



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2025 – 03:30 PM EDT

PDB ID : 1JJU  
Title : Structure of a Quinohemoprotein Amine Dehydrogenase with a Unique Redox Cofactor and Highly Unusual Crosslinking  
Authors : Datta, S.; Mori, Y.; Takagi, K.; Kawaguchi, K.; Chen, Z.-W.; Kano, K.; Ikeda, T.; Okajima, T.; Kuroda, S.; Tanizawa, K.; Mathews, F.S.  
Deposited on : 2001-07-09  
Resolution : 2.05 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

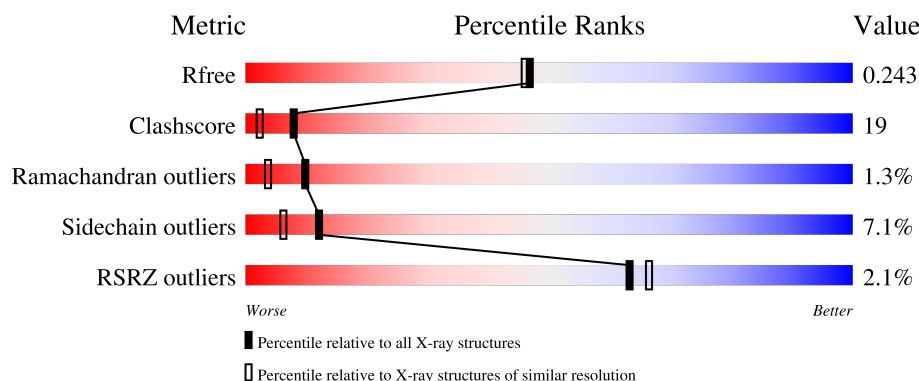
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*




The reported resolution of this entry is 2.05 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))
$R_{free}$	164625	2096 (2.04-2.04)
Clashscore	180529	2229 (2.04-2.04)
Ramachandran outliers	177936	2217 (2.04-2.04)
Sidechain outliers	177891	2217 (2.04-2.04)
RSRZ outliers	164620	2096 (2.04-2.04)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	489	
2	B	337	
3	C	79	

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7698 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 QUINOHEMOPROTEIN AMINE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	489	Total	C	N	O	S	0	2	0
			3718	2308	667	731	12			

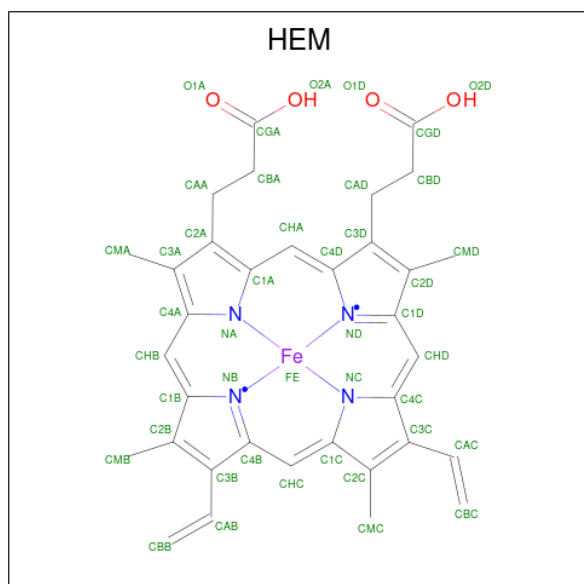
- Molecule 2 is a protein called QUINOHEMOPROTEIN AMINE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	337	Total	C	N	O	S	0	2	0
			2621	1656	443	511	11			

- Molecule 3 is a protein called QUINOHEMOPROTEIN AMINE DEHYDROGENASE.

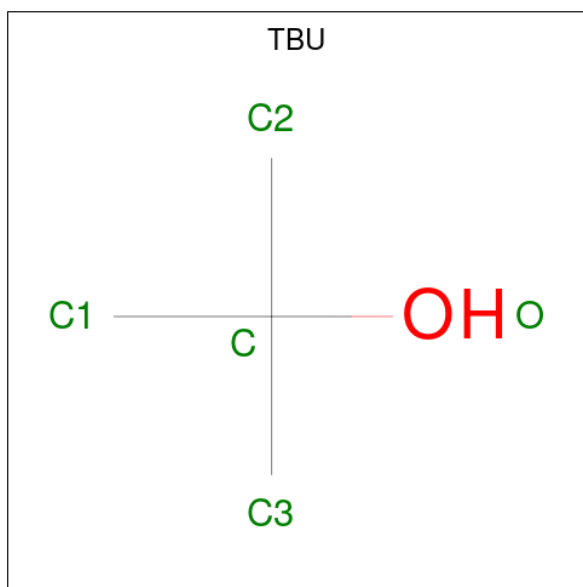
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	79	Total	C	N	O	S	0	1	0
			622	391	98	126	7			

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



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

- Molecule 5 is TERTIARY-BUTYL ALCOHOL (three-letter code: TBU) (formula: C<sub>4</sub>H<sub>10</sub>O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O		
			5	4	1		
5	B	1	Total	C	O		
			5	4	1		
5	C	1	Total	C	O		
			5	4	1		

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total	Na		
			1	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	330	Total	O		
			330	330		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	246	Total 246	O 246	0	0
7	C	59	Total 59	O 59	0	0

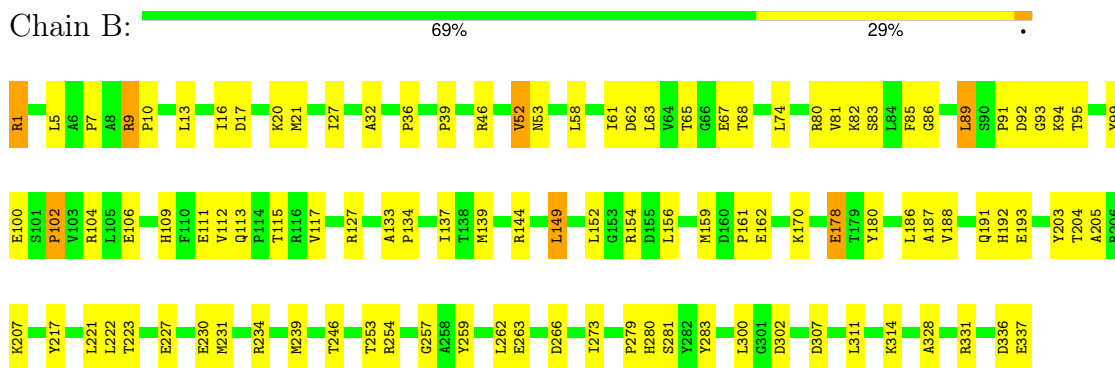
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: QUINOHEMOPROTEIN AMINE DEHYDROGENASE

