



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 12:33 PM UTC

PDB ID : 2BCN / pdb_00002bcn
Title : Solvent isotope effects on interfacial protein electron transfer between cytochrome c and cytochrome c peroxidase
Authors : Kang, S.A.; Crane, B.R.
Deposited on : 2005-10-19
Resolution : 1.70 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) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

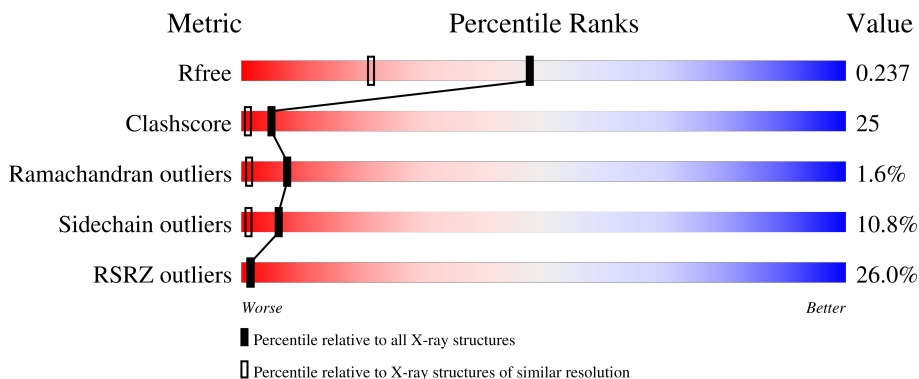
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	180053	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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 $\leq 5\%$. 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	296	
1	C	296	
2	B	108	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ZNH	A	295	X	-	-	-
3	ZNH	C	295	X	-	-	-

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 6265 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called cytochrome c peroxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	294	2371	1514	395	456	6	0	0	0
1	C	294	2371	1514	395	456	6	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	cloning artifact	GB 171177
A	0	ILE	-	cloning artifact	GB 171177
C	-1	MET	-	cloning artifact	GB 171177
C	0	ILE	-	cloning artifact	GB 171177

- Molecule 2 is a protein called Cytochrome c iso-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	108	847	534	151	158	4	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	107	SER	CYS	engineered mutation	UNP P00044

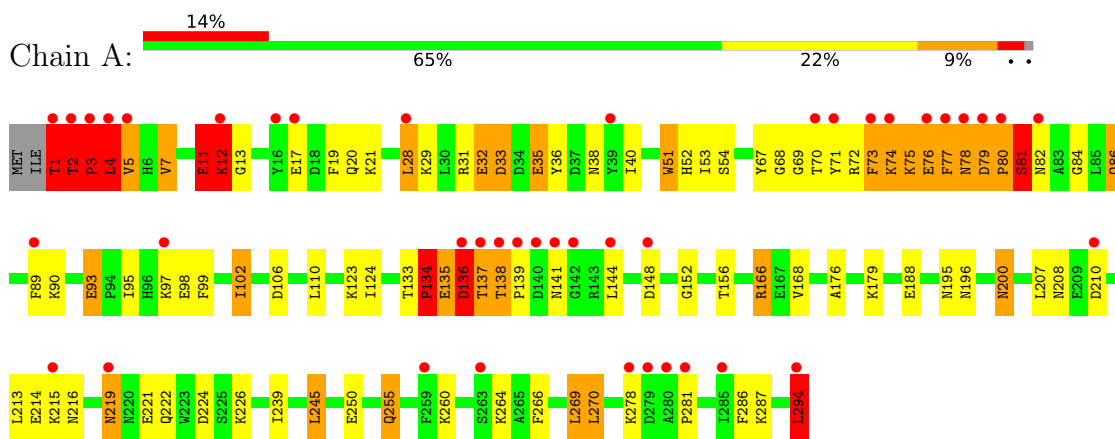
- Molecule 3 is PROTOPORPHYRIN IX CONTAINING ZN (CCD ID: ZNH) (formula: C₃₄H₃₂N₄O₄Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	286	Total 286	O 286	0	0
5	B	27	Total 27	O 27	0	0
5	C	234	Total 234	O 234	0	0

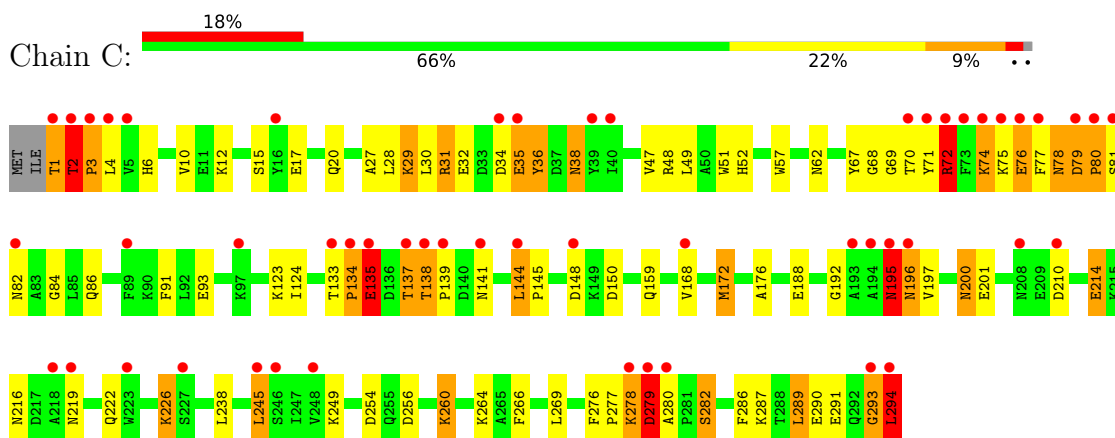
3 Residue-property plots [i](#)

