



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 07:36 pm GMT

PDB ID : 3WCU
Title : The structure of a deoxygenated 400 kda hemoglobin provides a more accurate description of the cooperative mechanism of giant hemoglobins: Deoxygenated form
Authors : Numoto, N.; Nakagawa, T.; Ohara, R.; Hasegawa, T.; Kita, A.; Yoshida, T.; Maruyama, T.; Imai, K.; Fukumori, Y.; Miki, K.
Deposited on : 2013-06-01
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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

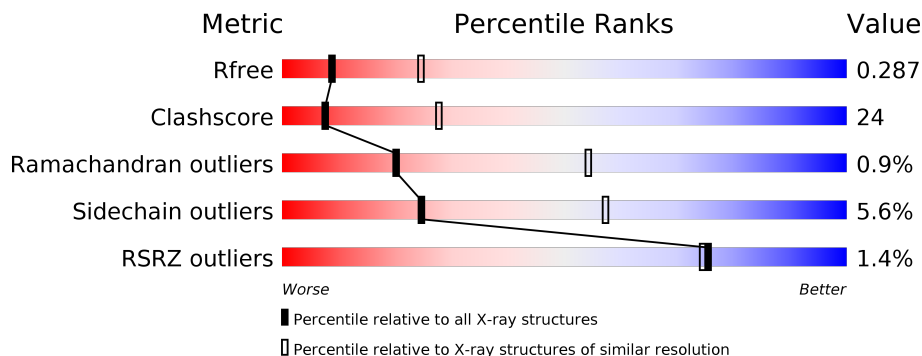
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	146	
1	E	146	
2	B	144	
2	F	144	
3	C	150	
3	G	150	

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Mol	Chain	Length	Quality of chain
4	D	149	<div><div><div>%</div><div><div></div><div>58%</div><div>39%</div><div></div></div><div></div></div></div>
4	H	149	<div><div><div>%</div><div><div></div><div>57%</div><div>40%</div><div></div></div><div></div></div></div>

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 9626 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 A1 globin chain of giant V2 hemoglobin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	0	0
			1151	736	198	209	8			
1	E	146	Total	C	N	O	S	0	0	0
			1151	736	198	209	8			

- Molecule 2 is a protein called A2 globin chain of giant V2 hemoglobin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	144	Total	C	N	O	S	0	0	0
			1120	697	206	210	7			
2	F	144	Total	C	N	O	S	0	0	0
			1120	697	206	210	7			

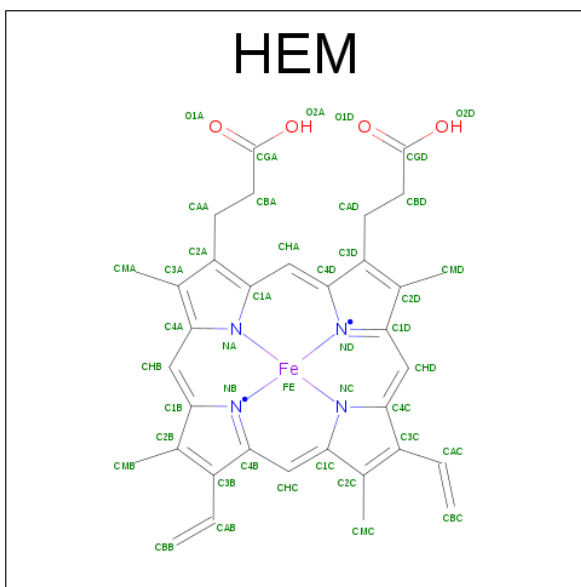
- Molecule 3 is a protein called B2 globin chain of giant V2 hemoglobin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	150	Total	C	N	O	S	0	0	0
			1158	718	211	221	8			
3	G	150	Total	C	N	O	S	0	0	0
			1158	718	211	221	8			

- Molecule 4 is a protein called B1 globin chain of giant V2 hemoglobin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	149	Total	C	N	O	S	0	0	0
			1140	727	196	210	7			
4	H	149	Total	C	N	O	S	0	0	0
			1140	727	196	210	7			

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C₃₄H₃₂FeN₄O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	E	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	F	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	G	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	H	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total Ca 1 1	0	0
6	F	1	Total Ca 1 1	0	0

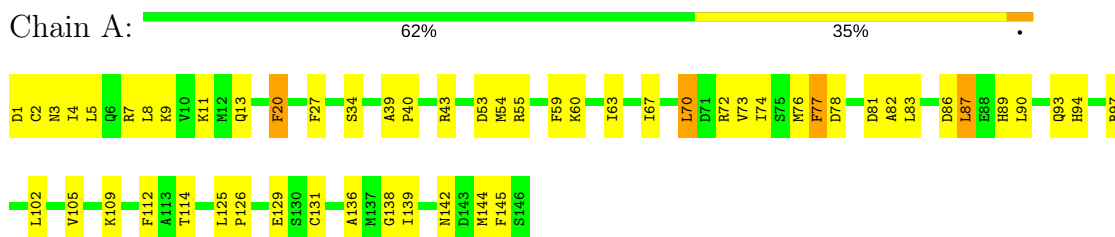
- Molecule 7 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	D	6	Total 71	C 40	N 2	O 29	0	0
7	H	6	Total 71	C 40	N 2	O 29	0	0

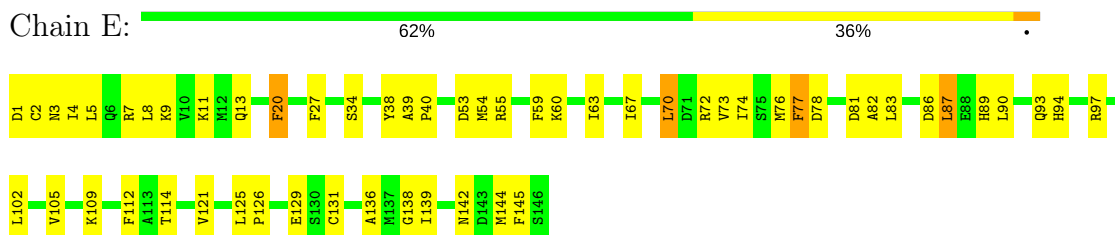
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

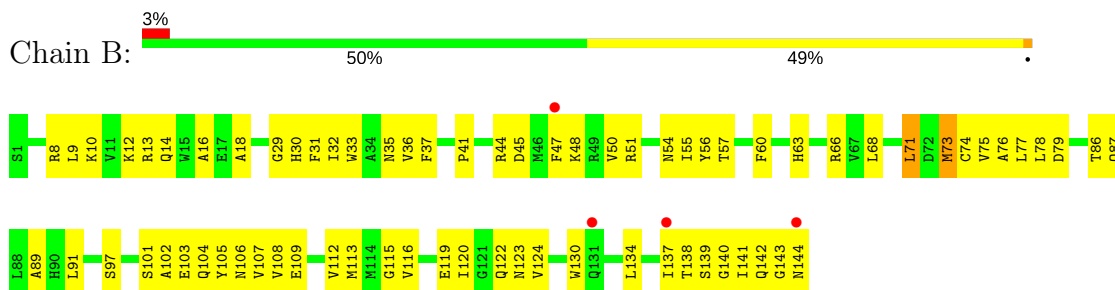
- Molecule 1: A1 globin chain of giant V2 hemoglobin



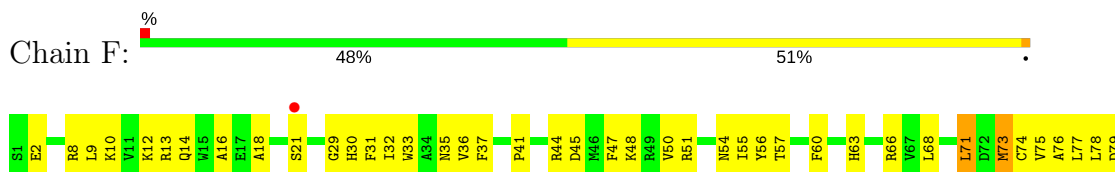
- Molecule 1: A1 globin chain of giant V2 hemoglobin

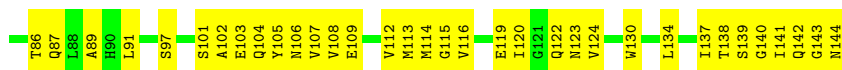


- Molecule 2: A2 globin chain of giant V2 hemoglobin

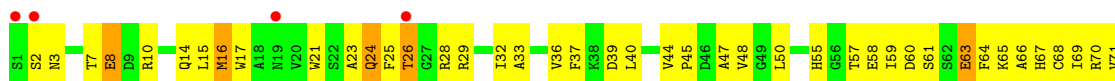


- Molecule 2: A2 globin chain of giant V2 hemoglobin





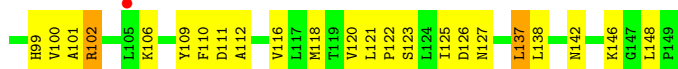
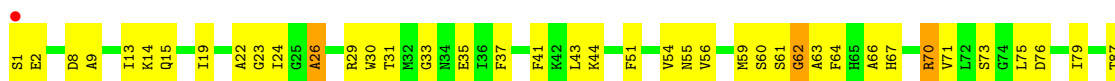
• Molecule 3: B2 globin chain of giant V2 hemoglobin