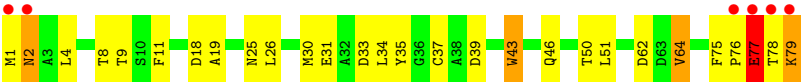


#### • Molecule 2: QUINOHEMOPROTEIN AMINE DEHYDROGENASE



#### • Molecule 3: QUINOHEMOPROTEIN AMINE DEHYDROGENASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	99.49Å 99.49Å 214.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.05 30.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	86.4 (30.00-2.05) 86.2 (30.00-2.05)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.53 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.202 , 0.251 0.194 , 0.243	Depositor DCC
$R_{free}$ test set	5994 reflections (10.17%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.4	Xtriage
Anisotropy	0.516	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 57.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7698	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TBU, NA, TRQ, 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	0.36	3/3803 (0.1%)	0.74	7/5174 (0.1%)
2	B	0.34	0/2675	0.66	0/3638
3	C	0.35	0/623	0.64	0/851
All	All	0.35	3/7101 (0.0%)	0.70	7/9663 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	HIS	CE1-NE2	5.61	1.45	1.32
1	A	15	HIS	CE1-NE2	5.56	1.45	1.32
1	A	104	HIS	CE1-NE2	5.12	1.44	1.32

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	467	GLY	N-CA-C	-9.82	88.54	113.10
1	A	393	GLN	N-CA-C	9.32	136.15	111.00
1	A	468	GLU	N-CA-C	8.90	135.04	111.00
1	A	15	HIS	ND1-CG-CD2	8.23	120.32	108.80
1	A	126	HIS	ND1-CG-CD2	8.13	120.19	108.80
1	A	104	HIS	ND1-CG-CD2	7.90	119.86	108.80
1	A	392	GLY	N-CA-C	-6.18	97.64	113.10

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3718	0	3584	165	0
2	B	2621	0	2609	89	0
3	C	622	0	550	37	0
4	A	86	0	60	1	0
5	A	5	0	10	0	0
5	B	5	0	10	0	0
5	C	5	0	9	0	0
6	C	1	0	0	0	0
7	A	330	0	0	12	0
7	B	246	0	0	12	0
7	C	59	0	0	0	0
All	All	7698	0	6832	269	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 19.

