



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:18 AM UTC

PDB ID : 1BVY / pdb_00001bvy
Title : COMPLEX OF THE HEME AND FMN-BINDING DOMAINS OF THE CYTOCHROME P450(BM-3)
Authors : Sevrioukova, I.F.; Li, H.; Zhang, H.; Peterson, J.A.; Poulos, T.L.
Deposited on : 1998-09-21
Resolution : 2.03 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

<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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

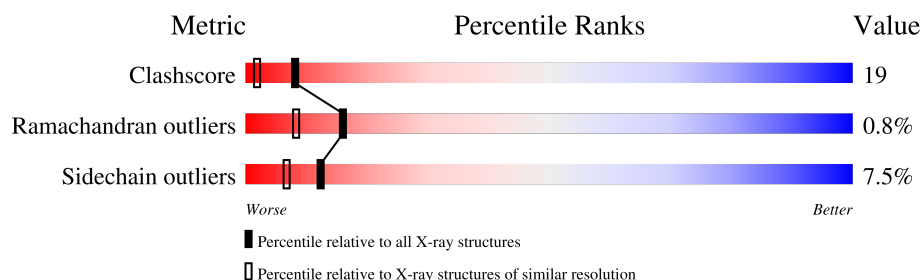
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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))
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)

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 ≥ 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	458	 66% 26% • •
1	B	458	 65% 26% 5% •
2	F	191	 39% 34% 7% 20%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9324 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.

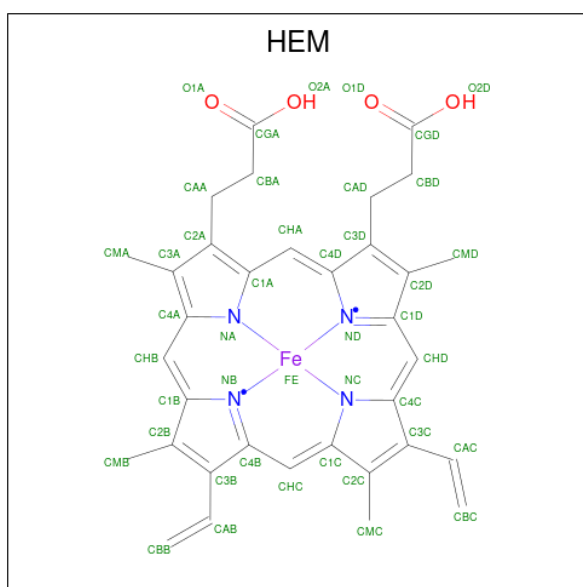
- Molecule 1 is a protein called PROTEIN (CYTOCHROME P450 BM-3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	439	Total	C	N	O	S	0	4	0
			3551	2263	606	665	17			
1	B	438	Total	C	N	O	S	0	0	0
			3527	2249	601	661	16			

- Molecule 2 is a protein called PROTEIN (CYTOCHROME P450 BM-3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	152	Total	C	N	O	S	0	0	0
			1162	728	195	235	4			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



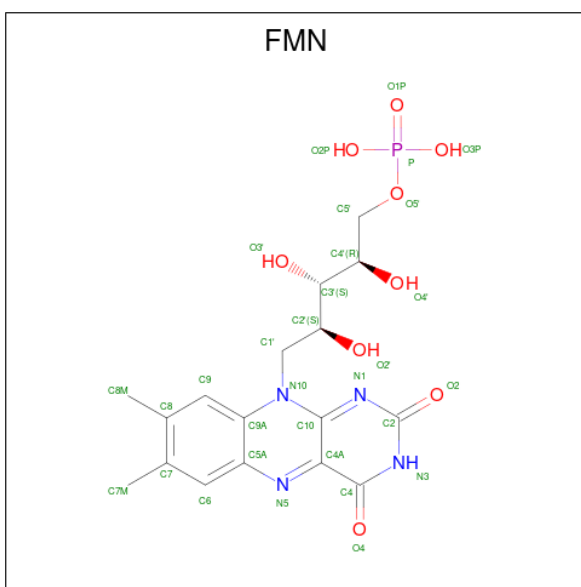
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
3	B	1	Total	C	Fe	N	O	
			43	34	1	4	4	

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O		
			4	2	2	0	0
4	A	1	Total	C	O		
			4	2	2	0	0
4	B	1	Total	C	O		
			4	2	2	0	0
4	B	1	Total	C	O		
			4	2	2	0	0
4	B	1	Total	C	O		
			4	2	2	0	0

- Molecule 5 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: C₁₇H₂₁N₄O₉P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	F	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 6 is water.

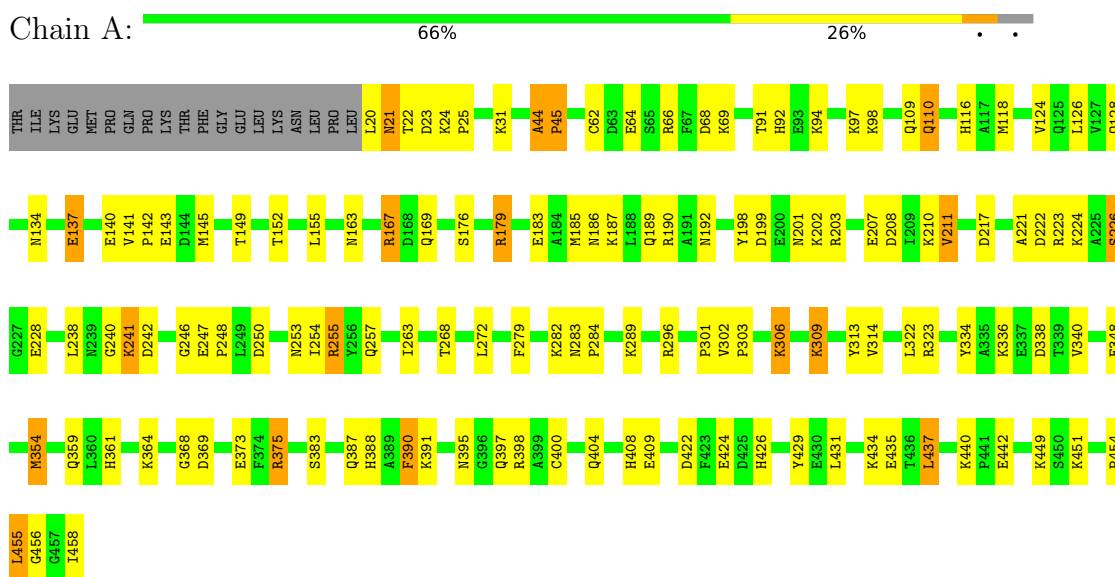
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	460	Total O 460 460	0	0
6	F	80	Total O 80 80	0	0
6	B	407	Total O 407 407	0	0

