₄O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			28	27	1		
4	B	1	Total	C	O	0	0
			28	27	1		

- Molecule 5 is UNKNOWN ATOM OR ION (three-letter code: UNX) (formula: X).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	4	Total	X	0	0
			4	4		
5	A	3	Total	X	0	0
			3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	59	Total	O	0	0
			59	59		

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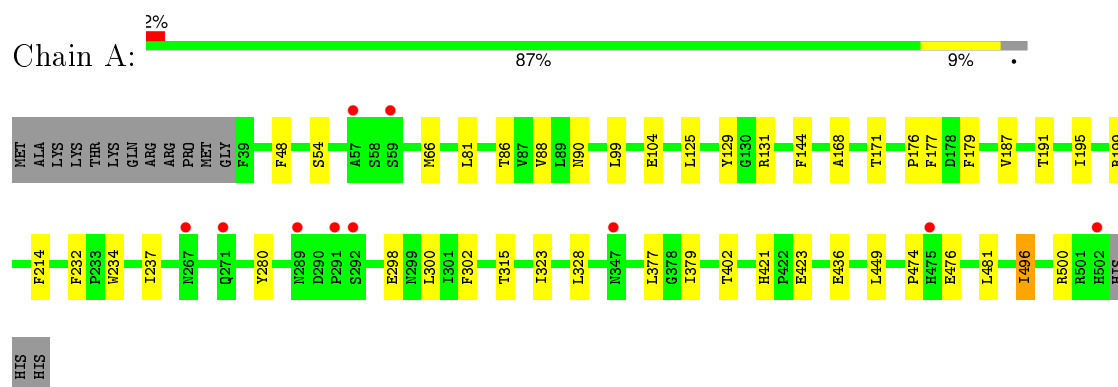
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	37	Total	O	0	0
			37	37		

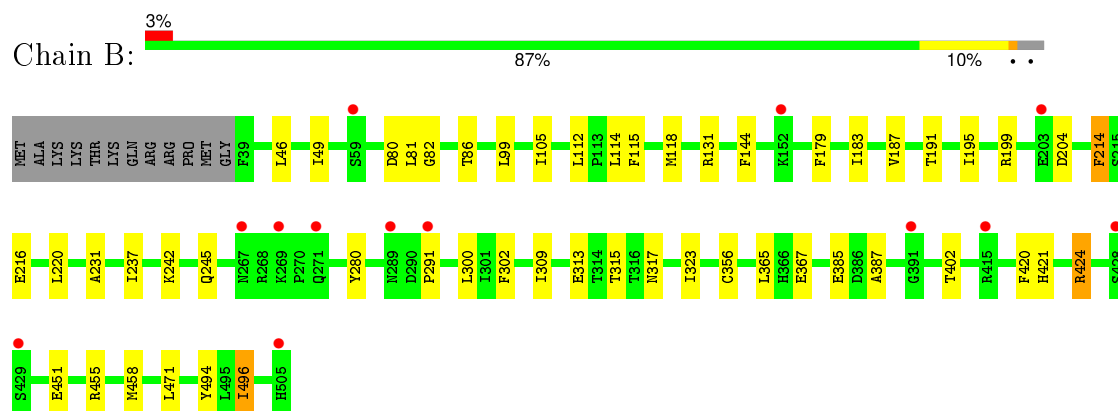
3 Residue-property plots [i](#)

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 2R1



• Molecule 1: Cytochrome P450 2R1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	137.68Å 163.05Å 152.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.34 – 2.80 29.30 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.34-2.80) 99.6 (29.30-2.80)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.77 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.3.0037	Depositor
R, R_{free}	0.193 , 0.231 0.203 , 0.237	Depositor DCC
R_{free} test set	2132 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	42.8	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 42391 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7955	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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: BCD, HEM, VD3, UNX

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	0.72	2/3872 (0.1%)	0.58	0/5247
1	B	0.67	1/3905 (0.0%)	0.58	0/5292
All	All	0.69	3/7777 (0.0%)	0.58	0/10539

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	298	GLU	CG-CD	5.21	1.59	1.51
1	A	436	GLU	CG-CD	5.12	1.59	1.51
1	B	356	CYS	CB-SG	-5.06	1.73	1.81

There are no bond angle outliers.