• Molecule 3: B2 globin chain of giant V2 hemoglobin



• Molecule 4: B1 globin chain of giant V2 hemoglobin



• Molecule 4: B1 globin chain of giant V2 hemoglobin



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	108.87Å 108.87Å 195.16Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.45 – 2.90 42.45 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.5 (42.45-2.90) 96.4 (42.45-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.67 (at 2.90Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.246 , 0.292 0.240 , 0.287	Depositor DCC
R_{free} test set	1437 reflections (5.41%)	DCC
Wilson B-factor (Å ²)	61.1	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 46.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.065 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9626	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% 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: FUC, HEM, CA, NAG, MAN

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	0.44	0/1178	0.60	0/1585
1	E	0.44	0/1178	0.60	0/1585
2	B	0.40	0/1143	0.59	0/1543
2	F	0.40	0/1143	0.59	0/1543
3	C	0.41	0/1181	0.63	0/1597
3	G	0.41	0/1181	0.63	0/1597
4	D	0.39	0/1164	0.60	0/1579
4	H	0.40	0/1164	0.60	0/1579
All	All	0.41	0/9332	0.60	0/12608

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 [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	1151	0	1123	50	0
1	E	1151	0	1123	50	0
2	B	1120	0	1077	75	0
2	F	1120	0	1077	76	0
3	C	1158	0	1115	81	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	1158	0	1115	83	0
4	D	1140	0	1140	59	0
4	H	1140	0	1140	61	0
5	A	43	0	30	1	0
5	B	43	0	30	1	0
5	C	43	0	30	1	0
5	D	43	0	30	3	0
5	E	43	0	30	1	0
5	F	43	0	30	1	0
5	G	43	0	30	1	0
5	H	43	0	30	3	0
6	B	1	0	0	0	0
6	F	1	0	0	0	0
7	D	71	0	61	0	0
7	H	71	0	61	0	0
All	All	9626	0	9272	461	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 24.