3 Residue-property plots [i](#)

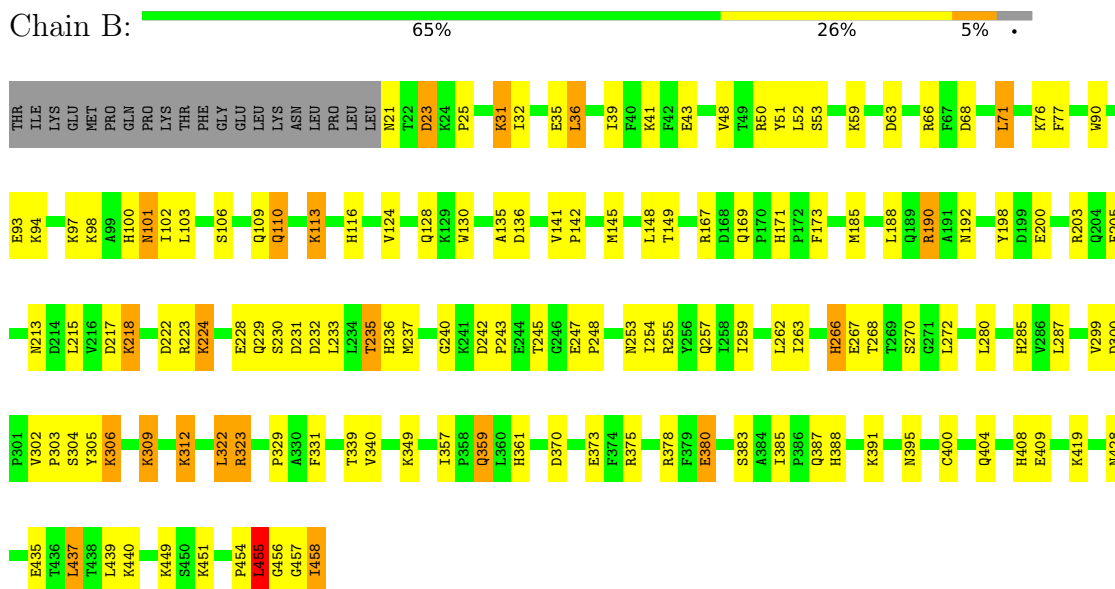
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

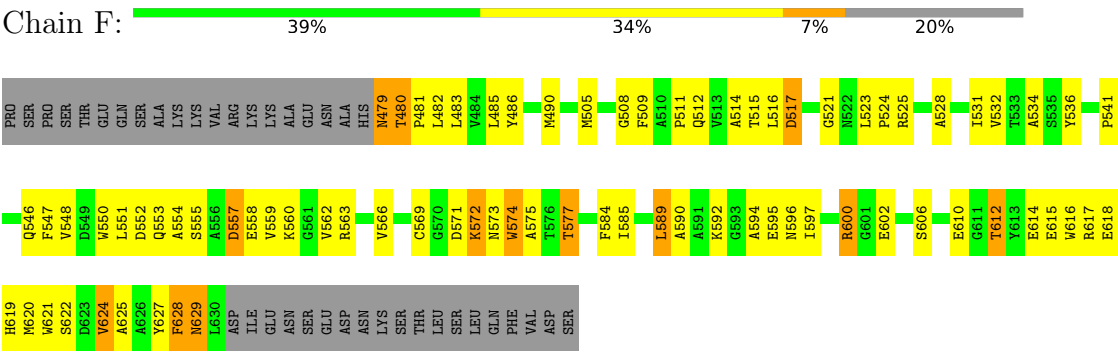
- Molecule 1: PROTEIN (CYTOCHROME P450 BM-3)



- Molecule 1: PROTEIN (CYTOCHROME P450 BM-3)



● Molecule 2: PROTEIN (CYTOCHROME P450 BM-3)



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.84Å 94.68Å 209.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.03	Depositor
% Data completeness (in resolution range)	83.0 (10.00-2.03)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.187 , 0.275	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9324	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.39	0/3652	1.26	7/4934 (0.1%)
1	B	0.40	0/3608	1.24	13/4876 (0.3%)
2	F	0.44	0/1190	1.30	6/1621 (0.4%)
All	All	0.40	0/8450	1.26	26/11431 (0.2%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	323	ARG	CD-NE-CZ	8.56	136.38	124.40
1	A	44	ALA	O-C-N	8.23	127.07	121.14
1	A	179	ARG	CD-NE-CZ	8.22	135.91	124.40
1	A	44	ALA	CA-C-O	-7.90	114.05	121.08
2	F	574	TRP	CA-C-N	7.89	136.60	121.54
2	F	574	TRP	C-N-CA	7.89	136.60	121.54
1	B	331	PHE	CA-CB-CG	7.65	121.45	113.80
1	B	455	LEU	CA-C-N	7.50	128.82	121.62
1	B	455	LEU	C-N-CA	7.50	128.82	121.62
2	F	517	ASP	CA-CB-CG	6.89	119.49	112.60
1	B	285	HIS	CA-CB-CG	-6.54	107.26	113.80
2	F	600	ARG	CA-C-N	6.41	126.01	121.65
2	F	600	ARG	C-N-CA	6.41	126.01	121.65
1	B	357	ILE	CA-C-O	-6.09	113.67	118.85
2	F	628	PHE	CA-CB-CG	5.67	119.47	113.80
1	B	266	HIS	CA-CB-CG	5.51	119.31	113.80
1	B	63	ASP	CA-CB-CG	5.34	117.94	112.60
1	B	68	ASP	CA-CB-CG	5.33	117.93	112.60
1	B	101	ASN	CA-CB-CG	-5.32	107.28	112.60
1	A	314	VAL	CA-C-N	5.23	125.78	119.98
1	A	314	VAL	C-N-CA	5.23	125.78	119.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	ARG	O-C-N	5.21	125.44	121.27
1	B	167	ARG	CA-C-N	5.18	129.28	120.88
1	B	167	ARG	C-N-CA	5.18	129.28	120.88
1	A	390	PHE	CA-CB-CG	5.15	118.95	113.80
1	B	323	ARG	NE-CZ-NH2	-5.06	114.65	119.20

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	3551	0	3512	119	0
1	B	3527	0	3483	135	0
2	F	1162	0	1077	75	0
3	A	43	0	30	4	0
3	B	43	0	30	3	0
4	A	8	0	11	1	0
4	B	12	0	17	1	0
5	F	31	0	19	1	0
6	A	460	0	0	30	1
6	B	407	0	0	27	1
6	F	80	0	0	6	0
All	All	9324	0	8179	322	1