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	3763	0	3704	36	0
1	B	3793	0	3724	35	0
2	B	154	0	140	0	0
3	A	43	0	30	8	0
3	B	43	0	30	6	0
4	A	28	0	44	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	28	0	44	10	0
5	A	3	0	0	0	0
5	B	4	0	0	0	0
6	A	59	0	0	0	0
6	B	37	0	0	0	0
All	All	7955	0	7716	88	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 6.

All (88) 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)
4:B:700:VD3:H183	4:B:700:VD3:H212	1.39	1.02
3:A:601:HEM:HMB2	3:A:601:HEM:HBB2	1.46	0.96
3:B:601:HEM:HMB2	3:B:601:HEM:HBB2	1.56	0.88
3:A:601:HEM:HMC2	3:A:601:HEM:HBC2	1.59	0.83
3:B:601:HEM:HBC2	3:B:601:HEM:HMC2	1.62	0.81
1:B:46:LEU:HD12	1:B:49:ILE:HD11	1.68	0.76
4:B:700:VD3:H213	4:B:700:VD3:H121	1.70	0.73
3:A:601:HEM:CMB	3:A:601:HEM:HBB2	2.14	0.73
1:A:99:LEU:HD11	1:A:402:THR:HG21	1.71	0.72
1:A:191:THR:HG22	1:A:195:ILE:HD12	1.71	0.71
1:B:99:LEU:HD11	1:B:402:THR:HG21	1.72	0.70
1:B:323:ILE:HG22	1:B:496:ILE:CD1	2.22	0.67
3:B:601:HEM:HBB2	3:B:601:HEM:CMB	2.24	0.67
1:B:191:THR:HG22	1:B:195:ILE:HD12	1.78	0.66
3:A:601:HEM:HBC2	3:A:601:HEM:CMC	2.27	0.65
1:A:66:MET:HE2	1:A:88:VAL:HG11	1.77	0.65
4:A:701:VD3:H212	4:A:701:VD3:H121	1.79	0.65
4:B:700:VD3:C21	4:B:700:VD3:H121	2.28	0.63
1:A:323:ILE:HG22	1:A:496:ILE:HD12	1.81	0.62
1:A:302:PHE:HD1	4:A:701:VD3:H12	1.64	0.61
1:A:81:LEU:HD12	1:A:86:THR:HG21	1.83	0.60
1:B:302:PHE:HD1	4:B:700:VD3:H12	1.65	0.60
1:A:144:PHE:CE2	3:A:601:HEM:HBC1	2.37	0.60
1:B:81:LEU:HD12	1:B:86:THR:HG21	1.84	0.59
1:A:48:PHE:CE1	1:A:232:PHE:CE2	2.90	0.59
3:B:601:HEM:HBC2	3:B:601:HEM:CMC	2.31	0.59
1:B:220:LEU:HD22	1:B:242:LYS:HB3	1.86	0.57
1:A:280:TYR:CG	1:A:300:LEU:HD13	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:ILE:HD11	1:B:237:ILE:HD11	1.86	0.57
4:B:700:VD3:C18	4:B:700:VD3:H212	2.23	0.56
1:A:66:MET:HE2	1:A:88:VAL:CG1	2.36	0.56
1:A:302:PHE:CD1	4:A:701:VD3:H12	2.40	0.56
1:A:66:MET:CE	1:A:88:VAL:CG1	2.84	0.56
1:A:104:GLU:OE2	1:A:129:TYR:OH	2.20	0.56
1:B:144:PHE:CE2	3:B:601:HEM:HBC1	2.42	0.55
1:B:421:HIS:O	1:B:424:ARG:HB3	2.07	0.54
1:B:82:GLY:HA3	1:B:231:ALA:HB1	1.90	0.54
1:A:66:MET:CE	1:A:88:VAL:HG13	2.38	0.54
1:A:421:HIS:CE1	1:A:423:GLU:HB3	2.43	0.54
1:A:323:ILE:CG2	1:A:496:ILE:HD12	2.37	0.53
1:A:176:PRO:HB3	1:A:474:PRO:HG3	1.91	0.53