All (461) 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)
2:B:45:ASP:HA	2:B:48:LYS:HE2	1.47	0.95
2:F:45:ASP:HA	2:F:48:LYS:HE2	1.47	0.93
1:E:86:ASP:HB2	4:H:66:ALA:HB1	1.59	0.84
3:C:77:SER:O	3:C:81:LEU:HD23	1.78	0.83
1:A:86:ASP:HB2	4:D:66:ALA:HB1	1.59	0.83
3:G:77:SER:O	3:G:81:LEU:HD23	1.78	0.82
2:B:10:LYS:O	2:B:14:GLN:HG3	1.79	0.81
2:B:8:ARG:O	2:B:12:LYS:HG3	1.80	0.81
2:F:8:ARG:O	2:F:12:LYS:HG3	1.80	0.81
2:F:10:LYS:O	2:F:14:GLN:HG3	1.79	0.81
2:B:36:VAL:HG12	2:B:37:PHE:HD1	1.48	0.78
2:F:36:VAL:HG12	2:F:37:PHE:HD1	1.48	0.78
3:G:25:PHE:HE1	4:H:24:ILE:HA	1.49	0.76
2:B:78:LEU:HD21	2:B:137:ILE:HD11	1.67	0.76
2:F:78:LEU:HD21	2:F:137:ILE:HD11	1.67	0.76
3:C:25:PHE:HE1	4:D:24:ILE:HA	1.50	0.75
3:G:21:TRP:HE1	3:G:29:ARG:HH21	1.34	0.75
3:G:16:MET:C	3:G:16:MET:HE2	2.06	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:21:TRP:HE1	3:C:29:ARG:HH21	1.34	0.74
3:C:16:MET:HE2	3:C:16:MET:C	2.08	0.73
4:D:15:GLN:HB3	4:D:125:ILE:HD13	1.70	0.73
4:H:15:GLN:HB3	4:H:125:ILE:HD13	1.70	0.73
3:C:16:MET:HE2	3:C:16:MET:O	1.89	0.73
2:B:55:ILE:HG22	2:B:60:PHE:CE2	2.24	0.73
3:G:16:MET:HE2	3:G:16:MET:O	1.88	0.73
2:F:55:ILE:HG22	2:F:60:PHE:CE2	2.24	0.72
4:H:19:ILE:HD13	4:H:125:ILE:HG13	1.72	0.72
4:D:19:ILE:HD13	4:D:125:ILE:HG13	1.72	0.71
2:F:36:VAL:HG12	2:F:37:PHE:CD1	2.25	0.71
2:B:36:VAL:HG12	2:B:37:PHE:CD1	2.25	0.70
3:C:144:MET:SD	5:C:200:HEM:HBB2	2.31	0.70
2:F:54:ASN:OD1	2:F:57:THR:HG23	1.92	0.70
4:D:26:ALA:HA	4:D:29:ARG:HG3	1.74	0.70
2:B:54:ASN:OD1	2:B:57:THR:HG23	1.92	0.70
4:H:26:ALA:HA	4:H:29:ARG:HG3	1.74	0.70
2:B:76:ALA:O	3:C:26:THR:OG1	2.10	0.70
2:F:76:ALA:O	3:G:26:THR:OG1	2.10	0.70
3:G:144:MET:SD	5:G:200:HEM:HBB2	2.32	0.69
3:G:107:LYS:HG2	3:G:145:THR:HB	1.74	0.69
1:A:11:LYS:HE2	4:D:22:ALA:O	1.94	0.68
3:C:107:LYS:HG2	3:C:145:THR:HB	1.74	0.68
1:A:13:GLN:HE21	3:C:126:SER:HB3	1.59	0.68
1:E:13:GLN:HE21	3:G:126:SER:HB3	1.59	0.68
1:E:11:LYS:HE2	4:H:22:ALA:O	1.94	0.68
3:C:64:PHE:O	3:C:67:HIS:HB3	1.95	0.67
3:G:33:ALA:HB1	3:G:68:CYS:HB3	1.77	0.67
3:G:64:PHE:O	3:G:67:HIS:HB3	1.95	0.67
2:B:41:PRO:O	2:B:44:ARG:HG2	1.95	0.66
3:C:40:LEU:HD22	3:C:113:ILE:HG13	1.76	0.66
3:C:33:ALA:HB1	3:C:68:CYS:HB3	1.77	0.66
2:F:41:PRO:O	2:F:44:ARG:HG2	1.94	0.66
3:G:40:LEU:HD22	3:G:113:ILE:HG13	1.76	0.66
4:D:67:HIS:O	4:D:71:VAL:HG23	1.96	0.65
2:B:101:SER:OG	2:B:104:GLN:HG3	1.97	0.65
2:F:73:MET:SD	3:G:69:ILE:HG22	2.36	0.65
4:H:67:HIS:O	4:H:71:VAL:HG23	1.96	0.65
3:G:95:LEU:HD13	3:G:144:MET:HE3	1.78	0.65
2:B:73:MET:SD	3:C:69:ILE:HG22	2.36	0.65
2:F:101:SER:OG	2:F:104:GLN:HG3	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:8:LEU:HD13	3:G:28:ARG:HD3	1.79	0.64
1:A:8:LEU:HD13	3:C:28:ARG:HD3	1.79	0.64
4:H:51:PHE:HB2	4:H:56:VAL:HB	1.80	0.64
4:D:100:VAL:C	4:D:102:ARG:H	2.00	0.64
3:C:92:LEU:HD21	3:C:140:ILE:HG23	1.80	0.64
4:D:51:PHE:HB2	4:D:56:VAL:HB	1.80	0.63
1:A:3:ASN:OD1	1:A:5:LEU:HB2	1.98	0.63
1:E:3:ASN:OD1	1:E:5:LEU:HB2	1.98	0.63
3:G:92:LEU:HD21	3:G:140:ILE:HG23	1.80	0.63
4:H:118:MET:CE	4:H:137:LEU:HD12	2.29	0.62
1:A:2:CYS:SG	1:A:2:CYS:O	2.57	0.62
1:E:2:CYS:SG	1:E:2:CYS:O	2.57	0.62
4:D:118:MET:CE	4:D:137:LEU:HD12	2.29	0.62
3:G:44:VAL:HG12	3:G:47:ALA:HB2	1.80	0.61
3:C:44:VAL:HG12	3:C:47:ALA:HB2	1.80	0.61
3:G:126:SER:O	3:G:127:CYS:HB2	1.99	0.61
3:C:126:SER:O	3:C:127:CYS:HB2	1.99	0.61
2:F:124:VAL:HB	4:H:126:ASP:HB3	1.83	0.61
4:H:100:VAL:C	4:H:102:ARG:H	2.00	0.61
3:G:81:LEU:HB3	3:G:88:LEU:HB2	1.82	0.61
3:C:81:LEU:HB3	3:C:88:LEU:HB2	1.82	0.61
2:B:124:VAL:HB	4:D:126:ASP:HB3	1.82	0.61
3:C:95:LEU:HD13	3:C:144:MET:HE3	1.81	0.61
3:G:32:ILE:O	3:G:36:VAL:HG23	2.00	0.61
3:G:59:ILE:HA	3:G:64:PHE:HD2	1.65	0.61
1:A:125:LEU:O	1:A:129:GLU:HG3	2.02	0.60
1:E:125:LEU:O	1:E:129:GLU:HG3	2.02	0.60
3:C:59:ILE:HA	3:C:64:PHE:HD2	1.65	0.60
3:C:32:ILE:O	3:C:36:VAL:HG23	2.00	0.60
3:C:121:MET:N	3:C:122:PRO:HD2	2.17	0.60
3:C:135:ARG:NH2	1:E:55:ARG:NH2	2.50	0.60
3:G:121:MET:N	3:G:122:PRO:HD2	2.17	0.59
3:C:66:ALA:O	3:C:70:ARG:HG3	2.02	0.59
2:F:86:THR:HG21	3:G:63:GLU:HA	1.83	0.59
2:B:86:THR:HG21	3:C:63:GLU:HA	1.83	0.59
1:A:81:ASP:HB3	4:D:62:GLY:HA3	1.85	0.59
1:E:60:LYS:HE3	4:H:87:THR:OG1	2.01	0.59
1:A:60:LYS:HE3	4:D:87:THR:OG1	2.01	0.59
2:F:76:ALA:HB2	3:G:29:ARG:NH1	2.17	0.59
3:C:55:HIS:CE1	3:C:63:GLU:HG3	2.38	0.59
1:E:81:ASP:HB3	4:H:62:GLY:HA3	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:66:ALA:O	3:G:70:ARG:HG3	2.02	0.59
3:G:55:HIS:CE1	3:G:63:GLU:HG3	2.38	0.58
2:B:76:ALA:HB2	3:C:29:ARG:NH1	2.17	0.58
3:G:39:ASP:OD2	3:G:116:SER:HA	2.03	0.58
2:F:105:TYR:O	2:F:109:GLU:HG3	2.03	0.58
2:B:105:TYR:O	2:B:109:GLU:HG3	2.03	0.58
3:C:25:PHE:CE1	4:D:24:ILE:HA	2.36	0.58
3:C:39:ASP:OD2	3:C:116:SER:HA	2.03	0.58
3:G:21:TRP:HE1	3:G:29:ARG:NH2	2.02	0.57
3:C:58:GLU:HB3	3:C:61:SER:HB3	1.87	0.57
3:G:58:GLU:HB3	3:G:61:SER:HB3	1.87	0.57
3:G:21:TRP:NE1	3:G:29:ARG:NH2	2.48	0.57
3:C:21:TRP:HE1	3:C:29:ARG:NH2	2.01	0.57
1:E:74:ILE:HA	1:E:77:PHE:CD2	2.40	0.57
2:F:45:ASP:O	2:F:48:LYS:HG2	2.05	0.57
3:G:139:ARG:HG3	3:G:139:ARG:HH11	1.70	0.57
1:A:74:ILE:HA	1:A:77:PHE:CD2	2.40	0.56
2:B:45:ASP:O	2:B:48:LYS:HG2	2.05	0.56
1:A:43:ARG:NH2	3:G:131:ASP:OD1	2.26	0.56
3:C:95:LEU:HD13	3:C:144:MET:CE	2.36	0.56
3:C:21:TRP:NE1	3:C:29:ARG:NH2	2.48	0.56
3:G:95:LEU:HD13	3:G:144:MET:CE	2.36	0.56
2:F:76:ALA:CB	3:G:29:ARG:HH11	2.18	0.56
3:C:139:ARG:HG3	3:C:139:ARG:HH11	1.70	0.56
4:D:9:ALA:O	4:D:13:ILE:HG13	2.06	0.56
2:B:76:ALA:CB	3:C:29:ARG:HH11	2.18	0.56
3:C:7:THR:HA	3:C:10:ARG:HD2	1.88	0.56
3:C:67:HIS:HA	3:C:70:ARG:CZ	2.36	0.55
1:E:125:LEU:HB3	1:E:126:PRO:HD3	1.88	0.55
3:G:25:PHE:CE1	4:H:24:ILE:HA	2.36	0.55
4:H:9:ALA:O	4:H:13:ILE:HG13	2.06	0.55
1:A:20:PHE:CD2	1:A:20:PHE:N	2.73	0.55
2:B:29:GLY:HA2	2:B:68:LEU:HD11	1.89	0.55
1:A:125:LEU:HB3	1:A:126:PRO:HD3	1.88	0.55
3:G:7:THR:HA	3:G:10:ARG:HD2	1.88	0.55
3:C:25:PHE:CE1	4:D:23:GLY:O	2.60	0.55
2:F:29:GLY:HA2	2:F:68:LEU:HD11	1.89	0.55
1:E:20:PHE:N	1:E:20:PHE:CD2	2.73	0.55
3:G:67:HIS:HA	3:G:70:ARG:CZ	2.36	0.55
4:D:61:SER:O	4:D:63:ALA:N	2.40	0.55
3:G:25:PHE:CE1	4:H:23:GLY:O	2.60	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:61:SER:O	4:H:63:ALA:N	2.40	0.55
3:G:55:HIS:CG	3:G:63:GLU:HG3	2.42	0.55
2:B:55:ILE:HA	2:B:60:PHE:CD2	2.42	0.54
3:C:55:HIS:CG	3:C:63:GLU:HG3	2.42	0.54
1:E:83:LEU:O	1:E:87:LEU:HB2	2.07	0.54
3:G:2:SER:O	3:G:3:ASN:HB2	2.08	0.54
3:C:106:THR:HB	3:C:150:GLU:O	2.07	0.54
3:G:106:THR:HB	3:G:150:GLU:O	2.07	0.54
2:F:55:ILE:HA	2:F:60:PHE:CD2	2.42	0.54
1:A:83:LEU:O	1:A:87:LEU:HB2	2.07	0.53
2:F:13:ARG:NH1	4:H:126:ASP:OD1	2.40	0.53
2:B:31:PHE:CE2	2:B:115:GLY:HA2	2.44	0.53
2:B:45:ASP:CA	2:B:48:LYS:HE2	2.31	0.53
3:C:2:SER:O	3:C:3:ASN:HB2	2.08	0.53
2:F:45:ASP:CA	2:F:48:LYS:HE2	2.31	0.53
4:H:100:VAL:C	4:H:102:ARG:N	2.61	0.53
2:F:31:PHE:CE2	2:F:115:GLY:HA2	2.44	0.53
1:A:89:HIS:O	1:A:93:GLN:HG3	2.09	0.52
1:E:89:HIS:O	1:E:93:GLN:HG3	2.09	0.52
2:F:76:ALA:CB	3:G:29:ARG:NH1	2.73	0.52
4:D:75:LEU:O	4:D:75:LEU:HG	2.09	0.52
1:A:27:PHE:CE2	1:A:114:THR:HG22	2.45	0.52
3:C:113:ILE:O	3:C:116:SER:HB3	2.10	0.52
3:G:113:ILE:O	3:G:116:SER:HB3	2.10	0.52
1:A:43:ARG:HH22	3:G:131:ASP:CG	2.13	0.52
1:A:9:LYS:HD2	3:C:123:GLN:O	2.10	0.52
4:D:106:LYS:H	4:D:109:TYR:HD2	1.57	0.52
1:E:9:LYS:HD2	3:G:123:GLN:O	2.10	0.51
4:H:75:LEU:O	4:H:75:LEU:HG	2.09	0.51
2:B:76:ALA:CB	3:C:29:ARG:NH1	2.73	0.51
2:F:51:ARG:HG3	2:F:51:ARG:HH11	1.75	0.51
4:H:106:LYS:H	4:H:109:TYR:HD2	1.58	0.51
2:B:143:GLY:O	2:B:144:ASN:CB	2.59	0.51
4:D:100:VAL:C	4:D:102:ARG:N	2.61	0.51
4:D:118:MET:HE2	4:D:137:LEU:HD12	1.91	0.51
1:A:54:MET:HA	1:A:59:PHE:CD1	2.46	0.51
1:E:27:PHE:CE2	1:E:114:THR:HG22	2.45	0.51
3:G:55:HIS:ND1	3:G:63:GLU:HG3	2.26	0.51
3:C:55:HIS:ND1	3:C:63:GLU:HG3	2.26	0.51
1:E:77:PHE:CE1	1:E:131:CYS:HB3	2.46	0.51
1:A:77:PHE:CE1	1:A:131:CYS:HB3	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:16:ALA:HB2	3:C:23:ALA:HB3	1.92	0.51
2:B:51:ARG:HG3	2:B:51:ARG:HH11	1.75	0.51
3:C:16:MET:HE1	3:C:17:TRP:HA	1.93	0.51
2:B:13:ARG:NH1	4:D:126:ASP:OD1	2.40	0.51
1:E:139:ILE:HD12	5:E:200:HEM:HBB2	1.93	0.51
1:E:74:ILE:HG12	1:E:77:PHE:HE2	1.76	0.51
1:A:74:ILE:HG12	1:A:77:PHE:HE2	1.76	0.50
2:F:143:GLY:O	2:F:144:ASN:CB	2.59	0.50
2:F:66:ARG:HD3	3:G:91:GLN:OE1	2.11	0.50
1:A:139:ILE:HD12	5:A:200:HEM:HBB2	1.93	0.50
1:E:7:ARG:HD3	1:E:77:PHE:O	2.12	0.50
2:F:16:ALA:HB2	3:G:23:ALA:HB3	1.92	0.50
1:E:87:LEU:HD22	1:E:138:GLY:HA3	1.94	0.50
1:E:142:ASN:OD1	1:E:144:MET:HB2	2.12	0.50
2:F:18:ALA:HB2	2:F:120:ILE:HG23	1.92	0.50
2:B:124:VAL:HG12	4:D:127:ASN:HB2	1.93	0.50
4:H:66:ALA:O	4:H:70:ARG:HG2	2.11	0.50
1:A:7:ARG:HD3	1:A:77:PHE:O	2.12	0.50
4:H:109:TYR:O	4:H:112:ALA:HB3	2.11	0.50
4:H:51:PHE:HB3	4:H:54:VAL:CG2	2.41	0.50
4:D:51:PHE:HB3	4:D:54:VAL:CG2	2.41	0.50
4:D:66:ALA:O	4:D:70:ARG:HG2	2.11	0.50
1:E:54:MET:HA	1:E:59:PHE:CD1	2.46	0.50
2:F:124:VAL:HG12	4:H:127:ASN:HB2	1.93	0.50
1:A:142:ASN:OD1	1:A:144:MET:HB2	2.12	0.49
4:D:109:TYR:O	4:D:112:ALA:HB3	2.11	0.49
2:F:47:PHE:HB3	2:F:50:VAL:HG23	1.94	0.49
2:B:18:ALA:HB2	2:B:120:ILE:HG23	1.92	0.49
2:B:47:PHE:HB3	2:B:50:VAL:HG23	1.94	0.49
4:D:41:PHE:HE2	4:D:51:PHE:HE1	1.60	0.49
1:E:13:GLN:NE2	3:G:126:SER:HB3	2.27	0.49
1:A:87:LEU:HD22	1:A:138:GLY:HA3	1.94	0.49
1:A:13:GLN:NE2	3:C:126:SER:HB3	2.27	0.49
1:E:136:ALA:HB1	1:E:145:PHE:CD1	2.47	0.49
2:B:138:THR:C	2:B:140:GLY:H	2.16	0.49
1:A:136:ALA:HB1	1:A:145:PHE:CD1	2.47	0.49
2:B:66:ARG:HD3	3:C:91:GLN:OE1	2.11	0.49
2:F:138:THR:C	2:F:140:GLY:H	2.16	0.49
4:H:30:TRP:CZ2	4:H:60:SER:HA	2.48	0.49
3:C:139:ARG:NH1	3:C:139:ARG:HG3	2.28	0.49
4:D:30:TRP:CZ2	4:D:60:SER:HA	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:118:MET:HE1	4:H:137:LEU:HD12	1.94	0.49
2:B:31:PHE:CD2	2:B:115:GLY:HA2	2.48	0.48
2:B:87:GLN:HE22	3:C:70:ARG:HD3	1.78	0.48
4:D:100:VAL:O	4:D:102:ARG:N	2.47	0.48
2:F:87:GLN:HE22	3:G:70:ARG:HD3	1.78	0.48
4:H:100:VAL:O	4:H:102:ARG:N	2.47	0.48
4:H:61:SER:C	4:H:63:ALA:N	2.66	0.48
2:B:76:ALA:HB2	3:C:29:ARG:HH11	1.78	0.48
4:D:61:SER:C	4:D:63:ALA:N	2.66	0.48
3:G:139:ARG:HG3	3:G:139:ARG:NH1	2.28	0.48
3:C:115:GLN:HA	3:C:115:GLN:HE21	1.79	0.48