All (269) 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:393:GLN:HE22	1:A:398:ASP:HB3	1.05	1.13
2:B:144:ARG:HG3	2:B:227[A]:GLU:HG2	1.35	1.05
1:A:393:GLN:NE2	1:A:398:ASP:HB3	1.81	0.95
1:A:444:ASN:HD22	1:A:453:ASN:HD21	1.13	0.94
1:A:456:ASN:HD21	1:A:475:HIS:HE1	1.15	0.94
1:A:387:LEU:HG	1:A:402:LEU:HD11	1.47	0.93
1:A:102:ARG:HH12	1:A:136:GLN:HE21	1.13	0.92
1:A:452:ASN:HD21	3:C:30:MET:H	0.94	0.90
3:C:1:MET:HB3	3:C:9:THR:HG22	1.54	0.89
1:A:452:ASN:ND2	3:C:30:MET:H	1.73	0.85
1:A:86:VAL:HG22	1:A:360:PRO:HG3	1.60	0.82
1:A:6:VAL:HG11	1:A:59:ILE:HG12	1.61	0.82
1:A:293:GLN:HG2	7:A:1099:HOH:O	1.79	0.82
1:A:364:ILE:HG23	3:C:4:LEU:HD22	1.62	0.82
2:B:102:PRO:HG2	2:B:113:GLN:HB2	1.60	0.81
1:A:452:ASN:HD21	3:C:30:MET:N	1.78	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:1247:HOH:O	3:C:79:LYS:HD2	1.82	0.78
1:A:444:ASN:ND2	1:A:453:ASN:HD21	1.83	0.76
1:A:229:TRP:HE1	1:A:242:GLN:HE21	1.34	0.75
1:A:456:ASN:ND2	1:A:475:HIS:HE1	1.82	0.75
1:A:393:GLN:HG3	1:A:396:THR:HG23	1.68	0.75
2:B:32:ALA:HB1	2:B:52:VAL:HG21	1.67	0.74
1:A:229:TRP:HE1	1:A:242:GLN:NE2	1.85	0.73
1:A:456:ASN:HD21	1:A:475:HIS:CE1	2.04	0.73
2:B:95:THR:HG22	7:B:996:HOH:O	1.87	0.73
2:B:254:ARG:HG3	2:B:254:ARG:HH11	1.53	0.73
1:A:306:THR:HG23	1:A:339:GLU:HB2	1.69	0.73
1:A:102:ARG:HH12	1:A:136:GLN:NE2	1.87	0.72
1:A:392:GLY:O	1:A:394:PRO:HD2	1.90	0.72
1:A:266:ARG:HD2	7:A:1112:HOH:O	1.89	0.72
2:B:149:LEU:HB2	2:B:161:PRO:HG3	1.72	0.71
1:A:407:ALA:HB2	1:A:463:VAL:HG12	1.71	0.71
1:A:243:ILE:HD11	3:C:79:LYS:HD3	1.71	0.71
1:A:45:ARG:HD2	1:A:46:ASN:ND2	2.05	0.70
2:B:156:LEU:HB3	2:B:170:LYS:HB2	1.74	0.70
1:A:388:ASN:HB3	1:A:392:GLY:O	1.92	0.69
1:A:393:GLN:HE21	1:A:396:THR:HG23	1.58	0.68
1:A:45:ARG:HD2	1:A:46:ASN:HD21	1.58	0.68
1:A:390:PRO:HD2	1:A:398:ASP:OD2	1.93	0.68
1:A:444:ASN:HD22	1:A:453:ASN:ND2	1.90	0.67
1:A:325:LEU:HB3	1:A:327:LEU:CD1	2.25	0.66
2:B:280:HIS:HD2	2:B:281:SER:O	1.77	0.66
3:C:1:MET:CB	3:C:9:THR:HG22	2.26	0.66
2:B:178:GLU:H	2:B:178:GLU:CD	1.98	0.65
1:A:466:GLU:HA	1:A:468:GLU:HG2	1.78	0.65
2:B:32:ALA:HB1	2:B:52:VAL:CG2	2.27	0.65
1:A:393:GLN:NE2	1:A:396:THR:OG1	2.31	0.64
2:B:191:GLN:HA	2:B:193:GLU:OE2	1.96	0.63
1:A:291:GLU:CD	1:A:326:LYS:HE2	2.18	0.63
2:B:100:GLU:OE2	2:B:127:ARG:NH2	2.29	0.63
1:A:393:GLN:HE21	1:A:396:THR:CG2	2.12	0.62
1:A:90:GLU:HB3	3:C:2:ASN:HD22	1.64	0.62
1:A:23:TRP:HB2	1:A:26:ILE:HB	1.81	0.62
1:A:285:ARG:HD3	1:A:402:LEU:O	1.99	0.62
2:B:1:ARG:HG3	2:B:1:ARG:HH11	1.64	0.62
2:B:81:VAL:HG22	2:B:102:PRO:HB3	1.82	0.61
1:A:376:LYS:HB3	1:A:438:PRO:HG2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:PRO:HD2	1:A:294:LEU:HD23	1.83	0.61
2:B:13:LEU:HD12	2:B:61:ILE:CD1	2.30	0.61
2:B:230:GLU:OE2	2:B:230:GLU:HA	2.02	0.60
1:A:56:ARG:HA	1:A:59:ILE:HD11	1.83	0.60
3:C:35:TYR:CE2	3:C:64:VAL:HG22	2.36	0.60
2:B:203:TYR:HB3	2:B:239:MET:HE1	1.84	0.60
1:A:386:TRP:HA	1:A:402:LEU:HD13	1.83	0.60
1:A:388:ASN:O	1:A:393:GLN:HB3	2.02	0.60
1:A:46:ASN:N	1:A:46:ASN:HD22	1.99	0.59
2:B:13:LEU:HD12	2:B:61:ILE:HD11	1.83	0.59
2:B:5:LEU:HD11	2:B:13:LEU:HD22	1.84	0.59
2:B:109:HIS:HE1	2:B:111:GLU:OE1	1.85	0.59
1:A:466:GLU:C	1:A:468:GLU:H	2.06	0.58
1:A:182:ARG:HG3	1:A:188:ASP:OD1	2.03	0.58
1:A:137:ALA:O	1:A:138:LEU:HB2	2.03	0.58
2:B:80:ARG:HD3	2:B:82:LYS:HE2	1.84	0.58
1:A:95:SER:O	1:A:99:THR:HG23	2.04	0.58
1:A:393:GLN:HE21	1:A:396:THR:CB	2.16	0.58
1:A:325:LEU:HB3	1:A:327:LEU:HD11	1.84	0.58
2:B:307:ASP:HA	2:B:314:LYS:HE3	1.84	0.58
1:A:460:ILE:HG13	1:A:460:ILE:O	2.03	0.58
3:C:1:MET:SD	3:C:4:LEU:HD12	2.44	0.57
1:A:388:ASN:O	1:A:393:GLN:CB	2.52	0.57
1:A:392:GLY:O	1:A:394:PRO:CD	2.51	0.57
2:B:32:ALA:CB	2:B:52:VAL:HG21	2.33	0.57
1:A:102:ARG:HA	3:C:9:THR:OG1	2.04	0.57
2:B:254:ARG:HG3	2:B:254:ARG:NH1	2.15	0.57
1:A:286:LEU:HD21	1:A:294:LEU:HD11	1.87	0.57
1:A:408:SER:HB2	1:A:462:THR:HG23	1.86	0.57
1:A:308:PRO:HG2	1:A:311:VAL:HG12	1.87	0.56
3:C:11:PHE:HB3	3:C:19:ALA:HB2	1.86	0.56
2:B:104:ARG:HD3	7:B:1032:HOH:O	2.05	0.56
1:A:396:THR:O	1:A:399:ASP:OD1	2.22	0.56
2:B:61:ILE:HG22	2:B:68:THR:HA	1.87	0.56
2:B:65:THR:HG22	2:B:67:GLU:H	1.70	0.56
1:A:413:ASN:HD22	1:A:419:GLU:HG2	1.71	0.55
1:A:256:HIS:HD2	1:A:257:ASP:O	1.89	0.55
1:A:392:GLY:C	1:A:394:PRO:HD2	2.27	0.55
2:B:134:PRO:HG2	2:B:137:ILE:HD11	1.88	0.55
1:A:267:LEU:C	1:A:267:LEU:HD12	2.27	0.55
1:A:31:LYS:HB3	1:A:112:GLN:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:GLU:CD	1:A:54:GLU:H	2.10	0.55
1:A:233:LEU:N	1:A:233:LEU:HD12	2.21	0.55
1:A:75:GLU:O	1:A:76:GLU:HB2	2.06	0.55
3:C:18[A]:ASP:OD2	3:C:26:LEU:HG	2.07	0.54
1:A:170:ALA:O	1:A:171:ASP:HB2	2.07	0.54
1:A:316:GLU:OE2	1:A:326:LYS:HD2	2.07	0.54
2:B:280:HIS:HE1	2:B:302:ASP:OD2	1.90	0.54
1:A:469:PRO:O	1:A:470:LEU:HD12	2.07	0.54
1:A:99:THR:HB	7:A:1014:HOH:O	2.08	0.54
3:C:31:GLU:CD	3:C:31:GLU:H	2.11	0.54
1:A:459:VAL:HB	1:A:474:ALA:HB3	1.90	0.53
1:A:182:ARG:HH22	1:A:382:GLU:CD	2.12	0.53
2:B:203:TYR:HB3	2:B:239:MET:CE	2.39	0.53
1:A:113:ARG:HG3	1:A:161:PRO:O	2.08	0.53
1:A:241:ARG:HD2	3:C:79:LYS:HG2	1.91	0.53
1:A:388:ASN:CG	1:A:394:PRO:HD3	2.29	0.53
1:A:6:VAL:HG11	1:A:59:ILE:CG1	2.37	0.52
1:A:56:ARG:O	1:A:59:ILE:HD12	2.09	0.52
2:B:39:PRO:HG2	2:B:331:ARG:HG3	1.90	0.52
1:A:194:VAL:HG22	7:A:1129:HOH:O	2.08	0.52
7:A:1299:HOH:O	3:C:2:ASN:HB2	2.08	0.52
2:B:46:ARG:HH21	2:B:46:ARG:HG3	1.75	0.52
1:A:95:SER:O	1:A:99:THR:CG2	2.58	0.52
1:A:425:LYS:HE3	1:A:426:TYR:CZ	2.45	0.52
1:A:259:ASP:OD1	3:C:77:GLU:OE1	2.27	0.52
2:B:1:ARG:HG3	2:B:1:ARG:NH1	2.25	0.52
2:B:192:HIS:HE1	7:B:1008:HOH:O	1.91	0.52
3:C:75:PHE:O	3:C:77:GLU:N	2.36	0.52
2:B:253:THR:O	2:B:254:ARG:HG3	2.10	0.51
1:A:330:THR:O	1:A:330:THR:HG22	2.10	0.51
1:A:6:VAL:HB	1:A:59:ILE:HG23	1.93	0.51
1:A:427:ALA:HB3	1:A:457:LEU:HD11	1.93	0.51
1:A:408:SER:HB2	1:A:462:THR:CG2	2.40	0.51
1:A:336:VAL:CG2	1:A:347:LEU:HB2	2.40	0.50
2:B:170:LYS:HE3	7:B:1227:HOH:O	2.10	0.50
1:A:409:TRP:CZ3	1:A:461:ALA:HB2	2.46	0.50
3:C:77:GLU:OE2	3:C:77:GLU:O	2.29	0.50
2:B:58:LEU:HD23	2:B:58:LEU:C	2.31	0.50
2:B:7:PRO:HB2	2:B:36:PRO:HG2	1.92	0.50