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 (322) 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:B:280:LEU:HB3	1:B:287:LEU:HD23	1.55	0.89
2:F:590:ALA:HA	2:F:594:ALA:HB2	1.53	0.88
1:B:272:LEU:HD13	1:B:322:LEU:HD13	1.53	0.88
1:B:66:ARG:HE	1:B:339:THR:HG21	1.41	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:GLN:HG3	6:A:1465:HOH:O	1.74	0.86
2:F:517:ASP:HB2	2:F:546:GLN:HB2	1.61	0.82
6:A:1114:HOH:O	1:B:458:ILE:HD11	1.80	0.82
1:B:361:HIS:HE1	1:B:391:LYS:H	1.28	0.81
1:B:387:GLN:HB3	6:B:1112:HOH:O	1.84	0.77
1:B:217:ASP:OD1	1:B:255:ARG:HD3	1.86	0.76
1:A:179:ARG:HD2	1:A:208:ASP:OD2	1.87	0.75
1:A:145:MET:O	1:A:149:THR:HG23	1.87	0.75
1:B:218:LYS:HE2	1:B:222:ASP:OD2	1.87	0.74
1:B:232:ASP:H	1:B:235:THR:HG23	1.52	0.74
1:B:245:THR:OG1	1:B:247:GLU:HG2	1.88	0.73
1:A:361:HIS:HE1	1:A:391:LYS:H	1.36	0.73
1:B:59:LYS:HD2	6:B:1339:HOH:O	1.88	0.73
1:A:364:LYS:HG2	1:A:369:ASP:HA	1.73	0.71
1:B:149:THR:HG22	1:B:409:GLU:OE1	1.91	0.70
2:F:550:TRP:O	2:F:554:ALA:HB3	1.91	0.70
2:F:482:LEU:O	2:F:511:PRO:HA	1.92	0.69
1:B:171:HIS:HD2	1:B:173:PHE:H	1.38	0.69
6:A:1433:HOH:O	1:B:454:PRO:HB3	1.92	0.69
2:F:523:LEU:HD23	2:F:562:VAL:HG11	1.74	0.69
2:F:548:VAL:HG13	2:F:584:PHE:CE2	2.26	0.69
1:A:223:ARG:HD2	1:A:228:GLU:HA	1.75	0.69
2:F:485:LEU:HD22	2:F:515:THR:O	1.95	0.67
1:B:437:LEU:HD23	6:B:1160:HOH:O	1.94	0.67
1:A:149:THR:HG22	6:A:1348:HOH:O	1.94	0.67
2:F:528:ALA:HA	2:F:563:ARG:O	1.93	0.67
1:B:53:SER:HB2	1:B:359:GLN:HB3	1.76	0.66
1:A:124:VAL:HG11	1:B:458:ILE:HG12	1.77	0.66
2:F:615:GLU:HB3	6:F:1070:HOH:O	1.95	0.65
1:A:217:ASP:OD1	1:A:255:ARG:HD3	1.97	0.65
1:B:223:ARG:NH1	1:B:235:THR:HG22	2.12	0.65
1:A:207:GLU:O	1:A:211:VAL:HG12	1.96	0.65
1:A:400:CYS:HB2	3:A:1000:HEM:NA	2.13	0.64
1:A:364:LYS:HG2	1:A:368:GLY:O	1.97	0.63
1:B:110:GLN:H	1:B:110:GLN:NE2	1.95	0.63
1:A:222:ASP:O	1:A:226:SER:HB3	1.98	0.63
2:F:563:ARG:HG2	2:F:595:GLU:CB	2.28	0.62
2:F:490:MET:HE3	2:F:536:TYR:HE2	1.65	0.62
2:F:516:LEU:HD21	2:F:531:ILE:HG23	1.82	0.61
2:F:482:LEU:HB2	2:F:509:PHE:CD2	2.35	0.61
2:F:595:GLU:OE2	2:F:627:TYR:OH	2.18	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:PRO:HD2	6:B:1371:HOH:O	2.00	0.61
2:F:571:ASP:OD1	2:F:573:ASN:HB2	2.00	0.60
1:A:369:ASP:HB2	6:A:1175:HOH:O	2.00	0.60
1:A:128:GLN:NE2	1:A:455:LEU:HD11	2.17	0.60
2:F:616:TRP:O	2:F:620:MET:HB2	2.02	0.60
1:A:118[A]:MET:SD	1:A:155:LEU:HD21	2.42	0.59
2:F:563:ARG:HG2	2:F:595:GLU:HB3	1.83	0.59
1:B:242:ASP:OD2	1:B:245:THR:HG23	2.02	0.59
2:F:629:ASN:HD22	2:F:629:ASN:H	1.51	0.59
1:B:391:LYS:NZ	1:B:395:ASN:HD22	2.01	0.58
1:B:437:LEU:HB2	6:B:1041:HOH:O	2.02	0.58
1:B:110:GLN:NE2	1:B:110:GLN:N	2.51	0.58
1:A:301:PRO:HB3	1:A:456:GLY:H	1.66	0.58
2:F:523:LEU:HD11	2:F:547:PHE:HE1	1.69	0.58
1:A:124:VAL:HG13	1:A:455:LEU:HD22	1.84	0.58
1:B:66:ARG:NE	1:B:339:THR:HG21	2.16	0.58
1:A:458:ILE:HD12	1:A:458:ILE:H	1.68	0.58
1:A:391:LYS:HZ3	1:A:395:ASN:HD22	1.52	0.57
1:B:98:LYS:O	1:B:102:ILE:HG13	2.04	0.57
1:A:118[B]:MET:HG3	6:A:1300:HOH:O	2.05	0.57
1:B:329:PRO:HG3	1:B:439:LEU:HG	1.87	0.57
1:A:69:LYS:HD3	1:A:398:ARG:CZ	2.34	0.57
1:A:124:VAL:HG11	1:B:458:ILE:CD1	2.36	0.56
1:B:171:HIS:CD2	1:B:173:PHE:H	2.21	0.56
1:B:242:ASP:HB3	1:B:245:THR:OG1	2.04	0.56
1:A:296[B]:ARG:HD3	6:A:1075:HOH:O	2.05	0.56
1:A:176:SER:OG	1:A:211:VAL:HG11	2.05	0.56
1:B:205:PHE:HD2	6:B:1157:HOH:O	1.88	0.56
1:A:25:PRO:HD2	1:A:435:GLU:OE1	2.06	0.56
1:A:92:HIS:HA	1:A:97:LYS:HE2	1.88	0.56
1:B:97:LYS:HZ2	1:B:97:LYS:HB3	1.71	0.56
1:B:266:HIS:NE2	1:B:267:GLU:HG3	2.21	0.56
1:B:35:GLU:HG3	1:B:36:LEU:N	2.20	0.56
2:F:541:PRO:HA	6:F:1024:HOH:O	2.05	0.55
1:B:100:HIS:HD2	6:B:1118:HOH:O	1.88	0.55
1:B:190:ARG:HD2	1:B:198:TYR:CE1	2.41	0.55
2:F:490:MET:HE3	2:F:536:TYR:CE2	2.41	0.55
1:A:21:ASN:N	1:A:21:ASN:OD1	2.39	0.55
1:A:253:ASN:O	1:A:257:GLN:HG2	2.07	0.55
1:A:409:GLU:HG2	6:A:1348:HOH:O	2.07	0.55
1:A:224:LYS:HG2	1:A:238:LEU:HD22	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:485:LEU:HA	2:F:514:ALA:O	2.05	0.55
1:B:124:VAL:HG13	1:B:455:LEU:CD1	2.36	0.55
1:B:231:ASP:CG	1:B:236:HIS:HE2	2.15	0.54
1:A:391:LYS:NZ	1:A:395:ASN:HD22	2.04	0.54
1:B:458:ILE:HG22	1:B:458:ILE:O	2.07	0.54
1:B:304:SER:HB2	1:B:306:LYS:HZ2	1.73	0.54
2:F:482:LEU:HB2	2:F:509:PHE:HD2	1.73	0.53
2:F:551:LEU:HD11	2:F:589:LEU:HD13	1.90	0.53
1:B:435:GLU:HG2	1:B:439:LEU:CD2	2.38	0.53
1:A:387:GLN:O	1:A:388:HIS:HB2	2.07	0.53
1:A:301:PRO:HB3	1:A:456:GLY:N	2.24	0.53
1:B:90:TRP:HB2	1:B:93:GLU:HG3	1.90	0.53
1:A:309:LYS:HE2	6:A:1053:HOH:O	2.08	0.53
2:F:629:ASN:H	2:F:629:ASN:ND2	2.06	0.53
1:B:50:ARG:HD2	6:B:1109:HOH:O	2.09	0.53
2:F:585:ILE:O	2:F:589:LEU:HB2	2.09	0.53
1:B:259:ILE:O	1:B:263:ILE:HG12	2.08	0.52
1:A:426:HIS:HE1	6:A:1349:HOH:O	1.92	0.52
1:B:323:ARG:HA	1:B:361:HIS:CD2	2.45	0.52