2:F:143:GLY:O	2:F:144:ASN:HB2	2.13	0.48
3:G:16:MET:HE1	3:G:17:TRP:HA	1.96	0.48
1:A:90:LEU:HD13	1:A:139:ILE:HD11	1.95	0.47
1:A:27:PHE:CZ	1:A:114:THR:HG22	2.49	0.47
2:B:143:GLY:O	2:B:144:ASN:HB2	2.13	0.47
1:A:78:ASP:OD1	4:D:26:ALA:HB2	2.14	0.47
2:F:31:PHE:CD2	2:F:115:GLY:HA2	2.48	0.47
2:F:76:ALA:HB2	3:G:29:ARG:HH11	1.78	0.47
3:C:25:PHE:HB3	3:C:28:ARG:HB2	1.96	0.47
2:F:37:PHE:HB3	2:F:44:ARG:HH21	1.79	0.47
1:E:78:ASP:OD1	4:H:26:ALA:HB2	2.14	0.47
1:E:90:LEU:HD13	1:E:139:ILE:HD11	1.95	0.47
1:E:27:PHE:CZ	1:E:114:THR:HG22	2.49	0.47
2:F:60:PHE:O	2:F:63:HIS:HB3	2.14	0.47
4:H:41:PHE:HE2	4:H:51:PHE:HE1	1.60	0.47
1:A:77:PHE:CZ	1:A:131:CYS:HB3	2.50	0.47
2:B:60:PHE:O	2:B:63:HIS:HB3	2.14	0.47
3:G:25:PHE:HB3	3:G:28:ARG:HB2	1.96	0.47
3:G:37:PHE:HE2	3:G:68:CYS:HA	1.80	0.47
3:C:37:PHE:HE2	3:C:68:CYS:HA	1.80	0.47
4:H:116:VAL:O	4:H:120:VAL:HG23	2.15	0.47
3:G:61:SER:HB2	3:G:63:GLU:HG2	1.97	0.47
2:B:37:PHE:HB3	2:B:44:ARG:HH21	1.79	0.47
4:H:33:GLY:O	4:H:37:PHE:HD2	1.98	0.47
4:D:71:VAL:HG13	5:D:201:HEM:C1B	2.50	0.47
4:D:33:GLY:O	4:D:37:PHE:HD2	1.98	0.47
2:B:141:ILE:CD1	5:B:201:HEM:HAB	2.44	0.46
4:H:100:VAL:HG22	4:H:148:LEU:HD23	1.98	0.46
3:G:115:GLN:HE21	3:G:115:GLN:HA	1.79	0.46
4:H:121:LEU:N	4:H:122:PRO:HD2	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:100:GLN:HG2	3:C:147:GLY:O	2.16	0.46
3:C:37:PHE:HZ	3:C:71:VAL:HG21	1.80	0.46
3:C:59:ILE:HA	3:C:64:PHE:CD2	2.49	0.46
3:C:61:SER:HB2	3:C:63:GLU:HG2	1.97	0.46
2:F:141:ILE:CD1	5:F:201:HEM:HAB	2.44	0.46
2:F:33:TRP:C	2:F:35:ASN:H	2.18	0.46
2:F:63:HIS:HA	2:F:66:ARG:CZ	2.46	0.46
3:G:100:GLN:HG2	3:G:147:GLY:O	2.16	0.46
2:B:30:HIS:HE1	2:B:56:TYR:CE1	2.33	0.46
1:E:77:PHE:CZ	1:E:131:CYS:HB3	2.50	0.46
1:E:72:ARG:HG3	1:E:73:VAL:N	2.30	0.46
3:G:37:PHE:HZ	3:G:71:VAL:HG21	1.80	0.46
1:A:105:VAL:O	1:A:109:LYS:HD3	2.16	0.46
1:E:105:VAL:O	1:E:109:LYS:HD3	2.16	0.46
2:F:137:ILE:O	2:F:141:ILE:HG13	2.16	0.46
2:F:32:ILE:O	2:F:35:ASN:HB3	2.16	0.46
2:F:30:HIS:HE1	2:F:56:TYR:CE1	2.33	0.46
1:E:1:ASP:CG	1:E:2:CYS:H	2.19	0.46
1:A:1:ASP:CG	1:A:2:CYS:H	2.19	0.46
4:H:118:MET:HE2	4:H:137:LEU:HD12	1.96	0.46
1:A:125:LEU:CD2	1:E:112:PHE:HB2	2.46	0.46
1:A:72:ARG:HG3	1:A:73:VAL:N	2.30	0.46
2:B:86:THR:O	2:B:86:THR:HG22	2.15	0.46
2:F:86:THR:O	2:F:86:THR:HG22	2.15	0.46
4:H:2:GLU:HA	4:H:2:GLU:OE1	2.16	0.46
1:A:94:HIS:HA	1:A:97:ARG:HD2	1.97	0.46
4:D:116:VAL:O	4:D:120:VAL:HG23	2.15	0.46
4:H:71:VAL:HG13	5:H:201:HEM:C1B	2.50	0.46
2:B:63:HIS:HA	2:B:66:ARG:CZ	2.46	0.45
4:D:121:LEU:N	4:D:122:PRO:HD2	2.30	0.45
1:E:94:HIS:HA	1:E:97:ARG:HD2	1.97	0.45
2:F:12:LYS:NZ	2:F:12:LYS:HB3	2.31	0.45
4:H:43:LEU:HD12	4:H:116:VAL:CG2	2.46	0.45
4:H:51:PHE:O	4:H:56:VAL:HB	2.16	0.45
4:H:55:ASN:OD1	4:H:55:ASN:O	2.34	0.45
4:H:13:ILE:HG12	4:H:79:ILE:O	2.16	0.45
4:D:13:ILE:HG12	4:D:79:ILE:O	2.16	0.45
4:D:51:PHE:O	4:D:56:VAL:HB	2.16	0.45
2:B:12:LYS:HB3	2:B:12:LYS:NZ	2.31	0.45
2:B:33:TRP:C	2:B:35:ASN:H	2.18	0.45
2:B:32:ILE:O	2:B:35:ASN:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:100:VAL:HG22	4:D:148:LEU:HD23	1.97	0.45
4:D:55:ASN:OD1	4:D:55:ASN:O	2.34	0.45
2:F:103:GLU:O	2:F:107:VAL:HG23	2.17	0.45
3:G:16:MET:CE	3:G:17:TRP:HA	2.46	0.45
2:B:137:ILE:O	2:B:141:ILE:HG13	2.16	0.45
2:F:71:LEU:HD11	2:F:112:VAL:HG21	1.98	0.45
2:B:103:GLU:O	2:B:107:VAL:HG23	2.17	0.45
4:D:2:GLU:HA	4:D:2:GLU:OE1	2.16	0.45
4:D:43:LEU:HD12	4:D:116:VAL:CG2	2.47	0.45
2:B:71:LEU:HD11	2:B:112:VAL:HG21	1.98	0.45
2:B:75:VAL:O	2:B:78:LEU:HG	2.17	0.45
2:F:13:ARG:HG3	3:G:24:GLN:HE22	1.82	0.45
2:F:50:VAL:HG11	2:F:66:ARG:HH12	1.82	0.45
3:G:59:ILE:HA	3:G:64:PHE:CD2	2.49	0.45
3:C:16:MET:CE	3:C:17:TRP:HA	2.46	0.45
2:B:122:GLN:OE1	2:B:122:GLN:C	2.56	0.44
1:E:82:ALA:HA	4:H:62:GLY:O	2.17	0.44
2:F:122:GLN:C	2:F:122:GLN:OE1	2.56	0.44
2:F:75:VAL:O	2:F:78:LEU:HG	2.17	0.44
2:B:50:VAL:HG11	2:B:66:ARG:HH12	1.82	0.44
2:F:71:LEU:CD1	2:F:112:VAL:HG21	2.48	0.44
3:G:136:CYS:O	3:G:140:ILE:HG13	2.17	0.44
2:B:8:ARG:HD3	2:B:78:LEU:O	2.17	0.44
4:D:99:HIS:CE1	5:D:201:HEM:NA	2.86	0.44
3:C:136:CYS:O	3:C:140:ILE:HG13	2.17	0.44
3:C:21:TRP:CZ3	3:C:72:VAL:HG11	2.52	0.44
2:F:8:ARG:HD3	2:F:78:LEU:O	2.17	0.44
2:B:13:ARG:HG3	3:C:24:GLN:HE22	1.82	0.44
4:H:54:VAL:HG23	4:H:64:PHE:HD1	1.83	0.44
3:C:76:ASP:C	3:C:76:ASP:OD2	2.56	0.44
4:D:31:THR:O	4:D:35:GLU:HG2	2.18	0.44
4:D:54:VAL:HG23	4:D:64:PHE:HD1	1.83	0.44
1:A:142:ASN:C	1:A:144:MET:H	2.21	0.44
2:B:71:LEU:CD1	2:B:112:VAL:HG21	2.48	0.44
2:B:102:ALA:CB	2:B:142:GLN:OE1	2.66	0.44
1:E:59:PHE:O	1:E:63:ILE:HG13	2.18	0.44
3:G:21:TRP:CZ3	3:G:72:VAL:HG11	2.52	0.44
1:E:142:ASN:C	1:E:144:MET:H	2.21	0.44
2:F:130:TRP:O	2:F:134:LEU:HB2	2.18	0.43
4:H:99:HIS:CE1	5:H:201:HEM:NA	2.86	0.43
2:B:130:TRP:O	2:B:134:LEU:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:10:LYS:HE3	4:D:123:SER:O	2.19	0.43
2:F:102:ALA:CB	2:F:142:GLN:OE1	2.66	0.43
4:H:31:THR:O	4:H:35:GLU:HG2	2.18	0.43
3:C:65:LYS:C	3:C:67:HIS:N	2.71	0.43
1:A:82:ALA:HA	4:D:62:GLY:O	2.17	0.43
1:E:76:MET:O	1:E:78:ASP:N	2.51	0.43
2:F:10:LYS:HE3	4:H:123:SER:O	2.19	0.43
4:H:14:LYS:HE3	4:H:14:LYS:HB2	1.87	0.43
4:H:43:LEU:O	4:H:44:LYS:HD3	2.19	0.43
3:G:76:ASP:C	3:G:76:ASP:OD2	2.56	0.43
1:A:59:PHE:O	1:A:63:ILE:HG13	2.18	0.43
1:A:76:MET:O	1:A:78:ASP:N	2.51	0.43
2:B:13:ARG:HH21	2:B:13:ARG:HG2	1.84	0.43
4:D:102:ARG:NE	5:D:201:HEM:HAD2	2.34	0.43
3:G:44:VAL:O	3:G:47:ALA:HB3	2.19	0.43
3:C:44:VAL:O	3:C:47:ALA:HB3	2.18	0.43
4:D:15:GLN:HB3	4:D:125:ILE:HG21	2.01	0.43
3:G:8:GLU:N	3:G:8:GLU:OE1	2.52	0.43
2:B:55:ILE:CG2	2:B:60:PHE:CE2	2.99	0.42
4:D:43:LEU:O	4:D:44:LYS:HD3	2.19	0.42
4:H:102:ARG:NE	5:H:201:HEM:HAD2	2.34	0.42
1:E:53:ASP:OD1	1:E:55:ARG:HD3	2.18	0.42
1:E:67:ILE:HA	1:E:70:LEU:HB2	2.02	0.42
2:F:89:ALA:C	2:F:91:LEU:N	2.72	0.42
1:A:67:ILE:HA	1:A:70:LEU:HB2	2.02	0.42
2:B:139:SER:HA	2:B:142:GLN:HE21	1.85	0.42
2:B:76:ALA:HB1	3:C:29:ARG:HH11	1.85	0.42
3:G:121:MET:O	3:G:123:GLN:N	2.53	0.42
3:C:8:GLU:N	3:C:8:GLU:OE1	2.52	0.42
2:F:139:SER:HA	2:F:142:GLN:HE21	1.85	0.42
2:F:13:ARG:HH21	2:F:13:ARG:HG2	1.84	0.42
2:B:79:ASP:CG	3:C:26:THR:HG21	2.40	0.42
3:C:28:ARG:HG2	3:C:28:ARG:HH21	1.85	0.42
2:F:79:ASP:CG	3:G:26:THR:HG21	2.40	0.42
2:F:76:ALA:HB1	3:G:29:ARG:HH11	1.85	0.42
4:H:15:GLN:HB3	4:H:125:ILE:HG21	2.01	0.42
1:E:112:PHE:CZ	1:E:129:GLU:HA	2.55	0.42
2:F:109:GLU:O	2:F:113:MET:HE2	2.19	0.42
2:F:55:ILE:CG2	2:F:60:PHE:CE2	2.99	0.42
4:H:99:HIS:CD2	4:H:110:PHE:HZ	2.38	0.42
1:A:112:PHE:CZ	1:A:129:GLU:HA	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:106:LYS:O	4:H:109:TYR:HB2	2.20	0.42
4:H:34:ASN:HA	4:H:34:ASN:HD22	1.67	0.42
4:D:61:SER:C	4:D:63:ALA:H	2.24	0.42
4:H:146:LYS:HE3	4:H:146:LYS:HB2	1.93	0.42
4:H:59:MET:HA	4:H:64:PHE:CD2	2.54	0.42
4:H:61:SER:C	4:H:63:ALA:H	2.23	0.42
2:F:37:PHE:N	2:F:37:PHE:CD1	2.88	0.42
2:F:87:GLN:HE22	3:G:70:ARG:CD	2.33	0.42
1:A:53:ASP:OD1	1:A:55:ARG:HD3	2.18	0.41
3:C:121:MET:O	3:C:123:GLN:N	2.53	0.41
4:D:106:LYS:O	4:D:109:TYR:HB2	2.20	0.41
4:D:118:MET:HE1	4:D:137:LEU:HD12	2.00	0.41
1:E:39:ALA:O	1:E:40:PRO:C	2.58	0.41
2:F:104:GLN:C	2:F:106:ASN:H	2.23	0.41
2:F:77:LEU:C	2:F:79:ASP:N	2.72	0.41
3:G:78:ALA:HB1	3:G:140:ILE:HD13	2.02	0.41
3:G:2:SER:OG	3:G:4:SER:OG	2.30	0.41
2:B:89:ALA:C	2:B:91:LEU:N	2.72	0.41
4:D:14:LYS:HB2	4:D:14:LYS:HE3	1.87	0.41
3:G:96:ALA:O	3:G:100:GLN:HG3	2.20	0.41
3:C:67:HIS:O	3:C:70:ARG:N	2.53	0.41
3:G:28:ARG:HH21	3:G:28:ARG:HG2	1.85	0.41
4:H:73:SER:O	4:H:76:ASP:HB3	2.20	0.41
1:A:39:ALA:O	1:A:40:PRO:C	2.58	0.41
2:B:75:VAL:HA	2:B:78:LEU:HG	2.02	0.41
2:B:9:LEU:HA	2:B:9:LEU:HD23	1.83	0.41
3:C:78:ALA:HB1	3:C:140:ILE:HD13	2.02	0.41
1:E:102:LEU:O	1:E:105:VAL:HB	2.20	0.41
2:F:116:VAL:O	2:F:120:ILE:HG12	2.21	0.41
3:G:65:LYS:C	3:G:67:HIS:N	2.72	0.41
3:G:67:HIS:O	3:G:70:ARG:N	2.53	0.41
1:A:1:ASP:CG	1:A:2:CYS:N	2.73	0.41
2:B:74:CYS:HB3	2:B:137:ILE:HD13	2.03	0.41
4:D:99:HIS:CD2	4:D:110:PHE:HZ	2.38	0.41
2:F:9:LEU:HD23	2:F:9:LEU:HA	1.83	0.41
1:A:114:THR:HA	1:E:121:VAL:HG13	2.03	0.41
2:F:75:VAL:HA	2:F:78:LEU:HG	2.02	0.41
2:F:79:ASP:OD1	3:G:26:THR:HG21	2.21	0.41
4:D:59:MET:HA	4:D:64:PHE:CD2	2.54	0.41
2:F:74:CYS:HB3	2:F:137:ILE:HD13	2.03	0.41
3:G:50:LEU:HD22	3:G:103:ALA:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:142:ASN:O	4:H:146:LYS:HE3	2.20	0.41
2:B:104:GLN:C	2:B:106:ASN:H	2.23	0.41
3:C:50:LEU:HD22	3:C:103:ALA:O	2.20	0.41
4:D:73:SER:O	4:D:76:ASP:HB3	2.20	0.41
2:B:109:GLU:O	2:B:113:MET:HE2	2.21	0.41
2:B:87:GLN:HE22	3:C:70:ARG:CD	2.33	0.41
3:C:45:PRO:O	3:C:48:VAL:HG23	2.21	0.41
4:D:142:ASN:O	4:D:146:LYS:HE3	2.20	0.41
1:A:126:PRO:HD2	1:E:38:TYR:CE1	2.56	0.41
2:B:116:VAL:O	2:B:120:ILE:HG12	2.21	0.41
2:B:77:LEU:C	2:B:79:ASP:N	2.72	0.41
3:C:92:LEU:HD23	3:C:92:LEU:HA	1.92	0.41
3:G:113:ILE:HA	3:G:116:SER:CB	2.51	0.41
1:A:102:LEU:O	1:A:105:VAL:HB	2.20	0.40
1:A:125:LEU:HD23	1:E:112:PHE:HB2	2.02	0.40
2:B:37:PHE:N	2:B:37:PHE:CD1	2.88	0.40
4:D:61:SER:O	4:D:62:GLY:C	2.60	0.40
1:E:1:ASP:CG	1:E:2:CYS:N	2.73	0.40
3:G:37:PHE:CE2	3:G:68:CYS:HA	2.56	0.40
1:E:89:HIS:CD2	4:H:70:ARG:NE	2.89	0.40
2:B:73:MET:HB3	2:B:73:MET:HE2	1.87	0.40
2:B:79:ASP:OD1	3:C:26:THR:HG21	2.21	0.40
3:C:113:ILE:HA	3:C:116:SER:CB	2.51	0.40
2:F:2:GLU:OE2	2:F:2:GLU:N	2.45	0.40
2:B:142:GLN:HE21	2:B:142:GLN:HB2	1.71	0.40
3:C:96:ALA:O	3:C:100:GLN:HG3	2.20	0.40
4:D:8:ASP:O	4:D:9:ALA:C	2.59	0.40
2:F:21:SER:OG	3:G:18:ALA:HB2	2.22	0.40
4:H:69:VAL:O	4:H:73:SER:N	2.54	0.40
1:A:109:LYS:HG2	1:E:126:PRO:HG3	2.04	0.40
2:F:114:MET:CE	2:F:114:MET:HA	2.52	0.40
3:G:14:GLN:HE21	3:G:14:GLN:HB2	1.64	0.40
3:C:149:ALA:O	3:C:150:GLU:O	2.39	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	144/146 (99%)	126 (88%)	17 (12%)	1 (1%)	25	60
1	E	144/146 (99%)	126 (88%)	17 (12%)	1 (1%)	25	60
2	B	142/144 (99%)	113 (80%)	29 (20%)	0	100	100
2	F	142/144 (99%)	113 (80%)	29 (20%)	0	100	100
3	C	148/150 (99%)	125 (84%)	22 (15%)	1 (1%)	25	60
3	G	148/150 (99%)	125 (84%)	22 (15%)	1 (1%)	25	60
4	D	147/149 (99%)	130 (88%)	14 (10%)	3 (2%)	9	31
4	H	147/149 (99%)	130 (88%)	14 (10%)	3 (2%)	9	31
All	All	1162/1178 (99%)	988 (85%)	164 (14%)	10 (1%)	20	54