1:A:66:MET:HE3	1:A:88:VAL:HG13	1.90	0.52
1:A:449:LEU:HD23	1:A:449:LEU:O	2.10	0.52
1:A:66:MET:HE1	1:A:377:LEU:HD22	1.92	0.52
1:A:323:ILE:HG22	1:A:496:ILE:CD1	2.39	0.51
1:B:82:GLY:HA3	1:B:231:ALA:CB	2.40	0.51
1:B:451:GLU:O	1:B:455:ARG:HG3	2.11	0.50
1:A:168:ALA:O	1:A:171:THR:HB	2.12	0.50
3:A:601:HEM:HBA2	3:A:601:HEM:HHA	1.92	0.49
1:B:112:LEU:HD12	1:B:115:PHE:CE1	2.48	0.49
1:B:302:PHE:CD1	4:B:700:VD3:H12	2.44	0.49
4:B:700:VD3:C21	4:B:700:VD3:H183	2.24	0.48
1:A:191:THR:HG22	1:A:195:ILE:CD1	2.43	0.48
1:B:216:GLU:OE1	1:B:245:GLN:NE2	2.47	0.47
1:B:187:VAL:HG11	1:B:315:THR:HB	1.97	0.46
1:A:187:VAL:HG11	1:A:315:THR:HB	1.96	0.46
1:B:105:ILE:HG23	1:B:385:GLU:HG2	1.97	0.46
1:B:105:ILE:HG21	1:B:387:ALA:HB2	1.97	0.46
3:A:601:HEM:HMB2	3:A:601:HEM:CBB	2.32	0.46
1:B:323:ILE:HG23	1:B:471:LEU:HD13	1.97	0.46
1:B:280:TYR:CG	1:B:300:LEU:HD13	2.52	0.45
1:B:191:THR:CG2	1:B:195:ILE:HD12	2.46	0.45
1:B:80:ASP:OD1	1:B:82:GLY:O	2.35	0.44
1:B:313:GLU:O	1:B:317:ASN:ND2	2.50	0.44
1:B:367:GLU:HG3	1:B:420:PHE:CD1	2.51	0.44
1:B:114:LEU:CD1	1:B:118:MET:HE3	2.48	0.44
1:B:214:PHE:HB3	1:B:309:ILE:HD13	1.99	0.44
1:A:237:ILE:HD11	1:B:237:ILE:CD1	2.48	0.44
1:A:144:PHE:CZ	3:A:601:HEM:HBC1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:494:TYR:HE1	1:B:496:ILE:HD12	1.84	0.43
1:A:125:LEU:HD21	1:A:379:ILE:HD13	2.01	0.43
1:A:177:PHE:CZ	1:A:179:PHE:CE2	3.07	0.43
1:A:302:PHE:HD1	4:A:701:VD3:C1	2.31	0.43
1:B:323:ILE:CG2	1:B:496:ILE:CD1	2.94	0.42
1:A:48:PHE:CZ	1:A:232:PHE:CE2	3.07	0.42
1:A:66:MET:HE3	1:A:88:VAL:CG1	2.49	0.42
1:A:232:PHE:HB3	1:A:234:TRP:CE2	2.55	0.42
1:B:144:PHE:HE2	3:B:601:HEM:HBC1	1.85	0.41
4:B:700:VD3:C12	4:B:700:VD3:C21	2.96	0.41
1:B:323:ILE:HG22	1:B:496:ILE:HD12	2.00	0.41
1:A:88:VAL:HG12	1:A:90:ASN:HD21	1.84	0.41
1:B:114:LEU:O	1:B:118:MET:HG2	2.20	0.41
1:A:328:LEU:CD2	1:A:481:LEU:HD11	2.50	0.41
1:B:365:LEU:HD22	1:B:458:MET:CE	2.51	0.40
4:B:700:VD3:H7	4:B:700:VD3:C15	2.51	0.40
1:B:179:PHE:HB3	1:B:183:ILE:CD1	2.51	0.40
4:B:700:VD3:H8	4:B:700:VD3:H42	1.90	0.40
1:A:179:PHE:CE2	1:A:496:ILE:HG22	2.57	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	462/479 (96%)	446 (96%)	15 (3%)	1 (0%)	52	84
1	B	465/479 (97%)	450 (97%)	14 (3%)	1 (0%)	52	84
All	All	927/958 (97%)	896 (97%)	29 (3%)	2 (0%)	52	84