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 and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

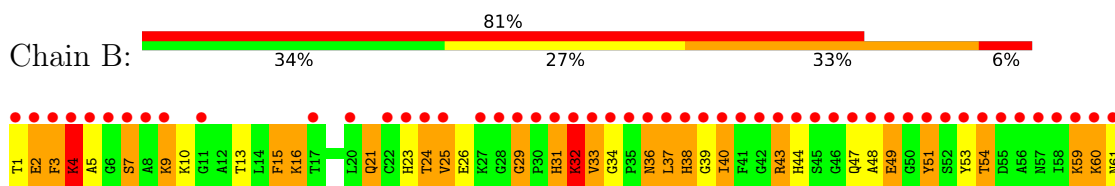
- Molecule 1: cytochrome c peroxidase



- Molecule 1: cytochrome c peroxidase



- Molecule 2: Cytochrome c iso-1



V62	L63	W64	D65	E66	N67	M68	R69	S70	E71	Y72	L73	T74	N75	P76	K77	K78	Y79	I80	F81	G82	T83	K84	M85	A86	F87	G88	G89	L90	K91	K92	E93	K94	D95	R96	N97	D98	L99	I100	T101	Y102	L103	K104	K105	A106	S107	E108
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	44.53Å 47.09Å 84.29Å 102.47° 98.74° 102.16°	Depositor
Resolution (Å)	19.61 – 1.70 19.61 – 1.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.61-1.70) 93.1 (19.61-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 1.70Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.219 , 0.230 0.225 , 0.237	Depositor DCC
R_{free} test set	3335 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	19.1	Xtrriage
Anisotropy	0.068	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6265	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.46% 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: HEC, ZNH

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	1.65	13/2438 (0.5%)	1.66	59/3302 (1.8%)
1	C	1.17	5/2438 (0.2%)	1.79	60/3302 (1.8%)
2	B	1.02	4/865 (0.5%)	1.82	50/1156 (4.3%)
All	All	1.38	22/5741 (0.4%)	1.74	169/7760 (2.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	8
1	C	1	6
All	All	1	14

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	79	ASP	C-N	36.68	1.74	1.34
1	A	1	THR	C-N	-33.12	0.65	1.33
1	A	11	GLU	C-N	28.84	1.74	1.33
1	A	2	THR	C-N	27.33	1.69	1.33
1	C	79	ASP	C-N	25.20	1.66	1.33
1	C	277	PRO	C-N	24.41	1.66	1.33
1	A	134	PRO	C-N	19.27	1.63	1.33
1	A	12	LYS	C-N	16.87	1.59	1.33
1	A	3	PRO	C-N	16.51	1.56	1.33
2	B	26	GLU	C-N	15.25	1.55	1.33
1	A	135	GLU	C-N	-13.67	1.14	1.33
1	A	73	PHE	C-N	-12.48	1.19	1.33
2	B	25	VAL	C-N	10.43	1.47	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	9	LYS	C-N	-9.69	1.20	1.33
1	A	74	LYS	C-N	9.52	1.45	1.33
1	A	81	SER	C-N	-8.98	1.21	1.33
1	A	4	LEU	C-N	-8.73	1.23	1.33
1	C	2	THR	C-N	8.50	1.44	1.33
2	B	10	LYS	C-N	7.58	1.43	1.33
1	A	54	SER	C-N	-6.61	1.25	1.33
1	C	3	PRO	N-CD	6.52	1.56	1.47
1	C	80	PRO	C-N	-5.81	1.26	1.33