2:B:85:PHE:CD2	2:B:188:VAL:HG13	2.47	0.50
1:A:6:VAL:CG1	1:A:59:ILE:HG12	2.39	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:291:GLU:HA	1:A:328:THR:HB	1.93	0.50
1:A:298:GLY:O	1:A:321:GLY:HA2	2.12	0.50
3:C:39:ASP:HA	3:C:50:THR:OG1	2.11	0.49
1:A:398:ASP:O	1:A:399:ASP:O	2.29	0.49
1:A:169:TYR:CZ	1:A:171:ASP:HB3	2.48	0.49
2:B:104:ARG:NH2	7:B:1234:HOH:O	2.45	0.49
1:A:36:TRP:O	1:A:40:VAL:HG23	2.13	0.49
1:A:45:ARG:HH11	1:A:46:ASN:HD21	1.59	0.49
1:A:292:THR:HG21	7:A:1266:HOH:O	2.12	0.49
1:A:193:LEU:HD22	1:A:195:LEU:CD1	2.43	0.49
1:A:241:ARG:HD2	3:C:79:LYS:CG	2.42	0.49
1:A:256:HIS:CE1	3:C:79:LYS:HB3	2.47	0.49
1:A:294:LEU:HD12	1:A:327:LEU:HD13	1.94	0.49
1:A:417:GLU:H	1:A:417:GLU:CD	2.14	0.49
1:A:356:ILE:HG21	1:A:463:VAL:HG21	1.95	0.48
2:B:52:VAL:HG13	2:B:53:ASN:N	2.28	0.48
2:B:154[B]:ARG:NH2	7:B:1202:HOH:O	2.45	0.48
3:C:78:THR:HG22	3:C:79:LYS:H	1.78	0.48
1:A:467:GLY:O	1:A:469:PRO:HD3	2.14	0.48
2:B:32:ALA:O	2:B:52:VAL:HG22	2.13	0.48
1:A:117:GLU:O	1:A:120:LYS:HB3	2.14	0.48
2:B:27:ILE:HG21	2:B:61:ILE:HD12	1.96	0.48
1:A:311:VAL:HG21	1:A:327:LEU:HD23	1.95	0.47
2:B:262:LEU:HD23	2:B:263:GLU:N	2.28	0.47
1:A:83:ARG:CZ	3:C:79:LYS:O	2.62	0.47
1:A:291:GLU:OE1	1:A:326:LYS:HE2	2.12	0.47
1:A:393:GLN:OE1	1:A:399:ASP:N	2.47	0.47
2:B:39:PRO:O	2:B:331:ARG:HD3	2.14	0.47
2:B:46:ARG:NH2	2:B:62:ASP:OD2	2.47	0.47
1:A:344:LYS:HE2	1:A:346:ASP:OD1	2.14	0.47
2:B:9:ARG:HD3	3:C:62:ASP:HA	1.95	0.47
2:B:180:TYR:HB3	2:B:204:THR:OG1	2.15	0.47
2:B:204:THR:OG1	2:B:205:ALA:N	2.47	0.47
1:A:116:PRO:HB3	1:A:157:ALA:HB1	1.97	0.47
2:B:17:ASP:HB3	2:B:20:LYS:HG3	1.97	0.47
1:A:192:ARG:NH2	7:A:1295:HOH:O	2.47	0.47
1:A:459:VAL:O	1:A:473:GLU:HA	2.15	0.47
1:A:34:GLU:HB3	1:A:488:ILE:HG13	1.96	0.46
2:B:13:LEU:CD1	2:B:61:ILE:HD11	2.45	0.46
2:B:39:PRO:O	2:B:331:ARG:CD	2.63	0.46
2:B:188:VAL:O	2:B:188:VAL:CG1	2.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:94:LYS:HE2	7:B:1084:HOH:O	2.15	0.46
1:A:94:THR:O	1:A:98:GLN:HG3	2.15	0.46
2:B:283:TYR:OH	3:C:33:ASP:OD1	2.30	0.46
1:A:72:ALA:O	1:A:75:GLU:HG3	2.15	0.46
1:A:327:LEU:N	1:A:327:LEU:HD12	2.31	0.46
1:A:98:GLN:O	1:A:102:ARG:HD3	2.15	0.46
1:A:47:HIS:HE1	4:A:991:HEM:O1D	1.99	0.46
1:A:364:ILE:CG2	3:C:4:LEU:HD22	2.39	0.46
1:A:376:LYS:HB3	1:A:438:PRO:CG	2.44	0.46
2:B:222:LEU:C	2:B:222:LEU:HD23	2.35	0.46
1:A:31:LYS:HG3	1:A:36:TRP:CE2	2.51	0.46
1:A:38:MET:HG2	1:A:487:PRO:HD2	1.98	0.46
1:A:70:SER:OG	1:A:73:GLU:HG3	2.16	0.46
1:A:458:LYS:HD2	1:A:473:GLU:OE1	2.16	0.46
1:A:346:ASP:C	1:A:347:LEU:HD22	2.37	0.45
1:A:337:SER:HB3	1:A:344:LYS:HE3	1.98	0.45
1:A:133:LEU:HD21	1:A:144:TRP:CE2	2.52	0.45
2:B:91:PRO:HG2	7:B:1091:HOH:O	2.17	0.45
2:B:100:GLU:O	2:B:102:PRO:HD3	2.16	0.45
1:A:83:ARG:NH2	1:A:489:ARG:CZ	2.80	0.45
2:B:7:PRO:HG3	2:B:328:ALA:HB1	1.98	0.45
2:B:246:THR:HG22	2:B:257:GLY:HA2	1.99	0.45
3:C:75:PHE:C	3:C:77:GLU:H	2.18	0.44
1:A:178:VAL:HG23	1:A:272:ALA:HA	1.99	0.44
1:A:356:ILE:HA	1:A:384:MET:O	2.17	0.44
1:A:362:LEU:CD2	3:C:4:LEU:HD21	2.48	0.44
2:B:180:TYR:O	2:B:207:LYS:HE3	2.18	0.44
1:A:393:GLN:HE22	1:A:398:ASP:CB	1.98	0.44
2:B:154[A]:ARG:NH2	7:B:1226:HOH:O	2.50	0.44
2:B:52:VAL:HG22	2:B:53:ASN:H	1.83	0.44
1:A:315:VAL:HG22	1:A:325:LEU:CD2	2.47	0.44
1:A:356:ILE:HG13	1:A:472:ALA:HB2	1.99	0.44
2:B:46:ARG:HH21	2:B:46:ARG:CG	2.30	0.44
1:A:153:ILE:HB	1:A:154:PRO:HD3	2.00	0.44
2:B:336:ASP:O	2:B:337:GLU:C	2.55	0.44
1:A:46:ASN:ND2	1:A:46:ASN:N	2.66	0.43
1:A:175:GLY:HA3	1:A:271:LYS:NZ	2.33	0.43
3:C:1:MET:HE3	3:C:8:THR:O	2.18	0.43
2:B:74:LEU:HD22	2:B:100:GLU:OE2	2.18	0.43
1:A:55:GLU:O	1:A:59:ILE:HG13	2.18	0.43
1:A:396:THR:O	1:A:398:ASP:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:ARG:HG3	7:A:1306:HOH:O	2.19	0.43
2:B:266:ASP:HB2	2:B:273:ILE:HD11	2.00	0.43
3:C:37:CYS:HB2	3:C:43:TRQ:HZ3	1.78	0.43
1:A:356:ILE:HD13	1:A:463:VAL:HG22	1.99	0.43
1:A:36:TRP:CE3	1:A:63:LEU:HD13	2.54	0.43
1:A:62:HIS:HE1	1:A:67:ARG:HH12	1.65	0.43
1:A:182:ARG:NH2	1:A:382:GLU:OE2	2.48	0.43
2:B:16:ILE:N	2:B:16:ILE:HD12	2.33	0.43
1:A:83:ARG:NH2	1:A:489:ARG:NE	2.66	0.43
1:A:192:ARG:HD3	7:A:1088:HOH:O	2.19	0.43
1:A:421:MET:O	1:A:422:GLN:HB2	2.19	0.43
2:B:109:HIS:HD2	7:B:1141:HOH:O	2.00	0.43
2:B:223:THR:O	2:B:231:MET:HA	2.19	0.43
1:A:75:GLU:O	1:A:76:GLU:CB	2.67	0.42
2:B:117:VAL:HG23	2:B:133:ALA:HB2	2.01	0.42
1:A:449:MET:HA	2:B:279:PRO:O	2.19	0.42
2:B:217:TYR:CD1	2:B:217:TYR:N	2.88	0.42
2:B:139:MET:HE1	2:B:187:ALA:C	2.40	0.42
3:C:39:ASP:O	3:C:51:LEU:HB2	2.19	0.42
2:B:283:TYR:OH	3:C:33:ASP:HA	2.19	0.42
1:A:45:ARG:HH11	1:A:46:ASN:ND2	2.16	0.42
1:A:116:PRO:HG2	7:A:1028:HOH:O	2.18	0.42
1:A:133:LEU:C	1:A:133:LEU:HD23	2.40	0.42
1:A:427:ALA:CB	1:A:457:LEU:HD11	2.50	0.42
3:C:79:LYS:H	3:C:79:LYS:HG3	1.52	0.42
1:A:338:LEU:C	1:A:338:LEU:HD12	2.40	0.42
1:A:358:ILE:HD11	1:A:472:ALA:HB3	2.01	0.42
2:B:83:SER:HA	2:B:99:TYR:O	2.20	0.42
7:A:1058:HOH:O	2:B:80:ARG:HD2	2.20	0.41
1:A:439:ALA:HB3	1:A:453:ASN:HB3	2.01	0.41
2:B:62:ASP:HB3	2:B:65:THR:HB	2.01	0.41
2:B:262:LEU:HD22	2:B:311:LEU:HD22	2.01	0.41
1:A:466:GLU:C	1:A:468:GLU:N	2.65	0.41
2:B:92:ASP:OD1	2:B:92:ASP:C	2.59	0.41
1:A:127:LEU:HB3	2:B:112:VAL:HG21	2.02	0.41
2:B:86:GLY:HA2	7:B:1024:HOH:O	2.21	0.41
2:B:89:LEU:CD2	2:B:93:GLY:HA2	2.50	0.41
1:A:241:ARG:HD3	3:C:79:LYS:HE2	2.02	0.41
1:A:468:GLU:O	1:A:470:LEU:N	2.53	0.41
2:B:9:ARG:CB	2:B:10:PRO:HA	2.51	0.41
2:B:62:ASP:CG	2:B:65:THR:HB	2.40	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:113:GLN:HG3	7:B:1234:HOH:O	2.21	0.40
1:A:241:ARG:CD	3:C:79:LYS:HE2	2.52	0.40
1:A:356:ILE:HD11	1:A:471:SER:HA	2.04	0.40
1:A:362:LEU:HD23	1:A:363:THR:N	2.37	0.40
1:A:91:GLY:HA2	1:A:106:TYR:CD2	2.56	0.40
1:A:143:ASP:HB2	2:B:154[B]:ARG:NH1	2.37	0.40
2:B:159:MET:O	2:B:161:PRO:HD3	2.22	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	489/489 (100%)	454 (93%)	27 (6%)	8 (2%)	8	2
2	B	337/337 (100%)	320 (95%)	15 (4%)	2 (1%)	22	13
3	C	77/79 (98%)	70 (91%)	5 (6%)	2 (3%)	4	0
All	All	903/905 (100%)	844 (94%)	47 (5%)	12 (1%)	10	4