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	476	GLU
1	B	204	ASP

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	409/422 (97%)	403 (98%)	6 (2%)	72	93
1	B	412/422 (98%)	406 (98%)	6 (2%)	72	93
All	All	821/844 (97%)	809 (98%)	12 (2%)	72	93

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	SER
1	A	131	ARG
1	A	199	ARG
1	A	214	PHE
1	A	496	ILE
1	A	500	ARG
1	B	131	ARG
1	B	199	ARG
1	B	214	PHE
1	B	291	PRO
1	B	424	ARG
1	B	496	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	271	GLN
1	A	286	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 ⓘ

Of 13 ligands modelled in this entry, 7 are unknown - leaving 6 for Mogul analysis.

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	HEM	A	601	1	30,50,50	2.35	5 (16%)	24,82,82	2.45	12 (50%)
4	VD3	A	701	-	28,30,30	1.34	5 (17%)	36,43,43	2.13	8 (22%)
3	HEM	B	601	1	30,50,50	2.17	6 (20%)	24,82,82	2.33	11 (45%)
4	VD3	B	700	-	28,30,30	1.32	3 (10%)	36,43,43	2.31	11 (30%)
2	BCD	B	800	-	84,84,84	0.64	0	126,126,126	1.18	15 (11%)
2	BCD	B	801	-	84,84,84	0.56	0	126,126,126	1.08	9 (7%)

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	HEM	A	601	1	-	0/10/54/54	0/0/8/8
4	VD3	A	701	-	-	0/15/56/56	0/3/3/3
3	HEM	B	601	1	-	0/10/54/54	0/0/8/8

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	VD3	B	700	-	-	0/15/56/56	0/3/3/3
2	BCD	B	800	-	-	0/42/182/182	0/0/8/8
2	BCD	B	801	-	-	0/42/182/182	0/0/8/8

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	HEM	C3B-C4B	-8.24	1.44	1.51
3	B	601	HEM	C3B-C4B	-7.11	1.45	1.51
3	A	601	HEM	C3D-C4D	-6.34	1.43	1.51
3	B	601	HEM	C3D-C4D	-5.28	1.44	1.51
3	A	601	HEM	C2C-C1C	-3.78	1.45	1.52
3	B	601	HEM	C2C-C1C	-2.97	1.46	1.52
3	A	601	HEM	C2D-C1D	-2.53	1.43	1.51
3	B	601	HEM	C2D-C1D	-2.42	1.43	1.51
3	A	601	HEM	C2B-C1B	-2.21	1.44	1.51
4	A	701	VD3	C12-C13	-2.08	1.50	1.54
4	A	701	VD3	C13-C14	-2.01	1.52	1.56
4	B	700	VD3	C14-C9	2.08	1.56	1.51
4	A	701	VD3	C4-C5	2.33	1.55	1.51
4	B	700	VD3	C4-C5	2.36	1.55	1.51
4	A	701	VD3	C19-C6	2.67	1.38	1.32
4	B	700	VD3	C8-C9	2.76	1.39	1.34
4	A	701	VD3	C8-C9	2.89	1.39	1.34
3	B	601	HEM	FE-NB	2.99	2.13	1.97
3	B	601	HEM	CAA-C2A	3.37	1.57	1.52