1:A:137:GLU:HB3	6:A:1437:HOH:O	2.10	0.52
1:A:250:ASP:O	1:A:254:ILE:HG13	2.09	0.52
1:A:289:LYS:HE3	1:A:313:TYR:CE1	2.45	0.52
1:B:124:VAL:O	1:B:128:GLN:HG3	2.10	0.52
1:B:230:SER:C	1:B:235:THR:HG21	2.34	0.52
1:A:116:HIS:HD2	1:A:408:HIS:NE2	2.08	0.52
1:B:240:GLY:O	1:B:248:PRO:HB3	2.09	0.52
1:A:296[B]:ARG:NH1	6:A:1464:HOH:O	2.42	0.52
1:B:94:LYS:HG3	6:B:1282:HOH:O	2.09	0.52
2:F:562:VAL:HA	6:F:1064:HOH:O	2.10	0.51
1:A:62:CYS:SG	1:A:391:LYS:HE2	2.50	0.51
1:A:185:MET:SD	1:A:437:LEU:HD13	2.50	0.51
1:A:109:GLN:HG2	6:A:1053:HOH:O	2.11	0.51
1:B:106:SER:HB3	1:B:233:LEU:HD23	1.92	0.51
1:B:145:MET:O	1:B:149:THR:HG23	2.10	0.51
2:F:479:ASN:HA	2:F:508:GLY:O	2.10	0.51
2:F:618:GLU:HA	2:F:618:GLU:OE1	2.11	0.51
1:A:223:ARG:HG3	1:A:228:GLU:HB3	1.93	0.51
1:B:266:HIS:CE1	1:B:267:GLU:HG3	2.45	0.51
1:A:400:CYS:HB2	3:A:1000:HEM:C4A	2.45	0.51
1:B:130:TRP:O	4:B:1007:EDO:H12	2.10	0.51
1:B:457:GLY:O	1:B:458:ILE:HB	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:77:PHE:HE2	1:B:188:LEU:HD23	1.76	0.51
1:B:185:MET:HE3	1:B:437:LEU:HD22	1.91	0.51
1:B:110:GLN:HB3	6:B:1176:HOH:O	2.11	0.51
1:B:116:HIS:HE1	1:B:303:PRO:O	1.92	0.51
1:B:116:HIS:HD2	1:B:408:HIS:NE2	2.08	0.51
1:A:124:VAL:HG11	1:B:458:ILE:CG1	2.41	0.51
1:A:134:ASN:O	1:A:137:GLU:HG3	2.11	0.51
1:B:388:HIS:HD2	1:B:391:LYS:HZ3	1.59	0.51
1:A:140:GLU:HB3	1:A:143:GLU:OE1	2.10	0.50
1:B:361:HIS:CE1	1:B:391:LYS:H	2.18	0.50
2:F:563:ARG:N	6:F:1064:HOH:O	2.44	0.50
2:F:563:ARG:HG2	2:F:595:GLU:CG	2.41	0.50
1:B:200:GLU:OE1	1:B:203:ARG:NH1	2.45	0.50
1:B:113:LYS:HG3	1:B:305:TYR:CE2	2.47	0.50
1:A:208:ASP:O	1:A:211:VAL:HG13	2.11	0.50
1:A:240:GLY:O	1:A:248:PRO:HB3	2.12	0.50
1:A:361:HIS:CE1	1:A:391:LYS:H	2.23	0.50
1:A:373:GLU:HA	1:A:373:GLU:OE1	2.10	0.50
1:A:458:ILE:HG22	1:A:458:ILE:OXT	2.11	0.50
1:A:169:GLN:OE1	1:B:169:GLN:OE1	2.30	0.49
2:F:571:ASP:HB3	2:F:574:TRP:HE3	1.76	0.49
2:F:597:ILE:HG22	2:F:597:ILE:O	2.11	0.49
1:A:98:LYS:HZ2	1:A:247:GLU:HG3	1.77	0.49
1:A:98:LYS:NZ	1:A:247:GLU:HG3	2.26	0.49
1:A:454:PRO:HB3	6:A:1395:HOH:O	2.12	0.49
1:B:110:GLN:O	1:B:113:LYS:HD3	2.12	0.49
1:B:43:GLU:CD	1:B:48:VAL:HG22	2.37	0.49
1:B:400:CYS:HA	3:B:1001:HEM:CHA	2.42	0.49
1:A:109:GLN:HG3	1:A:404[B]:GLN:HE21	1.77	0.49
1:A:128:GLN:HG3	1:A:455:LEU:HD21	1.94	0.49
1:B:440:LYS:HD3	6:B:1378:HOH:O	2.12	0.48
1:A:192:ASN:HB2	6:A:1257:HOH:O	2.12	0.48
1:B:230:SER:OG	1:B:231:ASP:N	2.41	0.48
2:F:486:TYR:HA	2:F:532:VAL:O	2.13	0.48
1:B:71:LEU:HD22	1:B:90:TRP:CE2	2.49	0.48
2:F:547:PHE:O	2:F:550:TRP:HE3	1.96	0.48
1:B:113:LYS:HD2	1:B:305:TYR:CZ	2.48	0.48
1:A:152:THR:HG21	1:A:409:GLU:OE2	2.12	0.48
2:F:483:LEU:HD12	2:F:512:GLN:O	2.14	0.48
1:A:241:LYS:HD2	1:A:246:GLY:O	2.13	0.48
1:B:185:MET:CE	1:B:437:LEU:HD22	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:419:LYS:O	1:B:451:LYS:HD2	2.12	0.48
1:B:391:LYS:HZ3	1:B:395:ASN:HD22	1.61	0.48
1:B:309:LYS:NZ	6:B:1044:HOH:O	2.47	0.48
1:A:272:LEU:HD13	1:A:322:LEU:HG	1.95	0.48
1:B:128:GLN:NE2	1:B:457:GLY:O	2.47	0.48
1:B:309:LYS:O	1:B:312:LYS:HE2	2.14	0.48
1:B:449:LYS:HE2	6:B:1172:HOH:O	2.13	0.48
2:F:523:LEU:HD11	2:F:547:PHE:CE1	2.49	0.47
1:B:103:LEU:HD21	1:B:237:MET:CG	2.44	0.47
1:B:224:LYS:HD3	6:B:1351:HOH:O	2.14	0.47
1:B:388:HIS:CD2	1:B:391:LYS:HZ3	2.32	0.47
2:F:534:ALA:HA	2:F:569:CYS:O	2.15	0.47
2:F:571:ASP:HB3	2:F:574:TRP:CE3	2.48	0.47
2:F:614:GLU:HA	2:F:617:ARG:NH2	2.28	0.47
1:A:309:LYS:NZ	6:A:1457:HOH:O	2.46	0.47
1:A:434:LYS:NZ	6:A:1185:HOH:O	2.47	0.47
1:B:349:LYS:NZ	6:B:1332:HOH:O	2.47	0.47
1:A:187:LYS:NZ	6:A:1161:HOH:O	2.47	0.47
1:A:361:HIS:CE1	1:A:390:PHE:HA	2.49	0.47
2:F:517:ASP:HA	2:F:547:PHE:HB2	1.97	0.47
2:F:559:VAL:HG22	2:F:592:LYS:O	2.15	0.47
1:B:173:PHE:CE1	1:B:262:LEU:HD13	2.50	0.47
2:F:479:ASN:OD1	2:F:479:ASN:N	2.48	0.47
1:B:280:LEU:HD22	1:B:287:LEU:HA	1.96	0.47
6:A:1465:HOH:O	2:F:490:MET:HE2	2.15	0.46
1:B:190:ARG:HD2	1:B:198:TYR:CZ	2.50	0.46
1:A:183:GLU:CD	1:A:190:ARG:HH22	2.24	0.46
1:A:306:LYS:HD3	6:A:1382:HOH:O	2.14	0.46
1:B:77:PHE:CE2	1:B:188:LEU:HD23	2.50	0.46
1:A:199:ASP:O	1:A:203:ARG:HG2	2.15	0.46
1:A:242:ASP:O	1:A:246:GLY:N	2.48	0.46
1:A:302:VAL:HA	1:A:303:PRO:HD3	1.82	0.46
2:F:505:MET:HG3	2:F:511:PRO:HG2	1.96	0.46
1:B:66:ARG:HH21	1:B:339:THR:CG2	2.28	0.46
2:F:509:PHE:C	2:F:511:PRO:HD3	2.41	0.46
1:A:221:ALA:HA	6:A:1455:HOH:O	2.16	0.46
1:A:223:ARG:HD2	1:A:228:GLU:CA	2.43	0.46
2:F:624:VAL:HG13	2:F:628:PHE:CE1	2.51	0.46
1:B:388:HIS:HD2	1:B:391:LYS:NZ	2.13	0.46
1:B:110:GLN:N	1:B:110:GLN:HE21	2.11	0.46
1:B:25:PRO:HD2	1:B:435:GLU:OE1	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:ASP:OD1	1:B:136:ASP:N	2.49	0.45
1:B:428:ASN:ND2	6:B:1353:HOH:O	2.49	0.45
1:A:388:HIS:HD2	1:A:391:LYS:NZ	2.13	0.45
1:B:223:ARG:HH12	1:B:235:THR:HG22	1.81	0.45
1:B:306:LYS:NZ	6:B:1209:HOH:O	2.47	0.45