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	393	GLN
1	A	394	PRO
1	A	399	ASP
1	A	468	GLU
1	A	397	GLY
2	B	259	TYR
3	C	76	PRO
3	C	77	GLU
1	A	395	GLY
2	B	102	PRO

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Mol	Chain	Res	Type
1	A	227	GLY
1	A	288	ILE

### 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	377/375 (100%)	348 (92%)	29 (8%)	10	4
2	B	281/279 (101%)	265 (94%)	16 (6%)	17	10
3	C	65/64 (102%)	58 (89%)	7 (11%)	5	1
All	All	723/718 (101%)	671 (93%)	52 (7%)	12	6

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	46	ASN
1	A	59	ILE
1	A	99	THR
1	A	118	ASP
1	A	127	LEU
1	A	138	LEU
1	A	162	LEU
1	A	192	ARG
1	A	193	LEU
1	A	230	ARG
1	A	246	LEU
1	A	292	THR
1	A	305	LEU
1	A	306	THR
1	A	309	GLU
1	A	328	THR
1	A	340	LEU
1	A	351[A]	ASP
1	A	351[B]	ASP

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Mol	Chain	Res	Type
1	A	352	ARG
1	A	370	ASN
1	A	382	GLU
1	A	393	GLN
1	A	394	PRO
1	A	396	THR
1	A	398	ASP
1	A	456	ASN
1	A	475	HIS
2	B	1	ARG
2	B	9	ARG
2	B	21	MET
2	B	52	VAL
2	B	63	LEU
2	B	89	LEU
2	B	106	GLU
2	B	115	THR
2	B	149	LEU
2	B	152	LEU
2	B	162	GLU
2	B	178	GLU
2	B	186	LEU
2	B	221	LEU
2	B	234	ARG
2	B	300	LEU
3	C	2	ASN
3	C	25	ASN
3	C	34	LEU
3	C	46	GLN
3	C	64	VAL
3	C	77	GLU
3	C	79	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	9	ASN
1	A	17	GLN
1	A	46	ASN
1	A	47	HIS
1	A	62	HIS