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	77	PHE
1	A	77	PHE
4	D	26	ALA
4	D	101	ALA
4	H	26	ALA
4	H	101	ALA
3	C	122	PRO
3	G	122	PRO
4	D	62	GLY
4	H	62	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	120/120 (100%)	115 (96%)	5 (4%)	34	69
1	E	120/120 (100%)	115 (96%)	5 (4%)	34	69
2	B	116/116 (100%)	110 (95%)	6 (5%)	27	61
2	F	116/116 (100%)	110 (95%)	6 (5%)	27	61
3	C	124/124 (100%)	114 (92%)	10 (8%)	14	38
3	G	124/124 (100%)	114 (92%)	10 (8%)	14	38
4	D	122/122 (100%)	116 (95%)	6 (5%)	29	63
4	H	122/122 (100%)	116 (95%)	6 (5%)	29	63
All	All	964/964 (100%)	910 (94%)	54 (6%)	25	57

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ILE
1	A	20	PHE
1	A	34	SER
1	A	70	LEU
1	A	87	LEU
2	B	71	LEU
2	B	73	MET
2	B	97	SER
2	B	108	VAL
2	B	119	GLU
2	B	123	ASN
3	C	8	GLU
3	C	14	GLN
3	C	15	LEU
3	C	16	MET
3	C	24	GLN
3	C	26	THR
3	C	57	THR
3	C	60	ASP
3	C	63	GLU
3	C	126	SER
4	D	1	SER
4	D	70	ARG
4	D	102	ARG