All (169) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	279	ASP	O-C-N	-47.62	59.26	122.59
1	A	79	ASP	CA-C-N	-25.54	93.31	119.87
1	A	79	ASP	C-N-CA	-25.54	93.31	119.87
1	C	294	LEU	CB-CA-C	-19.89	72.31	110.10
1	A	1	THR	O-C-N	-17.41	95.15	123.00
1	C	79	ASP	CA-C-N	-15.63	102.08	119.28
1	C	79	ASP	C-N-CA	-15.63	102.08	119.28
1	A	2	THR	O-C-N	-14.14	105.06	121.32
1	A	4	LEU	O-C-N	-14.11	103.82	122.59
1	C	279	ASP	CB-CA-C	12.29	134.88	110.42
1	A	1	THR	CA-C-N	11.39	149.59	121.80
1	A	1	THR	C-N-CA	11.39	149.59	121.80
2	B	77	LYS	CB-CA-C	11.05	130.04	110.37
1	A	134	PRO	CA-C-N	-10.91	102.45	121.66
1	A	134	PRO	C-N-CA	-10.91	102.45	121.66
1	A	12	LYS	CA-C-N	-10.66	102.76	122.05
1	A	12	LYS	C-N-CA	-10.66	102.76	122.05
2	B	69	MET	CB-CA-C	10.61	132.02	110.38
1	C	278	LYS	CA-C-N	-10.32	101.83	121.54
1	C	278	LYS	C-N-CA	-10.32	101.83	121.54
1	C	148	ASP	CA-CB-CG	10.26	122.86	112.60
2	B	24	THR	CB-CA-C	10.12	127.01	109.80
1	A	33	ASP	CB-CA-C	-10.06	97.70	111.88
1	A	81	SER	CB-CA-C	10.04	129.94	110.57
1	C	34	ASP	CB-CA-C	10.03	129.91	109.95
2	B	107	SER	CB-CA-C	10.02	130.35	110.42
1	C	293	GLY	O-C-N	-9.97	111.43	122.50
1	A	81	SER	CA-C-N	9.60	137.16	122.37
1	A	81	SER	C-N-CA	9.60	137.16	122.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	73	PHE	CA-C-N	9.60	136.45	121.19
1	A	73	PHE	C-N-CA	9.60	136.45	121.19
2	B	104	LYS	CB-CA-C	9.57	129.90	110.38
1	A	219	ASN	CB-CA-C	9.52	129.06	111.06
1	A	79	ASP	CB-CA-C	9.38	124.52	109.51
1	C	148	ASP	CA-C-N	9.17	138.53	122.67
1	C	148	ASP	C-N-CA	9.17	138.53	122.67
1	A	11	GLU	CA-C-N	-9.12	104.13	121.54
1	A	11	GLU	C-N-CA	-9.12	104.13	121.54
1	C	195	ASN	N-CA-C	9.08	122.16	111.71
2	B	25	VAL	N-CA-C	8.48	124.69	113.07
2	B	59	LYS	CB-CA-C	8.42	127.35	110.17
1	A	74	LYS	O-C-N	8.40	132.37	122.20
1	A	5	VAL	CB-CA-C	8.36	121.87	111.25
1	C	74	LYS	N-CA-C	8.36	121.44	111.33
1	C	279	ASP	CA-C-N	8.13	141.65	121.80
1	C	279	ASP	C-N-CA	8.13	141.65	121.80
1	C	277	PRO	O-C-N	-8.03	113.46	123.10
2	B	33	VAL	N-CA-C	-7.97	103.13	111.58
2	B	63	LEU	CB-CA-C	7.89	122.10	110.26
1	A	2	THR	CA-C-N	7.80	127.56	119.76
1	A	2	THR	C-N-CA	7.80	127.56	119.76
2	B	36	ASN	CB-CA-C	7.64	122.82	110.29
1	A	210	ASP	CB-CA-C	7.60	122.95	110.78
1	A	7	VAL	CB-CA-C	7.60	121.03	111.15
2	B	31	HIS	CB-CA-C	7.58	124.05	111.31
1	A	73	PHE	O-C-N	-7.50	114.66	123.06
1	A	13	GLY	O-C-N	7.50	131.95	122.43
1	A	38	ASN	CB-CA-C	7.43	122.44	111.89
1	C	2	THR	CA-CB-OG1	7.40	120.70	109.60
1	C	277	PRO	CA-C-N	-7.34	110.44	120.28
1	C	277	PRO	C-N-CA	-7.34	110.44	120.28
1	C	150	ASP	CA-CB-CG	7.30	119.90	112.60
1	C	3	PRO	CA-N-CD	-7.25	101.85	112.00
1	C	137	THR	N-CA-C	-7.24	104.42	113.18
2	B	7	SER	CB-CA-C	7.23	121.66	109.80
2	B	102	TYR	CB-CA-C	7.21	122.29	110.90
1	C	138	THR	CB-CA-C	7.20	121.50	109.69
2	B	32	LYS	CB-CA-C	7.18	124.72	110.42
2	B	38	HIS	CB-CA-C	7.11	121.42	109.84
1	A	195	ASN	CB-CA-C	-7.09	96.71	110.46
1	A	137	THR	CB-CA-C	-7.03	96.43	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	29	GLY	CA-C-N	7.00	126.99	119.78
2	B	29	GLY	C-N-CA	7.00	126.99	119.78
1	A	136	ASP	O-C-N	-6.99	113.29	122.59
1	C	2	THR	N-CA-CB	6.98	123.75	110.49
1	A	78	ASN	CB-CA-C	6.96	121.85	111.06
1	A	213	LEU	CB-CA-C	6.91	122.30	110.77
1	C	74	LYS	CB-CA-C	-6.86	99.01	110.68
2	B	83	THR	CB-CA-C	6.85	122.07	109.46
2	B	87	PHE	CB-CA-C	6.83	123.24	109.38
2	B	2	GLU	CB-CA-C	6.79	124.23	110.38
2	B	49	GLU	N-CA-C	6.77	119.56	110.35
2	B	100	ILE	CB-CA-C	6.77	121.56	112.22