1:A:449:LYS:HG3	6:A:1462:HOH:O	2.16	0.45
2:F:557:ASP:OD1	2:F:557:ASP:N	2.48	0.45
2:F:590:ALA:CA	2:F:594:ALA:HB2	2.38	0.45
1:B:400:CYS:HB2	3:B:1001:HEM:NA	2.31	0.45
1:A:91:THR:O	1:A:97:LYS:HD2	2.15	0.45
1:A:141:VAL:HB	1:A:142:PRO:HD3	1.99	0.45
1:B:370:ASP:OD2	1:B:375:ARG:NH1	2.50	0.45
2:F:572:LYS:NZ	6:F:1048:HOH:O	2.49	0.45
1:B:232:ASP:H	1:B:235:THR:CG2	2.24	0.45
1:B:135:ALA:HA	6:B:1048:HOH:O	2.17	0.45
1:B:455:LEU:HB2	1:B:456:GLY:H	1.62	0.45
1:B:124:VAL:HG13	1:B:455:LEU:HD12	1.98	0.45
1:B:339:THR:HG22	1:B:340:VAL:H	1.81	0.45
1:A:268:THR:HB	3:A:1000:HEM:C3B	2.52	0.45
1:B:268:THR:HB	3:B:1001:HEM:C3B	2.52	0.45
1:A:202:LYS:NZ	6:A:1460:HOH:O	2.50	0.44
1:A:422:ASP:OD1	1:A:451:LYS:NZ	2.50	0.44
2:F:483:LEU:HD23	2:F:524:PRO:CG	2.47	0.44
2:F:615:GLU:O	2:F:619:HIS:ND1	2.51	0.44
1:A:68:ASP:OD1	1:A:69:LYS:N	2.49	0.44
1:A:198:TYR:HA	1:A:201:ASN:OD1	2.17	0.44
1:A:309:LYS:CE	6:A:1053:HOH:O	2.66	0.44
1:B:339:THR:HG23	6:B:1100:HOH:O	2.16	0.44
1:A:118[B]:MET:HE2	6:A:1300:HOH:O	2.16	0.44
1:A:336:LYS:NZ	6:A:1106:HOH:O	2.49	0.44
2:F:602:GLU:O	2:F:612:THR:HG21	2.18	0.44
1:A:64:GLU:OE2	1:A:397:GLN:HG2	2.17	0.44
1:A:66:ARG:NH2	1:A:340:VAL:O	2.51	0.44
2:F:486:TYR:O	2:F:516:LEU:N	2.50	0.44
2:F:521:GLY:HA2	2:F:550:TRP:CZ2	2.53	0.44
2:F:548:VAL:HG13	2:F:584:PHE:HE2	1.79	0.44
1:B:101:ASN:ND2	6:B:1238:HOH:O	2.50	0.44
2:F:517:ASP:CB	2:F:546:GLN:HB2	2.41	0.44
2:F:600:ARG:NH2	6:F:1069:HOH:O	2.51	0.44
2:F:620:MET:HE3	2:F:620:MET:HB3	1.91	0.44
1:B:272:LEU:HD13	1:B:322:LEU:CD1	2.38	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:437:LEU:HD12	1:B:437:LEU:HA	1.81	0.44
1:B:102:ILE:HG23	1:B:240:GLY:HA3	2.00	0.43
1:A:186:ASN:O	1:A:189:GLN:HG2	2.18	0.43
1:A:208:ASP:HA	1:A:211:VAL:CG1	2.48	0.43
1:A:354:MET:HB3	1:A:354:MET:HE3	1.62	0.43
1:A:62:CYS:HB3	1:A:395:ASN:ND2	2.33	0.43
1:B:97:LYS:NZ	6:B:1347:HOH:O	2.48	0.43
1:B:253:ASN:O	1:B:257:GLN:HG2	2.19	0.43
1:B:349:LYS:HE3	6:B:1098:HOH:O	2.18	0.43
1:A:289:LYS:HE3	1:A:313:TYR:CZ	2.53	0.43
1:B:141:VAL:HB	1:B:142:PRO:HD3	1.99	0.43
1:B:243:PRO:HG3	6:B:1395:HOH:O	2.18	0.43
1:A:68:ASP:HB3	1:A:334:TYR:CE1	2.54	0.43
1:A:110:GLN:HB3	2:F:610:GLU:OE2	2.19	0.43
1:A:429:TYR:CE2	1:A:431:LEU:HA	2.54	0.43
1:A:458:ILE:O	1:B:456:GLY:HA3	2.19	0.43
2:F:480:THR:HA	2:F:481:PRO:HD3	1.70	0.43
2:F:483:LEU:HD23	2:F:524:PRO:HG3	2.00	0.43
2:F:555:SER:HB3	2:F:592:LYS:HE3	2.00	0.43
2:F:563:ARG:HD3	2:F:627:TYR:OH	2.19	0.43
2:F:620:MET:HE1	2:F:621:TRP:CZ2	2.54	0.42
1:A:323:ARG:HA	1:A:361:HIS:CD2	2.54	0.42
1:A:361:HIS:HE1	1:A:391:LYS:N	2.12	0.42
2:F:523:LEU:HA	2:F:524:PRO:HD3	1.89	0.42
1:A:296[A]:ARG:NH1	6:A:1378:HOH:O	2.50	0.42
2:F:628:PHE:HB2	2:F:629:ASN:H	1.37	0.42
1:B:373:GLU:O	1:B:378:ARG:NH2	2.49	0.42
1:B:306:LYS:CE	1:B:306:LYS:H	2.32	0.42
1:B:380:GLU:HG3	6:B:1216:HOH:O	2.20	0.42
1:B:449:LYS:CE	6:B:1172:HOH:O	2.68	0.42
1:B:32:ILE:HG23	1:B:36:LEU:HD22	2.02	0.42
1:A:126:LEU:HD21	1:A:145:MET:HE1	2.02	0.42
2:F:620:MET:HE1	2:F:621:TRP:CH2	2.54	0.42
1:A:20:LEU:HB3	6:A:1358:HOH:O	2.20	0.42
1:A:128:GLN:HE21	1:A:455:LEU:HD11	1.81	0.42
1:A:282:LYS:C	1:A:284:PRO:HD3	2.44	0.42
1:B:31:LYS:HD2	1:B:31:LYS:HA	1.63	0.42
2:F:563:ARG:HA	2:F:595:GLU:CB	2.50	0.42
2:F:624:VAL:O	2:F:628:PHE:HD1	2.03	0.42
1:B:109:GLN:HB3	1:B:110:GLN:NE2	2.35	0.42
1:B:299:VAL:HG23	1:B:300:ASP:OD1	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:ARG:HD2	1:A:228:GLU:CB	2.50	0.41
1:A:424:GLU:OE1	1:A:449:LYS:NZ	2.52	0.41
1:B:66:ARG:HH21	1:B:339:THR:HG22	1.85	0.41
2:F:485:LEU:HB3	2:F:516:LEU:HD23	2.02	0.41
1:A:24:LYS:N	1:A:25:PRO:HD3	2.36	0.41
1:B:242:ASP:HA	1:B:243:PRO:HD2	1.92	0.41
1:A:110:GLN:HE22	4:A:1006:EDO:H21	1.86	0.41
2:F:602:GLU:O	2:F:612:THR:CG2	2.68	0.41
1:B:39:ILE:HA	1:B:51:TYR:O	2.20	0.41
1:A:210:LYS:HG3	6:A:1318:HOH:O	2.20	0.41
1:B:213:ASN:HA	1:B:255:ARG:HD2	2.03	0.41
1:B:339:THR:HG22	1:B:340:VAL:N	2.36	0.41
1:A:44:ALA:HB1	1:A:45:PRO:HD2	2.02	0.41
1:A:163:ASN:O	1:A:167:ARG:HD2	2.21	0.41
1:A:279:PHE:O	1:A:283:ASN:HB2	2.20	0.41
1:B:192:ASN:OD1	1:B:192:ASN:N	2.52	0.41
1:A:437:LEU:HD12	1:A:437:LEU:HA	1.72	0.41
1:B:458:ILE:HG21	6:B:1194:HOH:O	2.20	0.41
1:A:224:LYS:HD2	6:A:1455:HOH:O	2.20	0.40
1:A:383:SER:HA	2:F:574:TRP:CZ2	2.56	0.40
1:B:113:LYS:HG3	1:B:305:TYR:CD2	2.56	0.40
1:B:237:MET:HE3	1:B:254:ILE:HG23	2.03	0.40
2:F:566:VAL:O	2:F:600:ARG:HA	2.20	0.40
1:B:23:ASP:C	1:B:25:PRO:HD3	2.46	0.40
1:B:32:ILE:HA	1:B:35:GLU:HG2	2.03	0.40
1:A:283:ASN:N	1:A:284:PRO:HD3	2.35	0.40
1:B:322:LEU:HD12	1:B:322:LEU:HA	1.88	0.40
2:F:577:THR:HB	5:F:1002:FMN:O4	2.22	0.40
1:A:400:CYS:HB2	3:A:1000:HEM:C1A	2.56	0.40
1:B:404:GLN:NE2	6:B:1372:HOH:O	2.55	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:1356:HOH:O	6:B:1206:HOH:O[4_566]	2.19	0.01