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Mol	Chain	Res	Type
4	D	111	ASP
4	D	137	LEU
4	D	138	LEU
1	E	4	ILE
1	E	20	PHE
1	E	34	SER
1	E	70	LEU
1	E	87	LEU
2	F	71	LEU
2	F	73	MET
2	F	97	SER
2	F	108	VAL
2	F	119	GLU
2	F	123	ASN
3	G	8	GLU
3	G	14	GLN
3	G	15	LEU
3	G	16	MET
3	G	24	GLN
3	G	26	THR
3	G	57	THR
3	G	60	ASP
3	G	63	GLU
3	G	126	SER
4	H	1	SER
4	H	70	ARG
4	H	102	ARG
4	H	111	ASP
4	H	137	LEU
4	H	138	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	23	ASN
2	B	30	HIS
2	B	106	ASN
3	C	3	ASN
3	C	14	GLN
3	C	24	GLN
3	C	34	GLN
3	C	100	GLN

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Mol	Chain	Res	Type
3	C	115	GLN
4	D	18	GLN
2	F	23	ASN
2	F	30	HIS
2	F	106	ASN
3	G	3	ASN
3	G	14	GLN
3	G	24	GLN
3	G	34	GLN
3	G	100	GLN
3	G	115	GLN
4	H	18	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