2	B	9	LYS	CB-CA-C	-6.74	96.63	110.38
1	C	210	ASP	CB-CA-C	6.66	122.62	111.30
1	C	290	GLU	CB-CA-C	6.66	121.84	110.79
2	B	108	GLU	CB-CA-C	-6.66	97.45	110.10
1	C	75	LYS	N-CA-C	6.63	118.17	111.07
1	C	38	ASN	CB-CA-C	6.56	121.04	111.73
1	C	29	LYS	CB-CA-C	6.50	123.15	110.67
1	A	138	THR	CB-CA-C	6.48	120.51	109.82
1	C	78	ASN	CB-CA-C	6.47	121.61	110.94
1	C	3	PRO	N-CA-CB	6.45	109.05	103.31
2	B	72	TYR	CB-CA-C	6.37	121.49	110.79
1	C	197	VAL	CB-CA-C	-6.36	101.04	110.33
1	C	72	ARG	CB-CA-C	6.35	122.38	110.63
1	C	282	SER	CB-CA-C	-6.32	99.10	109.27
2	B	84	LYS	CB-CA-C	6.26	121.64	110.81
1	C	214	GLU	CB-CA-C	-6.22	95.75	109.56
2	B	4	LYS	CB-CA-C	6.21	122.79	110.42
2	B	9	LYS	CA-C-N	6.21	131.21	120.58
2	B	9	LYS	C-N-CA	6.21	131.21	120.58
1	A	76	GLU	CB-CA-C	-6.21	97.72	110.38
1	C	77	PHE	CB-CA-C	6.18	121.36	110.85
1	C	134	PRO	N-CA-C	6.15	120.12	111.33
1	A	294	LEU	CB-CA-C	-6.12	98.47	110.10
2	B	64	TRP	CB-CA-C	-6.11	98.26	110.42
1	A	77	PHE	CB-CA-C	6.09	121.21	110.85
2	B	54	THR	CB-CA-C	6.09	120.49	109.83
2	B	68	ASN	CB-CA-C	6.08	122.05	109.95
1	A	195	ASN	N-CA-C	6.07	119.52	111.75
2	B	10	LYS	CB-CA-C	6.04	121.75	110.70
2	B	103	LEU	CB-CA-C	-6.02	97.74	110.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	THR	CB-CA-C	6.01	122.33	109.10
2	B	74	THR	CB-CA-C	6.01	122.64	110.38
2	B	43	ARG	CB-CA-C	-6.00	97.51	109.33
2	B	15	PHE	CB-CA-C	6.00	120.87	110.79
1	C	31	ARG	CB-CA-C	-5.98	98.17	110.38
1	A	176	ALA	N-CA-C	-5.97	105.25	112.54
1	C	135	GLU	CB-CA-C	-5.96	98.91	110.46
1	A	67	TYR	N-CA-C	5.92	118.24	111.02
2	B	21	GLN	CB-CA-C	-5.91	98.12	110.17
1	C	17	GLU	CB-CA-C	5.88	120.67	110.68
1	C	76	GLU	CB-CA-C	5.87	121.47	109.55
1	A	1	THR	CB-CA-C	5.85	121.97	109.10
1	C	172	MET	CB-CA-C	-5.82	98.30	110.17
2	B	70	SER	CB-CA-C	5.80	121.81	110.67
1	C	1	THR	O-C-N	-5.80	113.72	123.00
1	C	2	THR	O-C-N	5.80	127.11	121.66
1	C	135	GLU	N-CA-C	5.78	119.15	111.75
2	B	67	ASN	CB-CA-C	5.75	120.98	110.11
2	B	105	LYS	CB-CA-C	-5.71	98.19	109.67
1	A	98	GLU	CB-CA-C	-5.68	98.26	109.67
1	C	81	SER	CB-CA-C	5.67	122.29	110.32
1	C	159	GLN	N-CA-C	-5.60	105.32	111.82
1	A	255	GLN	CB-CA-C	5.59	121.41	110.67
2	B	37	LEU	CB-CA-C	5.53	120.22	110.37
1	A	79	ASP	C-N-CD	5.52	147.63	125.00
2	B	10	LYS	O-C-N	5.48	129.36	122.23
2	B	97	ASN	CB-CA-C	5.47	121.18	110.67
1	A	110	LEU	CB-CA-C	5.47	121.17	110.67
1	A	40	ILE	N-CA-C	-5.43	105.08	111.00
1	C	176	ALA	N-CA-C	-5.42	105.93	112.54
1	C	3	PRO	N-CA-C	5.40	119.57	111.14
1	C	70	THR	N-CA-C	5.35	120.76	113.37
1	C	2	THR	CA-CB-CG2	5.35	119.59	110.50
1	C	226	LYS	CB-CA-C	5.33	121.45	110.31
2	B	51	TYR	CB-CA-C	-5.32	98.26	109.38
1	A	28	LEU	CB-CA-C	5.32	119.89	110.85
2	B	24	THR	O-C-N	5.25	129.90	123.44
1	A	35	GLU	CB-CA-C	-5.25	99.03	109.99
1	C	219	ASN	CB-CA-C	5.23	120.94	111.06
1	A	4	LEU	CA-C-N	5.22	129.01	122.43
1	A	4	LEU	C-N-CA	5.22	129.01	122.43
1	C	47	VAL	CB-CA-C	-5.22	103.66	112.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	79	ASP	O-C-N	5.20	125.73	121.31
1	C	67	TYR	N-CA-C	5.19	119.89	111.37
1	A	135	GLU	CA-C-N	5.18	131.44	121.54
1	A	135	GLU	C-N-CA	5.18	131.44	121.54
1	A	208	ASN	N-CA-C	5.16	119.58	113.28
1	C	1	THR	CB-CA-C	5.15	120.42	109.10
1	A	136	ASP	CB-CA-C	5.15	120.66	110.42
2	B	26	GLU	O-C-N	-5.14	116.38	122.65
1	C	35	GLU	CB-CA-C	5.14	119.58	110.85
2	B	75	ASN	CA-C-N	5.12	124.91	119.28
2	B	75	ASN	C-N-CA	5.12	124.91	119.28
1	C	36	TYR	N-CA-C	5.08	117.84	110.42
1	A	68	GLY	N-CA-C	5.03	120.25	114.16
1	A	75	LYS	N-CA-C	5.01	117.12	111.11