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	441/458 (96%)	427 (97%)	13 (3%)	1 (0%)	43	40
1	B	436/458 (95%)	416 (95%)	18 (4%)	2 (0%)	24	17
2	F	150/191 (78%)	123 (82%)	22 (15%)	5 (3%)	3	1
All	All	1027/1107 (93%)	966 (94%)	53 (5%)	8 (1%)	16	8

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	558	GLU
2	F	575	ALA
1	B	228	GLU
1	B	229	GLN
2	F	596	ASN
2	F	552	ASP
1	A	22	THR
2	F	625	ALA

5.3.2 Protein sidechains [i](#)

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	386/400 (96%)	363 (94%)	23 (6%)	17	10
1	B	381/400 (95%)	352 (92%)	29 (8%)	12	6
2	F	117/153 (76%)	103 (88%)	14 (12%)	5	1

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	884/953 (93%)	818 (92%)	66 (8%)	12 7

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	31	LYS
1	A	45	PRO
1	A	94	LYS
1	A	110	GLN
1	A	137	GLU
1	A	167	ARG
1	A	211	VAL
1	A	226	SER
1	A	241	LYS
1	A	255	ARG
1	A	263	ILE
1	A	306	LYS
1	A	309	LYS
1	A	338	ASP
1	A	348	GLU
1	A	354	MET
1	A	359	GLN
1	A	375	ARG
1	A	437	LEU
1	A	440	LYS
1	A	442	GLU
1	A	455	LEU
2	F	479	ASN
2	F	480	THR
2	F	525	ARG
2	F	553	GLN
2	F	557	ASP
2	F	560	LYS
2	F	572	LYS
2	F	577	THR
2	F	589	LEU
2	F	606	SER
2	F	612	THR
2	F	622	SER
2	F	624	VAL
2	F	629	ASN

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Mol	Chain	Res	Type
1	B	21	ASN
1	B	23	ASP
1	B	31	LYS
1	B	36	LEU
1	B	41	LYS
1	B	52	LEU
1	B	71	LEU
1	B	76	LYS
1	B	110	GLN
1	B	113	LYS
1	B	148	LEU
1	B	190	ARG
1	B	215	LEU
1	B	218	LYS
1	B	224	LYS
1	B	235	THR
1	B	270	SER
1	B	302	VAL
1	B	306	LYS
1	B	309	LYS
1	B	312	LYS
1	B	322	LEU
1	B	359	GLN
1	B	380	GLU
1	B	383	SER
1	B	385	ILE
1	B	437	LEU
1	B	455	LEU
1	B	458	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	ASN
1	A	73	GLN
1	A	95	ASN
1	A	109	GLN
1	A	110	GLN
1	A	116	HIS
1	A	128	GLN
1	A	159	ASN
1	A	236	HIS

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Mol	Chain	Res	Type
1	A	319	ASN
1	A	359	GLN
1	A	361	HIS
1	A	381	ASN
1	A	388	HIS
1	A	395	ASN
1	A	426	HIS
2	F	479	ASN
2	F	519	HIS
2	F	546	GLN
2	F	573	ASN
2	F	629	ASN
1	B	27	GLN
1	B	70	ASN
1	B	73	GLN
1	B	100	HIS
1	B	109	GLN
1	B	110	GLN
1	B	116	HIS
1	B	159	ASN
1	B	171	HIS
1	B	189	GLN
1	B	204	GLN
1	B	310	GLN
1	B	361	HIS
1	B	387	GLN
1	B	388	HIS
1	B	395	ASN
1	B	397	GLN
1	B	428	ASN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	B	1007	-	3,3,3	0.47	0	2,2,2	0.34	0
3	HEM	A	1000	1,4	50,50,50	1.01	2 (4%)	67,82,82	0.99	3 (4%)
4	EDO	B	1005	-	3,3,3	0.56	0	2,2,2	0.46	0
3	HEM	B	1001	1,4	50,50,50	0.99	2 (4%)	67,82,82	0.97	2 (2%)
4	EDO	A	1003	3	3,3,3	0.47	0	2,2,2	0.39	0
4	EDO	B	1004	3	3,3,3	0.54	0	2,2,2	0.50	0
5	FMN	F	1002	-	33,33,33	1.39	4 (12%)	48,50,50	1.97	13 (27%)
4	EDO	A	1006	-	3,3,3	0.48	0	2,2,2	0.49	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 Torsions and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	1007	-	-	1/1/1/1	-
3	HEM	A	1000	1,4	-	2/14/54/54	-
4	EDO	B	1005	-	-	0/1/1/1	-
3	HEM	B	1001	1,4	-	2/14/54/54	-
4	EDO	A	1003	3	-	1/1/1/1	-
4	EDO	B	1004	3	-	1/1/1/1	-
5	FMN	F	1002	-	-	5/18/18/18	0/3/3/3
4	EDO	A	1006	-	-	0/1/1/1	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	1002	FMN	P-O5'	-4.33	1.46	1.60
3	B	1001	HEM	CAC-C3C	-3.27	1.38	1.47
5	F	1002	FMN	C5A-N5	-3.24	1.33	1.39
3	A	1000	HEM	CAC-C3C	-3.10	1.39	1.47
3	A	1000	HEM	CAB-C3B	-3.06	1.39	1.47
3	B	1001	HEM	CAB-C3B	-2.96	1.39	1.47
5	F	1002	FMN	O5'-C5'	2.63	1.54	1.44
5	F	1002	FMN	C10-N10	2.27	1.42	1.37