12 carbohydrates 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	$\# Z > 2$
7	NAG	D	202	4,7	14,14,15	0.94	1 (7%)	15,19,21	1.14	1 (6%)
7	NAG	D	203	7	14,14,15	0.66	0	15,19,21	1.33	2 (13%)
7	MAN	D	204	7	11,11,12	1.12	1 (9%)	13,15,17	1.48	2 (15%)
7	MAN	D	205	7	11,11,12	0.94	0	13,15,17	0.87	1 (7%)
7	MAN	D	206	7	11,11,12	0.82	1 (9%)	13,15,17	0.82	1 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	FUC	D	207	7	9,10,11	0.66	0	13,14,16	0.90	0
7	NAG	H	202	4,7	14,14,15	0.93	1 (7%)	15,19,21	1.14	1 (6%)
7	NAG	H	203	7	14,14,15	0.66	0	15,19,21	1.32	2 (13%)
7	MAN	H	204	7	11,11,12	1.12	1 (9%)	13,15,17	1.48	2 (15%)
7	MAN	H	205	7	11,11,12	0.95	0	13,15,17	0.86	1 (7%)
7	MAN	H	206	7	11,11,12	0.81	0	13,15,17	0.81	1 (7%)
7	FUC	H	207	7	9,10,11	0.66	0	13,14,16	0.90	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
7	NAG	D	202	4,7	-	0/6/23/26	0/1/1/1
7	NAG	D	203	7	-	0/6/23/26	0/1/1/1
7	MAN	D	204	7	-	0/2/19/22	0/1/1/1
7	MAN	D	205	7	-	0/2/19/22	0/1/1/1
7	MAN	D	206	7	-	0/2/19/22	0/1/1/1
7	FUC	D	207	7	-	0/0/17/20	0/1/1/1
7	NAG	H	202	4,7	-	0/6/23/26	0/1/1/1
7	NAG	H	203	7	-	0/6/23/26	0/1/1/1
7	MAN	H	204	7	-	0/2/19/22	0/1/1/1
7	MAN	H	205	7	-	0/2/19/22	0/1/1/1
7	MAN	H	206	7	-	0/2/19/22	0/1/1/1
7	FUC	H	207	7	-	0/0/17/20	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	202	NAG	O5-C1	-2.08	1.40	1.43
7	H	202	NAG	O5-C1	-2.04	1.40	1.43
7	D	206	MAN	C2-C3	2.02	1.55	1.52
7	D	204	MAN	C2-C3	3.14	1.56	1.52
7	H	204	MAN	C2-C3	3.14	1.56	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	203	NAG	C2-N2-C7	-3.18	118.30	122.94
7	H	203	NAG	C2-N2-C7	-3.14	118.36	122.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	H	202	NAG	O4-C4-C3	-2.97	103.90	110.36
7	D	202	NAG	O4-C4-C3	-2.96	103.93	110.36
7	D	203	NAG	C4-C3-C2	-2.93	106.72	111.02
7	H	203	NAG	C4-C3-C2	-2.92	106.74	111.02
7	H	206	MAN	C1-C2-C3	2.11	112.33	109.65
7	D	206	MAN	C1-C2-C3	2.14	112.36	109.65
7	H	205	MAN	C1-O5-C5	2.47	115.56	112.17
7	D	205	MAN	C1-O5-C5	2.47	115.58	112.17
7	D	204	MAN	C2-C3-C4	2.76	115.70	110.88
7	H	204	MAN	C2-C3-C4	2.77	115.70	110.88
7	D	204	MAN	C1-C2-C3	3.72	114.36	109.65
7	H	204	MAN	C1-C2-C3	3.73	114.38	109.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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$
5	HEM	A	200	1	28,50,50	1.82	6 (21%)	17,82,82	2.46	9 (52%)
5	HEM	B	201	2	28,50,50	1.86	4 (14%)	17,82,82	2.28	8 (47%)
5	HEM	C	200	3	28,50,50	1.87	6 (21%)	17,82,82	2.42	8 (47%)
5	HEM	D	201	4	28,50,50	1.80	6 (21%)	17,82,82	2.43	9 (52%)
5	HEM	E	200	1	28,50,50	1.81	6 (21%)	17,82,82	2.46	9 (52%)
5	HEM	F	201	2	28,50,50	1.86	4 (14%)	17,82,82	2.27	8 (47%)
5	HEM	G	200	3	28,50,50	1.87	6 (21%)	17,82,82	2.42	9 (52%)
5	HEM	H	201	4	28,50,50	1.80	6 (21%)	17,82,82	2.43	9 (52%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	HEM	A	200	1	-	0/6/54/54	0/0/8/8
5	HEM	B	201	2	-	0/6/54/54	0/0/8/8
5	HEM	C	200	3	-	0/6/54/54	0/0/8/8
5	HEM	D	201	4	-	0/6/54/54	0/0/8/8
5	HEM	E	200	1	-	0/6/54/54	0/0/8/8
5	HEM	F	201	2	-	0/6/54/54	0/0/8/8
5	HEM	G	200	3	-	0/6/54/54	0/0/8/8
5	HEM	H	201	4	-	0/6/54/54	0/0/8/8

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	201	HEM	C3C-CAC	-6.35	1.35	1.47
5	F	201	HEM	C3C-CAC	-6.34	1.35	1.47
5	E	200	HEM	C3C-CAC	-5.99	1.35	1.47
5	A	200	HEM	C3C-CAC	-5.99	1.35	1.47
5	G	200	HEM	C3C-CAC	-5.92	1.36	1.47
5	C	200	HEM	C3C-CAC	-5.91	1.36	1.47
5	H	201	HEM	C3C-CAC	-5.90	1.36	1.47
5	D	201	HEM	C3C-CAC	-5.89	1.36	1.47
5	A	200	HEM	C4A-CHB	-2.42	1.33	1.40
5	E	200	HEM	C4A-CHB	-2.39	1.33	1.40
5	G	200	HEM	C4A-CHB	-2.30	1.34	1.40
5	C	200	HEM	C4A-CHB	-2.29	1.34	1.40
5	D	201	HEM	C4A-CHB	-2.29	1.34	1.40
5	H	201	HEM	C4A-CHB	-2.29	1.34	1.40
5	A	200	HEM	C1B-NB	-2.12	1.34	1.36
5	E	200	HEM	C1B-NB	-2.08	1.34	1.36
5	D	201	HEM	C4D-ND	2.02	1.39	1.36
5	H	201	HEM	C4D-ND	2.02	1.39	1.36
5	A	200	HEM	C1C-NC	2.06	1.39	1.36
5	E	200	HEM	C1C-NC	2.08	1.39	1.36
5	G	200	HEM	C1C-NC	2.31	1.39	1.36
5	C	200	HEM	C1C-NC	2.33	1.39	1.36
5	B	201	HEM	C1A-NA	2.38	1.41	1.36
5	F	201	HEM	C1A-NA	2.40	1.41	1.36
5	D	201	HEM	C1A-NA	2.76	1.42	1.36
5	H	201	HEM	C1A-NA	2.77	1.42	1.36
5	G	200	HEM	C1A-NA	2.87	1.42	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	200	HEM	C1A-NA	2.88	1.42	1.36
5	H	201	HEM	C4B-NB	2.93	1.42	1.36
5	D	201	HEM	C4B-NB	2.95	1.42	1.36
5	E	200	HEM	C4B-NB	3.07	1.42	1.36
5	A	200	HEM	C4B-NB	3.08	1.42	1.36
5	B	201	HEM	C3B-C2B	3.22	1.44	1.40
5	F	201	HEM	C3B-C2B	3.23	1.44	1.40
5	B	201	HEM	C4B-NB	3.24	1.43	1.36
5	F	201	HEM	C4B-NB	3.26	1.43	1.36
5	D	201	HEM	C3B-C2B	3.36	1.44	1.40
5	H	201	HEM	C3B-C2B	3.37	1.44	1.40
5	C	200	HEM	C4B-NB	3.38	1.43	1.36
5	G	200	HEM	C4B-NB	3.41	1.43	1.36
5	E	200	HEM	C3B-C2B	3.59	1.45	1.40
5	A	200	HEM	C3B-C2B	3.62	1.45	1.40
5	C	200	HEM	C3B-C2B	3.69	1.45	1.40
5	G	200	HEM	C3B-C2B	3.71	1.45	1.40