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	C	279	ASP	CA

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	THR	Mainchain,Peptide
1	A	136	ASP	Mainchain
1	A	166	ARG	Sidechain
1	A	3	PRO	Mainchain,Peptide
1	A	80	PRO	Mainchain
1	A	81	SER	Mainchain
1	C	1	THR	Mainchain,Peptide
1	C	2	THR	Mainchain
1	C	279	ASP	Mainchain
1	C	293	GLY	Peptide
1	C	72	ARG	Sidechain

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	2371	0	2247	121	4
1	C	2371	0	2252	72	4
2	B	847	0	849	90	3
3	A	43	0	30	0	0
3	C	43	0	30	0	0
4	B	43	0	30	5	0
5	A	286	0	0	66	2
5	B	27	0	0	12	1
5	C	234	0	0	37	2
All	All	6265	0	5438	280	9

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 25.

All (280) 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:2:THR:C	1:A:3:PRO:N	1.69	1.49
1:A:11:GLU:C	1:A:12:LYS:N	1.74	1.42
1:A:79:ASP:C	1:A:80:PRO:N	1.74	1.41
2:B:100:ILE:O	2:B:104:LYS:HG3	1.22	1.34
1:C:76:GLU:HB3	5:C:1082:HOH:O	1.17	1.34
1:A:12:LYS:HE3	5:A:1377:HOH:O	1.16	1.30
1:C:141:ASN:HB2	5:C:1206:HOH:O	1.14	1.28
2:B:60:LYS:HE3	5:B:1525:HOH:O	1.34	1.27
1:A:12:LYS:CE	5:A:1377:HOH:O	1.69	1.22
1:C:260:LYS:HA	1:C:260:LYS:CE	1.69	1.22
2:B:40:ILE:H	2:B:40:ILE:CD1	1.51	1.19
1:C:260:LYS:HE3	1:C:260:LYS:CA	1.70	1.18
1:A:79:ASP:HA	5:A:1409:HOH:O	1.43	1.16
1:A:12:LYS:NZ	5:A:1378:HOH:O	1.69	1.14
1:A:74:LYS:HE3	5:A:1403:HOH:O	1.47	1.13
1:A:12:LYS:HG2	5:A:1377:HOH:O	1.49	1.12
1:C:278:LYS:O	1:C:279:ASP:C	1.78	1.12
2:B:40:ILE:HD12	2:B:40:ILE:N	1.55	1.10
1:A:141:ASN:ND2	5:A:1409:HOH:O	1.87	1.07
2:B:61:ASN:OD1	5:B:1510:HOH:O	1.71	1.06
2:B:100:ILE:O	2:B:104:LYS:CG	2.03	1.06
1:A:77:PHE:CD2	5:A:1408:HOH:O	2.10	1.05
2:B:4:LYS:NZ	2:B:4:LYS:HB2	1.71	1.04
1:C:139:PRO:HD2	5:C:1529:HOH:O	1.56	1.03
2:B:74:THR:HB	2:B:96:ARG:HE	1.23	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:101:THR:HA	2:B:104:LYS:HD3	1.40	1.01
1:A:77:PHE:CA	5:A:1405:HOH:O	2.10	1.00
1:A:79:ASP:C	1:A:80:PRO:CA	2.35	0.99
1:C:201:GLU:O	5:C:1524:HOH:O	1.78	0.99
1:A:156:THR:HG22	5:A:1085:HOH:O	1.60	0.99
1:C:82:ASN:ND2	5:C:1455:HOH:O	1.95	0.99
2:B:44:HIS:NE2	5:B:1501:HOH:O	1.96	0.96
2:B:62:VAL:HA	5:B:1535:HOH:O	1.63	0.96
2:B:105:LYS:HE3	2:B:105:LYS:HA	1.45	0.96
2:B:49:GLU:OE2	5:B:1185:HOH:O	1.82	0.96
1:A:77:PHE:HA	5:A:1405:HOH:O	1.64	0.96
1:C:141:ASN:CB	5:C:1206:HOH:O	1.82	0.96
1:A:221:GLU:CD	5:A:1180:HOH:O	2.07	0.95
1:A:139:PRO:HD2	5:A:1108:HOH:O	1.68	0.92
1:C:69:GLY:HA2	5:C:1457:HOH:O	1.70	0.90
2:B:33:VAL:O	2:B:51:TYR:OH	1.89	0.89
2:B:38:HIS:CD2	2:B:107:SER:HA	2.10	0.87
2:B:4:LYS:NZ	2:B:4:LYS:CB	2.38	0.86
2:B:4:LYS:HB2	2:B:4:LYS:HZ3	1.39	0.86
2:B:40:ILE:H	2:B:40:ILE:HD12	0.69	0.84
1:A:77:PHE:CG	5:A:1408:HOH:O	2.28	0.84
2:B:65:ASP:OD2	5:B:1536:HOH:O	1.96	0.84
1:A:70:THR:O	5:A:1511:HOH:O	1.95	0.83
1:A:76:GLU:C	5:A:1513:HOH:O	2.23	0.82
1:A:75:LYS:O	5:A:1513:HOH:O	1.96	0.82