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	1002	FMN	O4'-C4'-C3'	-5.60	96.15	109.25
5	F	1002	FMN	C4-N3-C2	-4.25	118.09	125.64
5	F	1002	FMN	C4'-C3'-C2'	-4.12	106.71	113.57
5	F	1002	FMN	C5A-N5-C4A	3.59	123.90	118.09
5	F	1002	FMN	O5'-C5'-C4'	-3.30	100.54	109.36
5	F	1002	FMN	C4-C4A-N5	3.21	122.64	118.21
5	F	1002	FMN	C4A-C4-N3	2.80	120.38	113.25
3	A	1000	HEM	CHD-C1D-C2D	-2.75	120.69	125.03
5	F	1002	FMN	O2P-P-O5'	2.73	113.79	106.67
5	F	1002	FMN	O3P-P-O5'	2.72	113.77	106.67
5	F	1002	FMN	C9A-C5A-N5	-2.55	119.75	122.45
5	F	1002	FMN	C10-C4A-N5	-2.47	119.77	124.81
5	F	1002	FMN	O5'-P-O1P	2.45	113.06	106.44
5	F	1002	FMN	C6-C5A-N5	2.28	122.23	118.44
3	A	1000	HEM	C3B-C4B-NB	2.26	111.09	109.47
3	A	1000	HEM	CMC-C2C-C1C	2.12	128.47	124.73
3	B	1001	HEM	O2D-CGD-CBD	2.08	120.58	114.00
3	B	1001	HEM	CHB-C4A-C3A	-2.05	121.48	127.43

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1007	EDO	O1-C1-C2-O2
5	F	1002	FMN	O3'-C3'-C4'-C5'
5	F	1002	FMN	C1'-C2'-C3'-O3'
5	F	1002	FMN	C2'-C3'-C4'-C5'
5	F	1002	FMN	O2'-C2'-C3'-O3'
4	A	1003	EDO	O1-C1-C2-O2
4	B	1004	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
5	F	1002	FMN	O3'-C3'-C4'-O4'
3	A	1000	HEM	CAD-CBD-CGD-O2D
3	B	1001	HEM	CAD-CBD-CGD-O1D
3	A	1000	HEM	CAD-CBD-CGD-O1D
3	B	1001	HEM	CAD-CBD-CGD-O2D

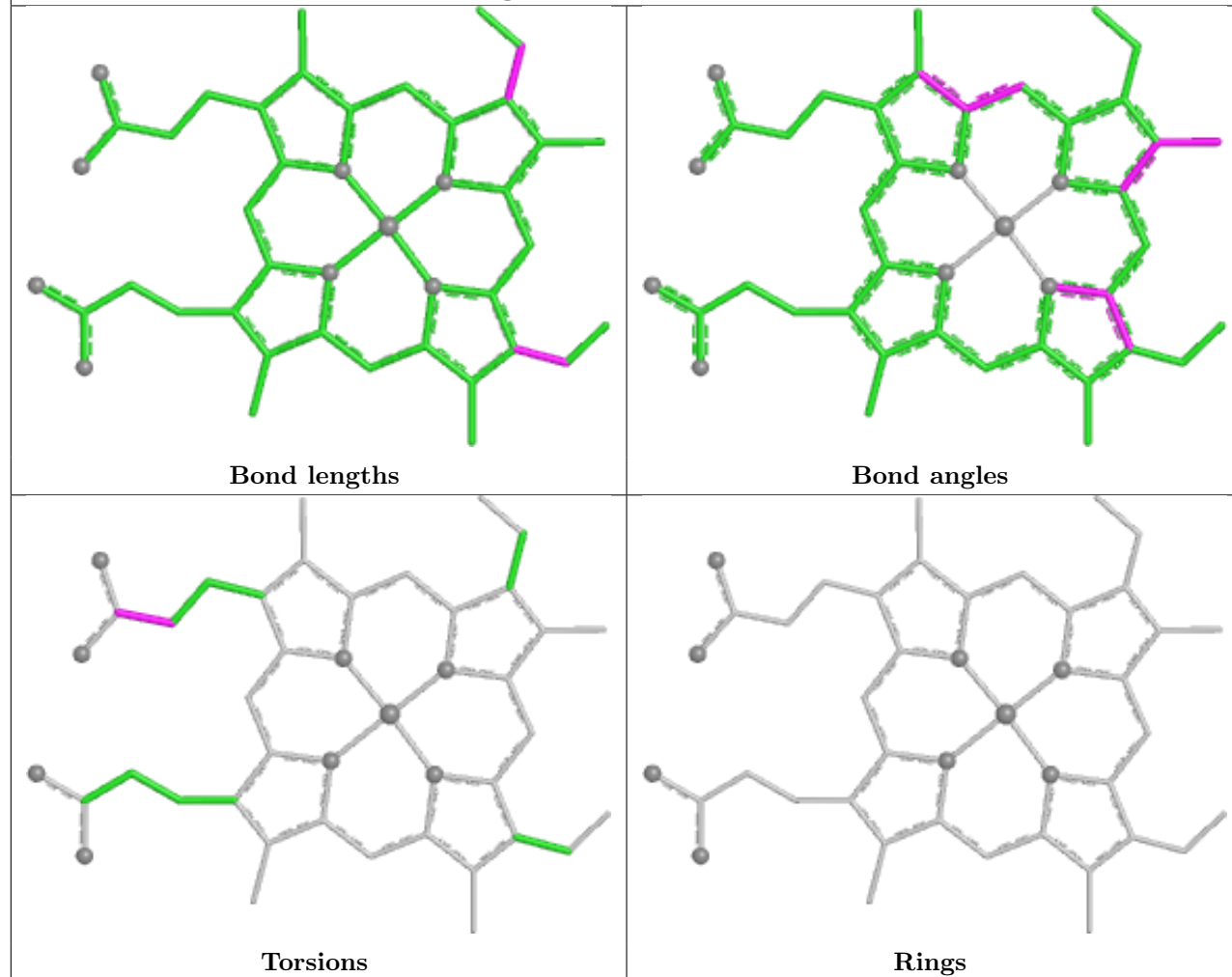
There are no ring outliers.