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	201	HEM	C4C-C3C-C2C	-5.02	103.39	106.90
5	D	201	HEM	C4C-C3C-C2C	-4.98	103.42	106.90
5	C	200	HEM	C4C-C3C-C2C	-4.93	103.45	106.90
5	G	200	HEM	C4C-C3C-C2C	-4.93	103.45	106.90
5	A	200	HEM	C4C-C3C-C2C	-4.77	103.57	106.90
5	E	200	HEM	C4C-C3C-C2C	-4.74	103.59	106.90
5	B	201	HEM	C4C-C3C-C2C	-4.42	103.81	106.90
5	F	201	HEM	C4C-C3C-C2C	-4.41	103.82	106.90
5	E	200	HEM	C1D-C2D-C3D	-3.96	104.24	107.00
5	A	200	HEM	C1D-C2D-C3D	-3.92	104.27	107.00
5	B	201	HEM	C1D-C2D-C3D	-3.66	104.45	107.00
5	F	201	HEM	C1D-C2D-C3D	-3.66	104.45	107.00
5	E	200	HEM	CMA-C3A-C4A	-3.58	122.97	128.46
5	A	200	HEM	CMA-C3A-C4A	-3.57	122.98	128.46
5	C	200	HEM	CMA-C3A-C4A	-3.32	123.36	128.46
5	G	200	HEM	CMA-C3A-C4A	-3.32	123.36	128.46
5	F	201	HEM	CMA-C3A-C4A	-3.27	123.44	128.46
5	B	201	HEM	CMA-C3A-C4A	-3.25	123.47	128.46
5	G	200	HEM	C1D-C2D-C3D	-3.24	104.74	107.00
5	C	200	HEM	C1D-C2D-C3D	-3.23	104.75	107.00
5	H	201	HEM	C1D-C2D-C3D	-2.94	104.95	107.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	201	HEM	C1D-C2D-C3D	-2.92	104.97	107.00
5	C	200	HEM	CMD-C2D-C1D	-2.81	124.15	128.46
5	G	200	HEM	CMD-C2D-C1D	-2.79	124.18	128.46
5	H	201	HEM	CMD-C2D-C1D	-2.74	124.25	128.46
5	D	201	HEM	CMA-C3A-C4A	-2.74	124.26	128.46
5	D	201	HEM	CMD-C2D-C1D	-2.72	124.28	128.46
5	H	201	HEM	CMA-C3A-C4A	-2.72	124.29	128.46
5	A	200	HEM	CMD-C2D-C1D	-2.32	124.90	128.46
5	E	200	HEM	CMD-C2D-C1D	-2.31	124.92	128.46
5	B	201	HEM	CMD-C2D-C1D	-2.18	125.11	128.46
5	F	201	HEM	CMD-C2D-C1D	-2.15	125.15	128.46
5	G	200	HEM	C3B-C4B-NB	2.02	111.82	109.21
5	D	201	HEM	CMB-C2B-C3B	2.14	128.86	124.89
5	H	201	HEM	CMB-C2B-C3B	2.15	128.89	124.89
5	F	201	HEM	C3B-C4B-NB	2.18	112.03	109.21
5	B	201	HEM	C3B-C4B-NB	2.20	112.05	109.21
5	E	200	HEM	CMB-C2B-C3B	2.23	129.03	124.89
5	A	200	HEM	CMB-C2B-C3B	2.25	129.06	124.89
5	G	200	HEM	CMB-C2B-C3B	2.25	129.07	124.89
5	C	200	HEM	CMB-C2B-C3B	2.27	129.11	124.89
5	E	200	HEM	C4A-C3A-C2A	2.29	108.59	107.00
5	A	200	HEM	C4A-C3A-C2A	2.33	108.62	107.00
5	A	200	HEM	C3B-C4B-NB	2.42	112.34	109.21
5	E	200	HEM	C3B-C4B-NB	2.43	112.35	109.21
5	D	201	HEM	C3B-C4B-NB	2.50	112.44	109.21
5	B	201	HEM	C4A-C3A-C2A	2.50	108.73	107.00
5	H	201	HEM	C3B-C4B-NB	2.52	112.47	109.21
5	F	201	HEM	C4A-C3A-C2A	2.55	108.77	107.00
5	C	200	HEM	CBA-CAA-C2A	2.63	117.51	112.48
5	G	200	HEM	CBA-CAA-C2A	2.65	117.55	112.48
5	F	201	HEM	CMD-C2D-C3D	2.89	130.40	124.94
5	B	201	HEM	CMD-C2D-C3D	2.91	130.44	124.94
5	H	201	HEM	C4A-C3A-C2A	2.95	109.05	107.00
5	D	201	HEM	C4A-C3A-C2A	3.04	109.11	107.00
5	D	201	HEM	CMD-C2D-C3D	3.08	130.75	124.94
5	H	201	HEM	CMD-C2D-C3D	3.10	130.79	124.94
5	A	200	HEM	CMD-C2D-C3D	3.12	130.82	124.94
5	E	200	HEM	CMD-C2D-C3D	3.13	130.84	124.94
5	C	200	HEM	C4A-C3A-C2A	3.13	109.17	107.00
5	G	200	HEM	C4A-C3A-C2A	3.14	109.18	107.00
5	G	200	HEM	CMD-C2D-C3D	3.26	131.08	124.94
5	C	200	HEM	CMD-C2D-C3D	3.27	131.10	124.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	201	HEM	CBA-CAA-C2A	3.44	119.06	112.48
5	B	201	HEM	CBA-CAA-C2A	3.46	119.10	112.48
5	D	201	HEM	CBA-CAA-C2A	3.56	119.30	112.48
5	H	201	HEM	CBA-CAA-C2A	3.57	119.31	112.48
5	E	200	HEM	CBA-CAA-C2A	3.83	119.81	112.48
5	A	200	HEM	CBA-CAA-C2A	3.85	119.85	112.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	200	HEM	1	0
5	B	201	HEM	1	0
5	C	200	HEM	1	0
5	D	201	HEM	3	0
5	E	200	HEM	1	0
5	F	201	HEM	1	0
5	G	200	HEM	1	0
5	H	201	HEM	3	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	146/146 (100%)	-0.24	0 100 100	34, 55, 83, 136	0
1	E	146/146 (100%)	-0.34	0 100 100	34, 55, 83, 136	0
2	B	144/144 (100%)	0.06	4 (2%) 53 48	40, 71, 107, 164	0
2	F	144/144 (100%)	-0.08	1 (0%) 87 86	40, 71, 107, 164	0
3	C	150/150 (100%)	0.12	4 (2%) 55 50	39, 69, 113, 170	0
3	G	150/150 (100%)	0.01	3 (2%) 65 62	39, 69, 113, 170	0
4	D	149/149 (100%)	-0.09	2 (1%) 77 76	40, 66, 104, 140	0
4	H	149/149 (100%)	-0.17	2 (1%) 77 76	40, 66, 104, 140	0
All	All	1178/1178 (100%)	-0.09	16 (1%) 75 74	34, 65, 107, 170	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	1	SER	9.1
3	G	1	SER	9.0
4	D	1	SER	4.6
4	H	1	SER	4.1
2	B	144	ASN	3.2
3	G	107	LYS	3.1
3	C	2	SER	3.0
3	C	26	THR	2.6
4	D	105	LEU	2.6
2	B	137	ILE	2.5
2	B	131	GLN	2.4
3	C	19	ASN	2.3
2	B	47	PHE	2.2
3	G	2	SER	2.2
2	F	21	SER	2.2
4	H	30	TRP	2.1

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](#)

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
7	FUC	D	207	10/11	0.89	0.28	-	106,106,106,106	0
7	MAN	D	206	11/12	0.54	0.81	-	218,220,221,222	0
7	MAN	H	206	11/12	0.38	0.60	-	218,220,221,222	0
7	MAN	D	205	11/12	0.33	0.88	-	236,237,238,239	0
7	NAG	D	203	14/15	0.84	0.38	-	153,153,153,153	0
7	MAN	D	204	11/12	0.74	0.45	-	230,233,236,237	0
7	MAN	H	205	11/12	0.03	1.13	-	236,237,238,239	0
7	NAG	H	202	14/15	0.73	0.40	-	156,156,156,156	0
7	FUC	H	207	10/11	0.83	0.29	-	106,106,106,106	0
7	NAG	D	202	14/15	0.62	0.30	-	156,156,156,156	0
7	NAG	H	203	14/15	0.72	0.43	-	153,153,153,153	0
7	MAN	H	204	11/12	0.61	0.66	-	230,233,236,237	0

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	HEM	G	200	43/43	0.93	0.24	0.43	67,67,67,67	0
5	HEM	H	201	43/43	0.93	0.23	0.32	67,67,67,67	0
5	HEM	C	200	43/43	0.92	0.23	0.23	67,67,67,67	0
5	HEM	A	200	43/43	0.95	0.21	0.18	67,67,67,67	0
5	HEM	F	201	43/43	0.95	0.22	0.07	67,67,67,67	0
5	HEM	D	201	43/43	0.94	0.22	0.02	67,67,67,67	0
5	HEM	E	200	43/43	0.94	0.19	0.02	67,67,67,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	HEM	B	201	43/43	0.93	0.23	-0.28	67,67,67,67	0
6	CA	B	202	1/1	0.80	0.27	-	67,67,67,67	0
6	CA	F	202	1/1	0.95	0.04	-	67,67,67,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.