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Mol	Chain	Res	Type
1	A	136	GLN
1	A	149	GLN
1	A	242	GLN
1	A	256	HIS
1	A	277	GLN
1	A	293	GLN
1	A	320	ASN
1	A	343	GLN
1	A	370	ASN
1	A	393	GLN
1	A	422	GLN
1	A	452	ASN
1	A	453	ASN
1	A	456	ASN
1	A	475	HIS
2	B	109	HIS
2	B	192	HIS
2	B	280	HIS
3	C	25	ASN
3	C	28	GLN
3	C	46	GLN
3	C	72	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	TRQ	C	43	3,6	13,17,18	4.46	4 (30%)	13,24,26	2.46	4 (30%)

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	TRQ	C	43	3,6	-	0/4/19/21	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	43	TRQ	CH2-CZ2	-13.29	1.37	1.53
3	C	43	TRQ	CE2-CZ2	-7.86	1.40	1.50
3	C	43	TRQ	CZ3-CE3	3.42	1.40	1.34
3	C	43	TRQ	CD2-CE3	-2.00	1.40	1.44

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	43	TRQ	CZ2-CE2-NE1	6.37	130.11	119.94
3	C	43	TRQ	O7-CZ2-CH2	3.42	122.80	119.01
3	C	43	TRQ	CD1-CG-CD2	2.46	106.20	104.79
3	C	43	TRQ	CD2-CE3-CZ3	-2.16	118.63	121.05

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	43	TRQ	1	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 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	TBU	B	994	-	4,4,4	0.50	0	6,6,6	0.55	0
5	TBU	A	995	-	4,4,4	0.46	0	6,6,6	0.51	0
4	HEM	A	991	1	42,50,50	1.88	10 (23%)	46,82,82	1.22	6 (13%)
5	TBU	C	993	6	4,4,4	0.46	0	6,6,6	0.57	0
4	HEM	A	992	1	42,50,50	1.84	11 (26%)	46,82,82	1.25	6 (13%)

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
4	HEM	A	992	1	-	4/12/54/54	-
4	HEM	A	991	1	-	6/12/54/54	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	991	HEM	CBB-CAB	5.22	1.55	1.30
4	A	992	HEM	CBB-CAB	4.96	1.54	1.30
4	A	992	HEM	C3C-C2C	-4.78	1.33	1.40
4	A	991	HEM	C3C-C2C	-4.36	1.34	1.40
4	A	992	HEM	CBC-CAC	4.02	1.54	1.29
4	A	991	HEM	CBC-CAC	3.98	1.54	1.29
4	A	991	HEM	C3C-C4C	3.47	1.46	1.41
4	A	992	HEM	C3C-C4C	3.47	1.46	1.41
4	A	991	HEM	C3C-CAC	3.30	1.55	1.47
4	A	992	HEM	C3C-CAC	2.82	1.54	1.47
4	A	991	HEM	O2A-CGA	-2.47	1.22	1.30
4	A	992	HEM	CAB-C3B	2.47	1.54	1.47
4	A	991	HEM	CAB-C3B	2.27	1.53	1.47
4	A	992	HEM	O2A-CGA	-2.20	1.23	1.30
4	A	991	HEM	CMD-C2D	2.16	1.55	1.50
4	A	992	HEM	CMD-C2D	2.15	1.55	1.50
4	A	992	HEM	CMC-C2C	2.12	1.56	1.51
4	A	991	HEM	O1A-CGA	2.11	1.29	1.22

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	991	HEM	C2C-C1C	2.09	1.47	1.42
4	A	992	HEM	C2C-C1C	2.03	1.47	1.42
4	A	992	HEM	O1A-CGA	2.01	1.28	1.22

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	992	HEM	C4C-CHD-C1D	3.85	127.64	122.56
4	A	991	HEM	C4C-CHD-C1D	3.78	127.55	122.56
4	A	992	HEM	CBB-CAB-C3B	-3.10	112.01	127.53
4	A	991	HEM	C4B-CHC-C1C	2.92	126.42	122.56
4	A	991	HEM	CBB-CAB-C3B	-2.89	113.09	127.53
4	A	992	HEM	C4B-CHC-C1C	2.86	126.34	122.56
4	A	992	HEM	CMA-C3A-C4A	-2.86	124.27	128.46
4	A	991	HEM	O1A-CGA-CBA	-2.57	114.95	123.09
4	A	991	HEM	CMA-C3A-C4A	-2.52	124.77	128.46
4	A	991	HEM	CMC-C2C-C3C	2.39	129.46	124.68
4	A	992	HEM	CMC-C2C-C3C	2.35	129.38	124.68
4	A	992	HEM	O1A-CGA-CBA	-2.30	115.80	123.09

There are no chirality outliers.

All (10) torsion outliers are listed below:

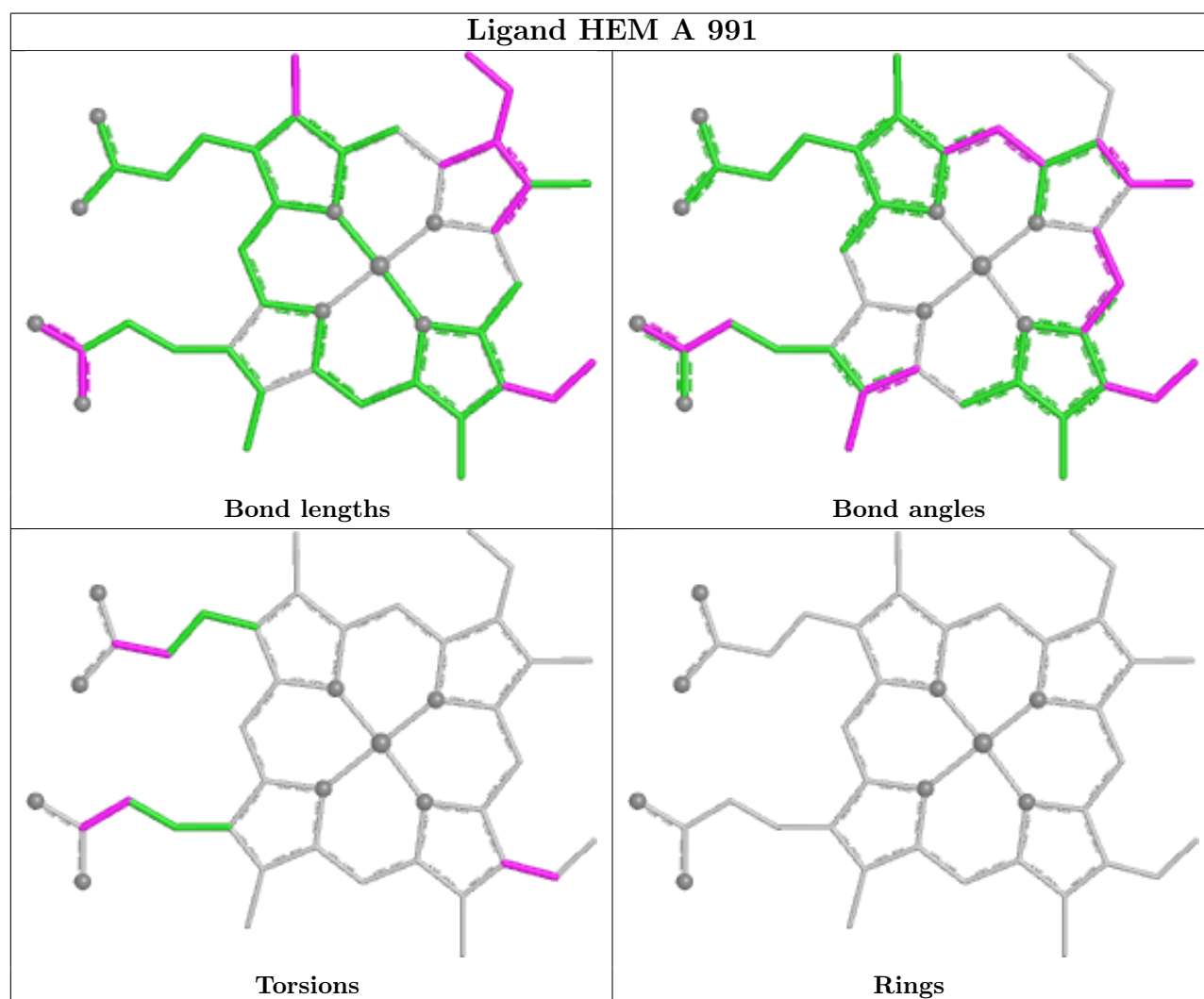
Mol	Chain	Res	Type	Atoms
4	A	991	HEM	C2B-C3B-CAB-CBB
4	A	991	HEM	C4B-C3B-CAB-CBB
4	A	992	HEM	C2B-C3B-CAB-CBB
4	A	992	HEM	C4B-C3B-CAB-CBB
4	A	992	HEM	CAD-CBD-CGD-O2D
4	A	991	HEM	CAA-CBA-CGA-O2A
4	A	992	HEM	CAD-CBD-CGD-O1D
4	A	991	HEM	CAA-CBA-CGA-O1A
4	A	991	HEM	CAD-CBD-CGD-O1D
4	A	991	HEM	CAD-CBD-CGD-O2D