5 monomers are involved in 10 short contacts:

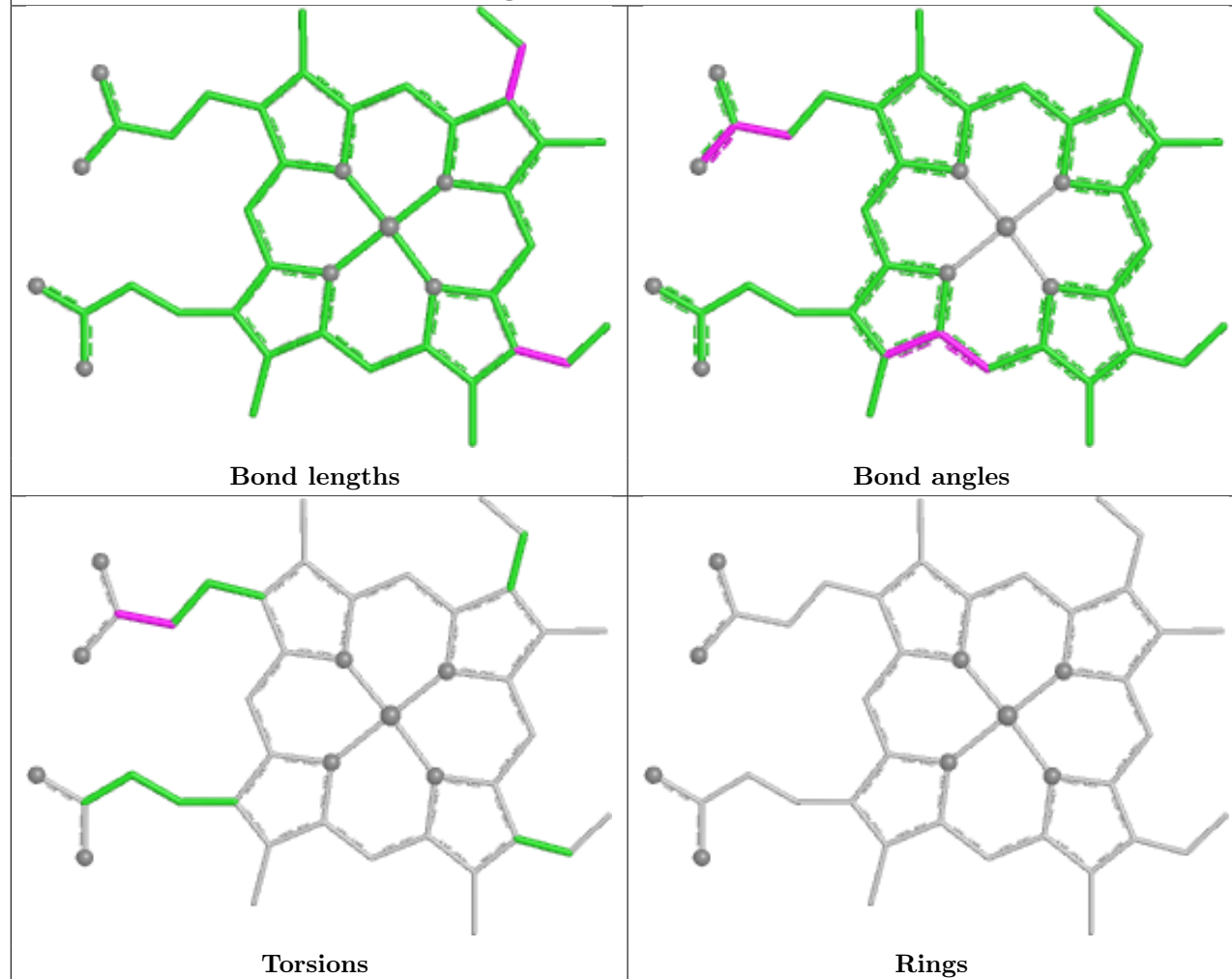
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1007	EDO	1	0
3	A	1000	HEM	4	0
3	B	1001	HEM	3	0
5	F	1002	FMN	1	0
4	A	1006	EDO	1	0

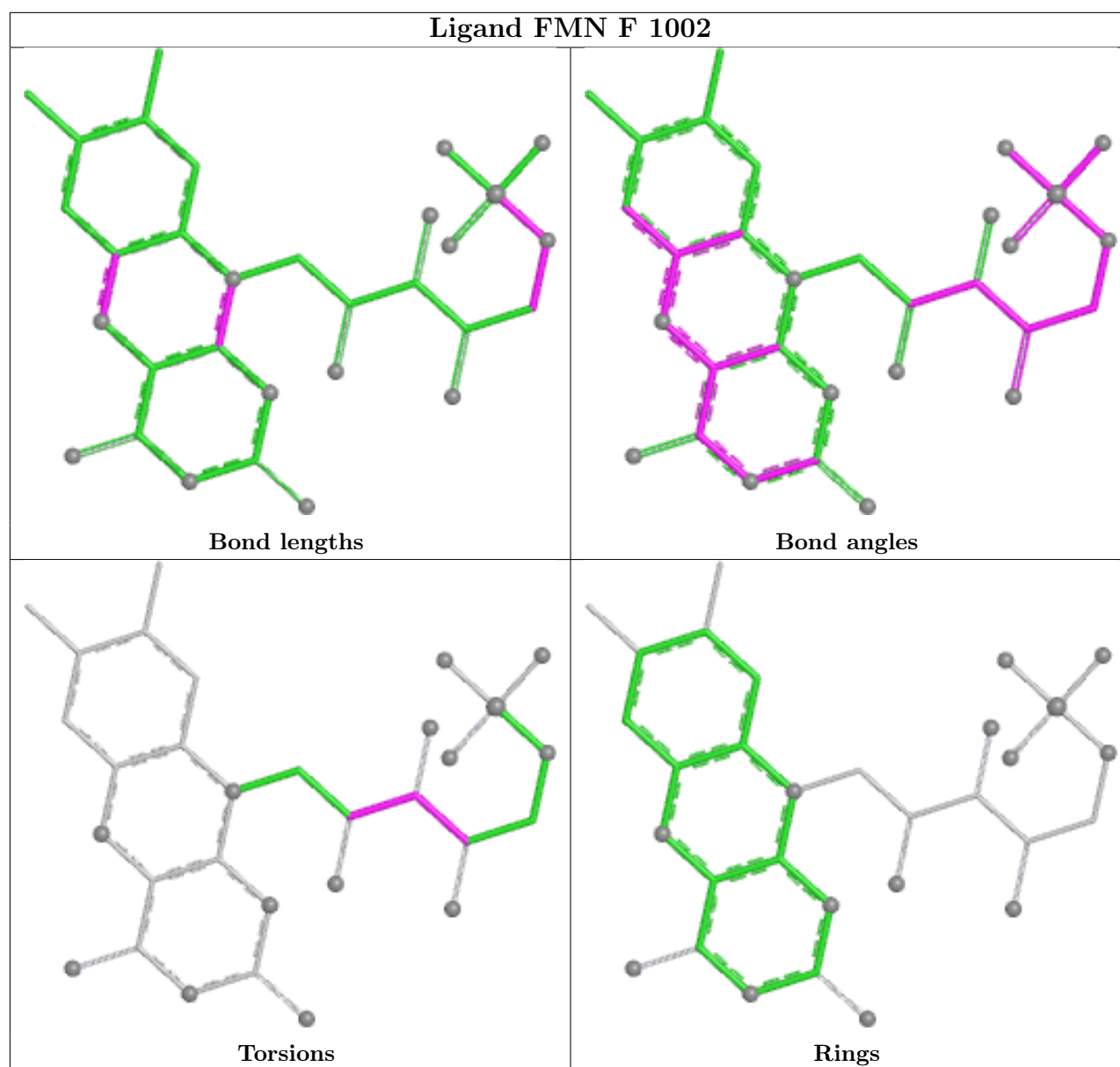
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

Ligand HEM A 1000



Ligand HEM B 1001





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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.