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	700	VD3	C10-C9-C8	-9.12	112.96	125.36
4	A	701	VD3	C10-C9-C8	-8.06	114.39	125.36
4	B	700	VD3	C13-C17-C20	-3.80	112.86	119.46
4	A	701	VD3	C2-C1-C6	-3.61	103.95	111.65
4	B	700	VD3	C12-C13-C17	-3.53	110.29	116.56
3	A	601	HEM	C3B-CAB-CBB	-3.43	119.20	124.46
4	A	701	VD3	C16-C15-C14	-3.40	99.75	105.39
4	A	701	VD3	C11-C12-C13	-3.35	108.82	113.17
4	B	700	VD3	C8-C7-C5	-3.00	121.30	126.81
3	A	601	HEM	CAA-CBA-CGA	-2.93	107.38	112.75
4	A	701	VD3	C13-C17-C20	-2.73	114.73	119.46
3	B	601	HEM	CBD-CAD-C3D	-2.68	105.75	113.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	701	VD3	C12-C13-C17	-2.52	112.07	116.56
4	B	700	VD3	C16-C15-C14	-2.50	101.24	105.39
3	B	601	HEM	C3B-CAB-CBB	-2.48	120.65	124.46
4	B	700	VD3	C11-C12-C13	-2.41	110.04	113.17
4	B	700	VD3	C21-C20-C22	-2.41	106.33	110.35
3	A	601	HEM	CBD-CAD-C3D	-2.41	106.54	113.55
3	A	601	HEM	C3C-CAC-CBC	-2.30	120.92	124.46
4	A	701	VD3	C8-C7-C5	-2.24	122.70	126.81
4	B	700	VD3	C13-C14-C9	-2.21	108.89	112.85
3	A	601	HEM	CAA-C2A-C3A	-2.18	122.78	129.00
3	B	601	HEM	C3C-CAC-CBC	-2.15	121.16	124.46
4	B	700	VD3	C15-C14-C13	-2.10	102.31	104.21
2	B	800	BCD	O44-C15-C25	2.06	113.12	108.10
2	B	800	BCD	O47-C11-C21	2.08	113.16	108.10
2	B	800	BCD	C26-C36-C46	2.13	114.28	109.60
3	B	601	HEM	CBA-CAA-C2A	2.15	116.37	112.53
2	B	801	BCD	O47-C11-C21	2.16	113.37	108.10
2	B	801	BCD	O55-C55-C65	2.18	111.86	106.36
3	B	601	HEM	C2C-C1C-CHC	2.19	127.02	123.68
2	B	800	BCD	C13-O53-C53	2.20	118.02	113.75
2	B	801	BCD	O51-C51-C41	2.21	114.41	109.75
2	B	801	BCD	O57-C57-C47	2.26	114.52	109.75
2	B	800	BCD	O53-C53-C43	2.31	114.63	109.75
2	B	800	BCD	C22-C32-C42	2.35	114.76	109.60
3	A	601	HEM	CAA-C2A-C1A	2.49	129.71	127.01
2	B	800	BCD	C16-C26-C36	2.51	114.92	109.97
4	B	700	VD3	C2-C3-C4	2.52	114.80	110.32
2	B	801	BCD	C37-C47-C57	2.52	116.55	110.84
2	B	800	BCD	O56-C56-C66	2.57	112.86	106.36
2	B	800	BCD	C21-C31-C41	2.58	115.26	109.60
2	B	800	BCD	C11-O51-C51	2.74	119.07	113.75
2	B	800	BCD	C14-C24-C34	2.75	115.39	109.97
3	B	601	HEM	C2D-C3D-C4D	2.76	106.18	101.50
3	A	601	HEM	C2D-C3D-C4D	2.79	106.23	101.50
2	B	800	BCD	C11-C21-C31	2.87	115.62	109.97
2	B	801	BCD	C25-C35-C45	2.90	115.97	109.60
3	A	601	HEM	CMD-C2D-C3D	2.92	127.29	114.35