There are no ring outliers.

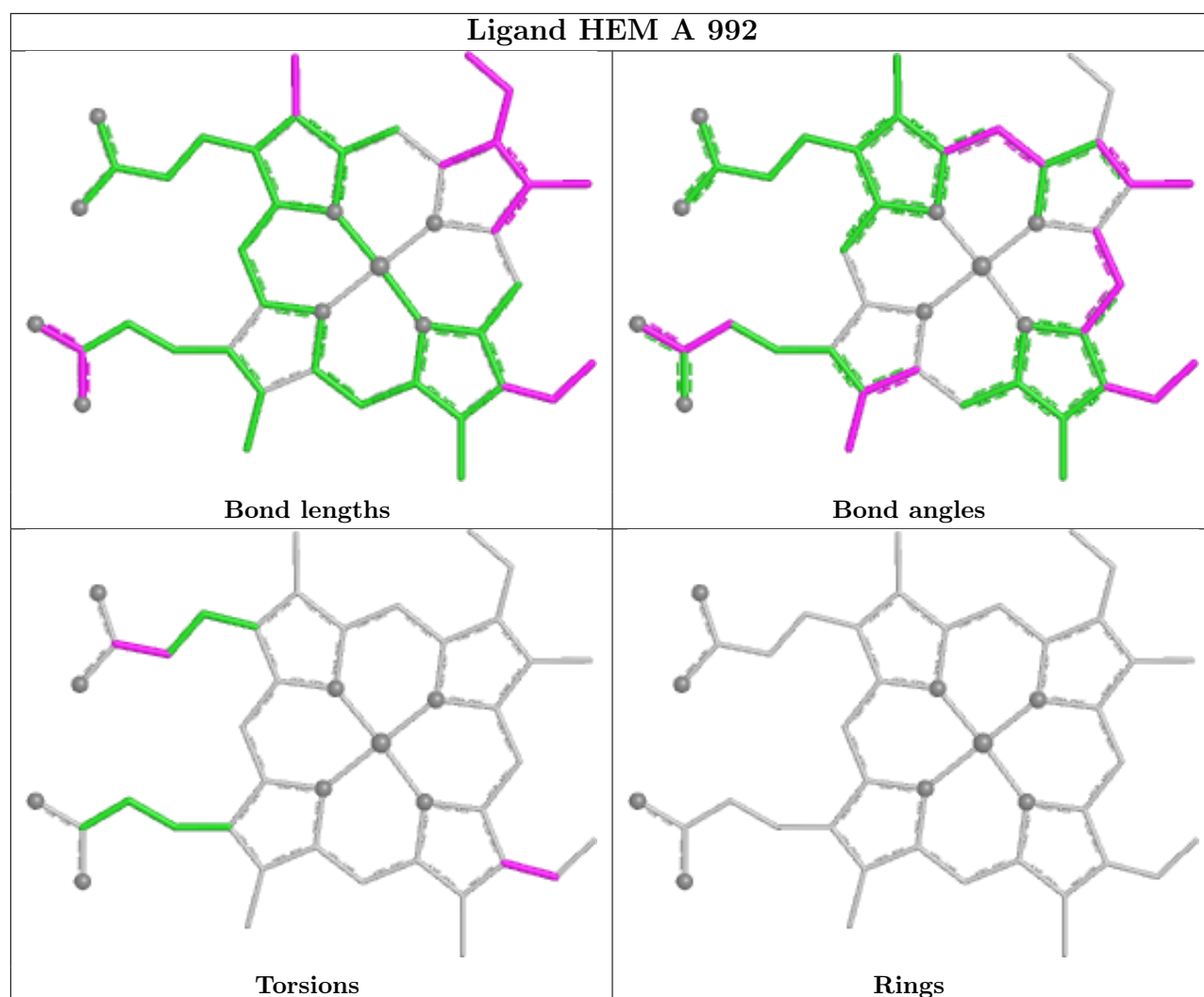
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	991	HEM	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	489/489 (100%)	-0.01	13 (2%) 56 57	23, 36, 67, 93	2 (0%)
2	B	337/337 (100%)	-0.38	0 100 100	14, 32, 49, 90	2 (0%)
3	C	78/79 (98%)	0.05	6 (7%) 21 23	16, 32, 68, 92	1 (1%)
All	All	904/905 (99%)	-0.15	19 (2%) 63 66	14, 34, 61, 93	5 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	1	MET	7.1
3	C	79	LYS	6.7
3	C	78	THR	6.1
1	A	469	PRO	5.8
1	A	470	LEU	4.3
1	A	467	GLY	3.5
3	C	2	ASN	3.2
1	A	395	GLY	3.1
3	C	77	GLU	3.0
1	A	465	ALA	3.0
1	A	392	GLY	2.9
1	A	351[A]	ASP	2.9
1	A	460	ILE	2.6
3	C	76	PRO	2.6
1	A	466	GLU	2.4
1	A	170	ALA	2.3
1	A	59	ILE	2.2
1	A	390	PRO	2.2
1	A	468	GLU	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	TRQ	C	43	16/17	0.93	0.07	24,28,30,31	0