1:A:78:ASN:N	5:A:1513:HOH:O	2.12	0.81
2:B:105:LYS:HE3	2:B:105:LYS:CA	2.11	0.81
2:B:3:PHE:HE1	2:B:97:ASN:CB	1.95	0.80
1:C:10:VAL:O	5:C:1445:HOH:O	1.99	0.80
1:C:20:GLN:HE22	1:C:287:LYS:H	1.29	0.80
1:A:52:HIS:NE2	5:A:1143:HOH:O	2.14	0.80
1:A:11:GLU:C	1:A:12:LYS:CA	2.54	0.80
1:C:134:PRO:HD2	5:C:1471:HOH:O	1.82	0.79
2:B:64:TRP:CZ2	4:B:109:HEC:HAA1	2.18	0.79
1:A:135:GLU:OE2	5:A:1250:HOH:O	2.01	0.78
2:B:77:LYS:HD2	2:B:82:GLY:H	1.48	0.78
1:A:80:PRO:HB3	5:A:1514:HOH:O	1.82	0.78
1:A:36:TYR:OH	5:A:1241:HOH:O	2.02	0.78
1:A:77:PHE:N	5:A:1405:HOH:O	2.15	0.77
2:B:74:THR:HB	2:B:96:ARG:NE	1.99	0.77
2:B:4:LYS:HB2	2:B:4:LYS:HZ2	1.50	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:ASN:OD1	5:C:1506:HOH:O	2.04	0.75
1:A:73:PHE:O	5:A:1511:HOH:O	2.05	0.74
1:A:188:GLU:H	1:A:222:GLN:HE22	1.35	0.74
2:B:96:ARG:HG3	2:B:96:ARG:HH11	1.51	0.74
2:B:96:ARG:HG3	2:B:96:ARG:NH1	2.03	0.74
2:B:3:PHE:CE1	2:B:97:ASN:CB	2.71	0.73
2:B:4:LYS:CB	2:B:4:LYS:HZ3	2.00	0.73
1:C:137:THR:HG23	5:C:1471:HOH:O	1.88	0.73
1:A:71:TYR:CD1	5:A:1405:HOH:O	2.41	0.73
1:A:137:THR:N	5:A:1417:HOH:O	2.21	0.73
1:A:216:ASN:HD22	1:A:222:GLN:HE21	1.37	0.72
2:B:105:LYS:HA	2:B:105:LYS:CE	2.16	0.72
1:C:144:LEU:HD12	1:C:145:PRO:HD2	1.71	0.72
1:A:264:LYS:CD	5:A:1225:HOH:O	2.38	0.71
1:A:166:ARG:HH22	1:A:250:GLU:CD	1.98	0.70
1:A:139:PRO:CD	5:A:1108:HOH:O	2.33	0.70
1:A:12:LYS:CG	5:A:1377:HOH:O	2.13	0.69
1:C:27:ALA:HB1	1:C:289:LEU:HD22	1.73	0.69
1:C:134:PRO:CD	5:C:1471:HOH:O	2.37	0.69
1:A:29:LYS:HE2	5:A:1189:HOH:O	1.93	0.68
2:B:96:ARG:HH11	2:B:96:ARG:CG	2.06	0.68
1:C:254:ASP:OD2	5:C:1226:HOH:O	2.11	0.68
2:B:4:LYS:HG2	2:B:5:ALA:H	1.58	0.68
1:C:123:LYS:NZ	1:C:286:PHE:O	2.27	0.68
1:A:221:GLU:CG	5:A:1180:HOH:O	2.40	0.67
1:C:4:LEU:O	1:C:62:ASN:OD1	2.12	0.67
1:C:76:GLU:HB2	5:C:1529:HOH:O	1.96	0.66
1:A:281:PRO:CB	5:A:1438:HOH:O	2.42	0.66
1:A:20:GLN:HE22	1:A:287:LYS:H	1.41	0.66
1:A:134:PRO:O	5:A:1417:HOH:O	2.14	0.65
1:A:221:GLU:OE2	5:A:1180:HOH:O	2.07	0.65
1:C:52:HIS:NE2	5:C:1454:HOH:O	2.28	0.65
1:A:264:LYS:HD2	5:A:1225:HOH:O	1.97	0.64
1:A:239:ILE:HG22	1:A:245:LEU:HD23	1.79	0.64
1:A:270:LEU:HD23	1:A:270:LEU:N	2.12	0.64
1:A:32:GLU:HB3	5:A:1543:HOH:O	1.96	0.64
1:A:79:ASP:N	5:A:1513:HOH:O	2.30	0.64
1:A:32:GLU:CB	5:A:1543:HOH:O	2.45	0.64
1:A:200:ASN:H	1:A:200:ASN:HD22	1.46	0.63
1:C:260:LYS:HA	1:C:260:LYS:HE3	0.80	0.63
1:A:166:ARG:NH2	1:A:250:GLU:OE1	2.31	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:62:VAL:HG12	5:B:1535:HOH:O	1.99	0.63
1:A:264:LYS:HD3	5:A:1225:HOH:O	1.99	0.63
1:C:80:PRO:O	5:C:1314:HOH:O	2.16	0.63
2:B:3:PHE:CE1	2:B:97:ASN:HB3	2.34	0.62
2:B:39:GLY:O	2:B:43:ARG:HG3	1.99	0.62
1:C:196:ASN:HD22	1:C:196:ASN:H	1.45	0.62
1:C:289:LEU:HG	1:C:294:LEU:HD22	1.80	0.62
2:B:7:SER:HB3	5:B:1491:HOH:O	1.99	0.62
1:A:21:LYS:HE3	5:A:1382:HOH:O	1.98	0.62
1:A:216:ASN:HD22	1:A:222:GLN:NE2	1.98	0.62