2	B	801	BCD	C27-C37-C47	2.93	116.03	109.60
2	B	800	BCD	O57-C57-C47	3.01	116.11	109.75
3	B	601	HEM	CMD-C2D-C3D	3.04	127.81	114.35
2	B	800	BCD	C24-C34-C44	3.05	116.30	109.60
2	B	800	BCD	C17-O57-C57	3.08	119.72	113.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	BCD	C15-C25-C35	3.38	116.62	109.97
2	B	801	BCD	C11-O51-C51	3.39	120.32	113.75
4	B	700	VD3	C14-C9-C8	3.80	130.21	123.25
3	A	601	HEM	CMB-C2B-C3B	3.97	126.44	116.53
3	B	601	HEM	CMC-C2C-C3C	4.08	126.71	116.53
3	B	601	HEM	CMB-C2B-C3B	4.11	126.79	116.53
3	A	601	HEM	CAD-C3D-C2D	4.24	125.42	113.22
3	B	601	HEM	CAD-C3D-C2D	4.35	125.72	113.22
3	B	601	HEM	CAD-C3D-C4D	4.43	128.08	112.47
4	A	701	VD3	C14-C9-C8	4.46	131.43	123.25
3	A	601	HEM	CAD-C3D-C4D	4.50	128.35	112.47
3	A	601	HEM	CMC-C2C-C3C	4.53	127.84	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	HEM	8	0
4	A	701	VD3	4	0
3	B	601	HEM	6	0
4	B	700	VD3	10	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	464/479 (96%)	-0.34	10 (2%) 65 54	17, 33, 57, 77	0
1	B	467/479 (97%)	-0.22	13 (2%) 56 44	21, 41, 61, 79	0
All	All	931/958 (97%)	-0.28	23 (2%) 61 48	17, 37, 60, 79	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	289	ASN	4.5
1	B	203	GLU	3.9
1	B	428	SER	3.4
1	A	291	PRO	3.4
1	A	267	ASN	3.4
1	A	57	ALA	3.2
1	A	502	HIS	3.2
1	A	289	ASN	3.1
1	B	429	SER	3.0
1	B	505	HIS	2.7
1	B	291	PRO	2.4
1	B	415	ARG	2.4
1	A	475	HIS	2.4
1	A	347	ASN	2.4
1	B	391	GLY	2.2
1	B	59	SER	2.2
1	B	271	GLN	2.2
1	B	152	LYS	2.1
1	A	292	SER	2.1
1	A	271	GLN	2.1
1	A	59	SER	2.0
1	B	267	ASN	2.0
1	B	269	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. 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
5	UNX	B	4	1/1	0.55	2.58	49.06	2,2,2,2	1
5	UNX	B	6	1/1	-0.38	2.25	35.26	2,2,2,2	1
4	VD3	A	701	28/28	0.94	0.24	2.71	31,35,37,38	0
2	BCD	B	800	77/77	0.80	0.27	2.03	62,72,76,78	0
2	BCD	B	801	77/77	0.75	0.29	1.63	85,90,99,100	0
4	VD3	B	700	28/28	0.95	0.20	1.14	28,34,39,40	0
3	HEM	B	601	43/43	0.98	0.18	-0.06	22,26,31,35	0
3	HEM	A	601	43/43	0.98	0.15	-0.37	19,24,26,28	0
5	UNX	A	3	1/1	0.38	1.20	-	2,2,2,2	1
5	UNX	B	5	1/1	0.22	1.35	-	2,2,2,2	1
5	UNX	A	2	1/1	0.55	1.61	-	2,2,2,2	1
5	UNX	B	7	1/1	0.35	1.77	-	2,2,2,2	1
5	UNX	A	1	1/1	0.61	1.62	-	2,2,2,2	1

6.5 Other polymers [i](#)

There are no such residues in this entry.