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

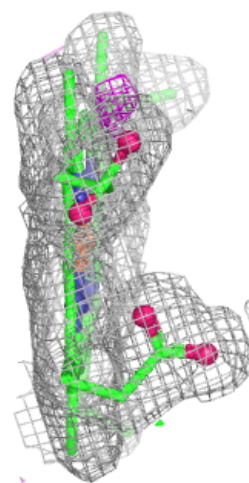
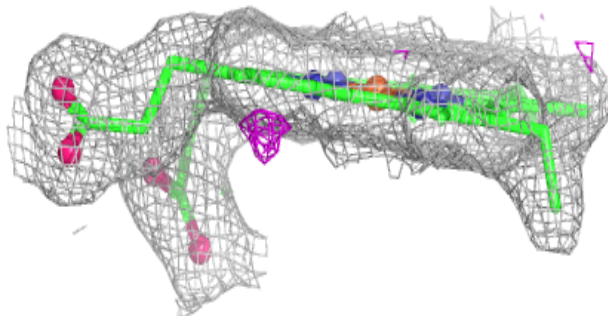
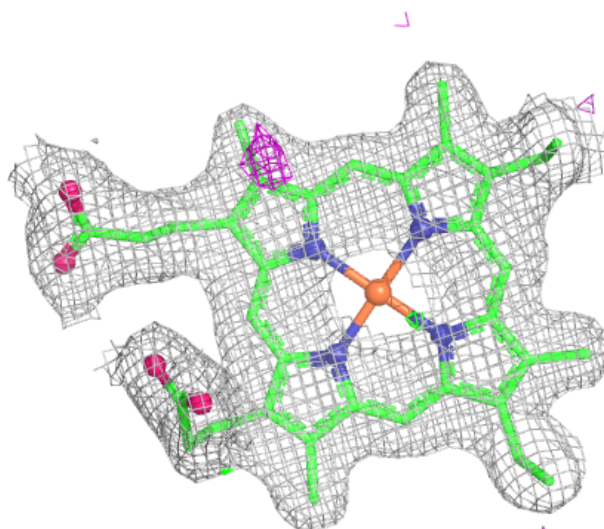
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

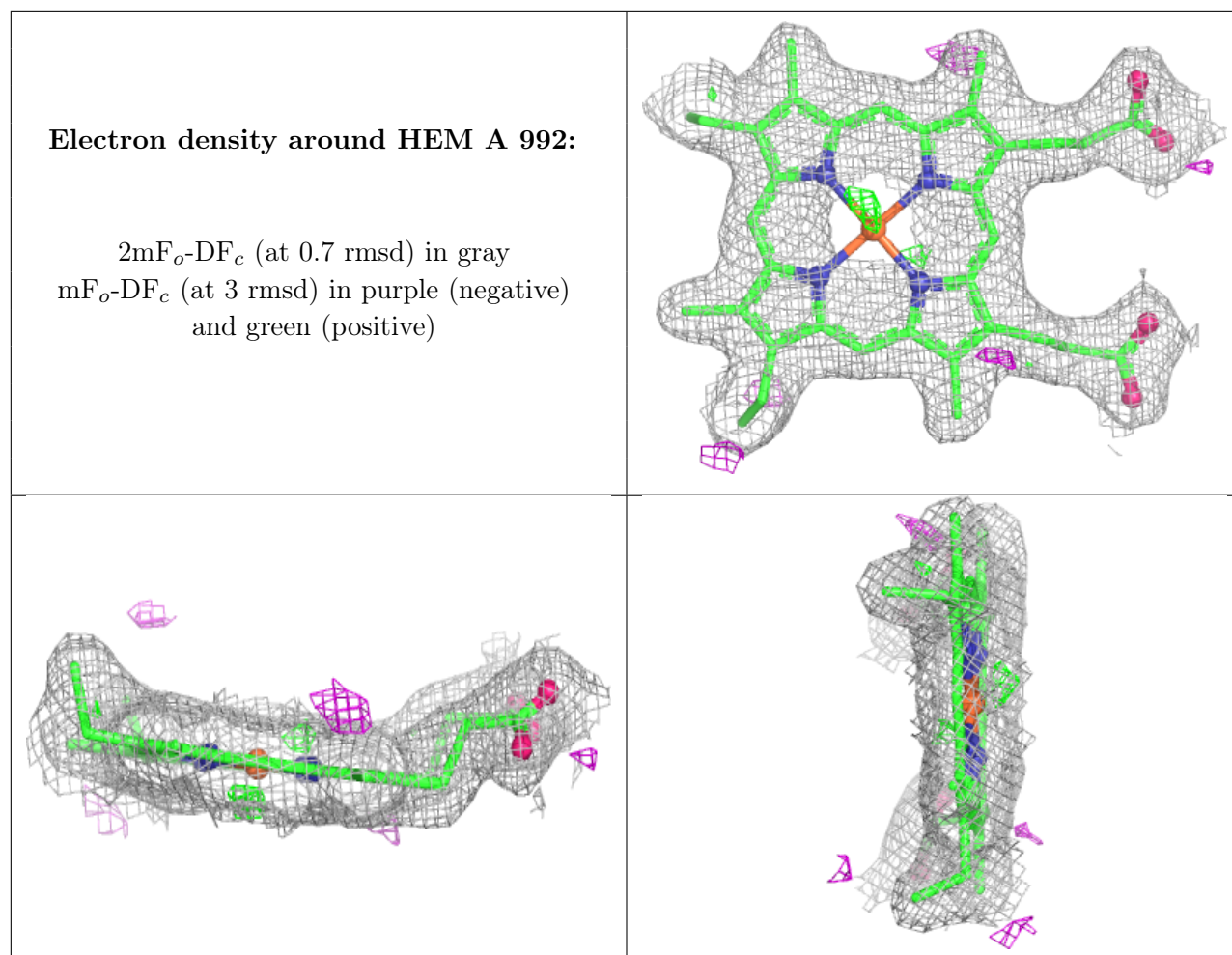
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	TBU	B	994	5/5	0.83	0.16	59,60,63,64	0
5	TBU	A	995	5/5	0.92	0.13	44,45,50,53	0
5	TBU	C	993	5/5	0.93	0.12	43,48,49,54	0
6	NA	C	996	1/1	0.95	0.15	51,51,51,51	0
4	HEM	A	991	43/43	0.97	0.07	21,29,33,36	0
4	HEM	A	992	43/43	0.97	0.07	22,27,33,34	0

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

**Electron density around HEM A 991:**

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





## 6.5 Other polymers ⓘ

There are no such residues in this entry.