1:C:15:SER:HB2	5:C:1161:HOH:O	1.99	0.62
2:B:77:LYS:CD	2:B:82:GLY:H	2.12	0.61
2:B:65:ASP:CG	5:B:1536:HOH:O	2.43	0.61
1:C:200:ASN:HD22	1:C:200:ASN:H	1.46	0.61
1:A:5:VAL:HG12	1:A:7:VAL:HG23	1.82	0.61
2:B:3:PHE:CD1	2:B:97:ASN:HB3	2.36	0.60
1:C:138:THR:HG23	5:C:1529:HOH:O	2.00	0.60
2:B:16:LYS:HB3	2:B:16:LYS:HZ3	1.67	0.60
1:C:134:PRO:HB2	5:C:1471:HOH:O	2.01	0.60
1:C:188:GLU:H	1:C:222:GLN:HE22	1.48	0.60
1:A:281:PRO:CA	5:A:1438:HOH:O	2.49	0.59
2:B:3:PHE:HE1	2:B:97:ASN:HA	1.67	0.59
1:A:99:PHE:O	1:A:102:ILE:HG22	2.02	0.59
2:B:48:ALA:HB3	2:B:53:TYR:OH	2.02	0.59
1:A:138:THR:HG23	5:A:1108:HOH:O	2.03	0.59
1:C:135:GLU:O	1:C:138:THR:HB	2.01	0.59
1:A:31:ARG:NH2	2:B:91:LYS:HB3	2.18	0.59
1:A:89:PHE:O	1:A:93:GLU:HG2	2.03	0.59
2:B:59:LYS:HG2	1:C:196:ASN:OD1	2.03	0.58
1:A:281:PRO:HA	5:A:1438:HOH:O	2.03	0.58
2:B:16:LYS:CB	2:B:16:LYS:NZ	2.67	0.58
1:A:156:THR:HG21	5:A:1421:HOH:O	2.02	0.58
2:B:93:GLU:HA	2:B:96:ARG:NH1	2.19	0.58
2:B:64:TRP:HZ2	4:B:109:HEC:HAA1	1.68	0.57
2:B:76:PRO:HD2	5:B:1495:HOH:O	2.03	0.57
2:B:3:PHE:HE1	2:B:97:ASN:HB2	1.70	0.57
1:C:144:LEU:HD12	1:C:145:PRO:CD	2.34	0.57
1:C:172:MET:HG2	1:C:238:LEU:HD11	1.87	0.56
1:A:77:PHE:CE2	5:A:1408:HOH:O	2.47	0.56
2:B:40:ILE:CD1	2:B:107:SER:HB2	2.36	0.56
1:A:152:GLY:O	1:A:156:THR:HG23	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:LYS:HD2	5:A:1404:HOH:O	2.05	0.56
1:A:73:PHE:HE2	5:A:1401:HOH:O	1.88	0.56
2:B:51:TYR:HB3	2:B:53:TYR:CE2	2.41	0.56
1:C:76:GLU:CB	5:C:1082:HOH:O	2.02	0.55
1:C:289:LEU:HD12	1:C:294:LEU:HD13	1.87	0.55
2:B:3:PHE:CE1	2:B:97:ASN:HA	2.42	0.55
1:C:216:ASN:HD22	1:C:222:GLN:HE21	1.54	0.55
1:A:69:GLY:HA2	5:A:1391:HOH:O	2.07	0.55
2:B:9:LYS:HE2	2:B:9:LYS:N	2.21	0.55
1:C:57:TRP:HE1	1:C:62:ASN:ND2	2.05	0.55
1:A:69:GLY:O	1:A:72:ARG:HG2	2.07	0.55
1:C:134:PRO:CB	5:C:1471:HOH:O	2.55	0.54
1:A:74:LYS:CD	5:A:1404:HOH:O	2.54	0.54
1:A:139:PRO:N	5:A:1108:HOH:O	2.40	0.54
1:A:2:THR:O	1:A:3:PRO:N	2.34	0.53
2:B:3:PHE:HE1	2:B:97:ASN:CA	2.22	0.53
1:C:6:HIS:CD2	5:C:1443:HOH:O	2.61	0.53
2:B:72:TYR:CE2	2:B:80:ILE:HD12	2.44	0.53
1:A:76:GLU:C	5:A:1405:HOH:O	2.49	0.53
1:A:134:PRO:HB2	1:A:136:ASP:OD1	2.09	0.53
1:A:216:ASN:HB2	5:A:1142:HOH:O	2.08	0.53
2:B:7:SER:CB	5:B:1491:HOH:O	2.57	0.53
1:C:134:PRO:N	5:C:1471:HOH:O	2.41	0.53
1:A:260:LYS:HG3	5:A:1437:HOH:O	2.09	0.53
1:A:221:GLU:HG3	5:A:1180:HOH:O	2.05	0.53
2:B:40:ILE:HD11	2:B:107:SER:HB2	1.91	0.53
2:B:40:ILE:CD1	2:B:40:ILE:N	2.31	0.52
1:A:214:GLU:OE1	1:A:224:ASP:HB2	2.10	0.52
1:C:57:TRP:HE1	1:C:62:ASN:HD22	1.57	0.52
2:B:33:VAL:C	2:B:51:TYR:OH	2.51	0.52
2:B:37:LEU:HA	2:B:40:ILE:HG13	1.91	0.52
1:C:36:TYR:OH	5:C:1075:HOH:O	2.19	0.52
1:A:79:ASP:C	1:A:80:PRO:HA	2.30	0.52
2:B:3:PHE:CE1	2:B:97:ASN:HB2	2.44	0.51
1:A:269:LEU:HB3	1:A:270:LEU:HD23	1.91	0.51
1:A:11:GLU:C	1:A:12:LYS:C	2.79	0.51
2:B:93:GLU:HA	2:B:96:ARG:HH12	1.75	0.51
1:A:52:HIS:ND1	1:A:144:LEU:HD13	2.25	0.51
1:A:138:THR:CG2	5:A:1108:HOH:O	2.58	0.51
1:C:86:GLN:CD	1:C:86:GLN:H	2.19	0.51
1:A:79:ASP:O	1:A:80:PRO:CA	2.59	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:ASN:H	1:A:255:GLN:HE21	1.59	0.51
1:A:138:THR:N	5:A:1417:HOH:O	2.44	0.50
1:A:200:ASN:HD22	1:A:200:ASN:N	2.09	0.50
1:A:123:LYS:NZ	1:A:286:PHE:O	2.45	0.50
1:C:245:LEU:O	1:C:249:LYS:HG3	2.11	0.50
2:B:37:LEU:HB3	2:B:40:ILE:HG12	1.93	0.49
1:C:276:PHE:O	5:C:1348:HOH:O	2.18	0.49
1:C:200:ASN:HD22	1:C:200:ASN:N	2.08	0.49
1:A:124:ILE:HD13	1:A:266:PHE:HD2	1.76	0.49
1:C:124:ILE:HD13	1:C:266:PHE:HD2	1.76	0.49
1:C:29:LYS:HG2	1:C:91:PHE:CE2	2.47	0.49
2:B:39:GLY:N	2:B:40:ILE:HD12	2.28	0.49
2:B:16:LYS:HB3	2:B:16:LYS:NZ	2.27	0.48
1:C:260:LYS:CE	1:C:260:LYS:CA	2.47	0.48
2:B:48:ALA:HB3	2:B:53:TYR:HH	1.76	0.48
1:A:73:PHE:HB2	5:A:1511:HOH:O	2.14	0.48
1:C:2:THR:HG22	1:C:3:PRO:HD2	1.94	0.48
1:C:68:GLY:HA2	1:C:133:THR:HG23	1.95	0.48
1:C:192:GLY:O	1:C:195:ASN:ND2	2.47	0.48
1:A:3:PRO:O	1:A:4:LEU:C	2.56	0.48
2:B:108:GLU:CG	2:B:108:GLU:OXT	2.59	0.48
1:A:2:THR:OG1	1:A:3:PRO:HD2	2.14	0.47
1:A:11:GLU:CA	1:A:12:LYS:N	2.71	0.47
1:A:294:LEU:N	1:A:294:LEU:HD23	2.29	0.47
1:C:245:LEU:HD22	1:C:249:LYS:HE3	1.96	0.47
2:B:34:GLY:HA2	2:B:51:TYR:OH	2.14	0.47
2:B:108:GLU:OXT	2:B:108:GLU:HG2	2.15	0.47
1:A:80:PRO:O	1:A:81:SER:C	2.57	0.47
1:A:84:GLY:H	1:A:86:GLN:HE22	1.62	0.46
2:B:73:LEU:HD13	2:B:99:LEU:HD23	1.97	0.46
1:C:76:GLU:CB	5:C:1529:HOH:O	2.60	0.46
1:A:226:LYS:HG2	5:A:1290:HOH:O	2.16	0.46
2:B:59:LYS:HB2	5:B:1507:HOH:O	2.15	0.46
1:C:80:PRO:HB3	5:C:1314:HOH:O	2.14	0.46
1:A:77:PHE:N	5:A:1513:HOH:O	2.46	0.46
2:B:38:HIS:HA	2:B:107:SER:HB3	1.97	0.46
2:B:69:MET:HG3	2:B:100:ILE:HD13	1.97	0.46
1:A:52:HIS:ND1	1:A:82:ASN:OD1	2.41	0.46
1:A:270:LEU:N	1:A:270:LEU:CD2	2.76	0.46
1:A:28:LEU:HG	1:A:32:GLU:OE1	2.16	0.45
1:A:19:PHE:HE1	1:A:102:ILE:HD12	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:TYR:N	5:C:1456:HOH:O	2.50	0.45
1:C:260:LYS:HE2	5:C:1031:HOH:O	2.17	0.45
1:A:2:THR:C	1:A:3:PRO:CD	2.78	0.45
1:A:86:GLN:O	1:A:90:LYS:HG3	2.17	0.45
1:A:102:ILE:HD11	1:A:106:ASP:HB3	1.98	0.45
1:C:79:ASP:HA	1:C:141:ASN:ND2	2.31	0.45
1:C:216:ASN:HD22	1:C:222:GLN:NE2	2.15	0.45
2:B:4:LYS:HG2	2:B:5:ALA:N	2.27	0.45
1:A:51:TRP:CD1	1:A:52:HIS:HD2	2.34	0.44
2:B:29:GLY:HA3	2:B:36:ASN:ND2	2.33	0.44
1:A:75:LYS:C	5:A:1513:HOH:O	2.52	0.44
1:A:148:ASP:O	1:A:148:ASP:OD2	2.36	0.44
2:B:29:GLY:HA3	2:B:36:ASN:HD21	1.83	0.44
1:C:80:PRO:CA	5:C:1314:HOH:O	2.65	0.44
1:A:156:THR:CG2	5:A:1085:HOH:O	2.39	0.43
1:C:134:PRO:CA	5:C:1471:HOH:O	2.65	0.43
1:A:73:PHE:CD1	1:A:135:GLU:HG3	2.54	0.43
2:B:34:GLY:CA	2:B:51:TYR:OH	2.66	0.43
2:B:70:SER:O	2:B:74:THR:HG22	2.18	0.43
1:A:29:LYS:NZ	1:A:33:ASP:OD2	2.51	0.43
1:C:48:ARG:NH2	5:C:1460:HOH:O	2.50	0.43
2:B:77:LYS:O	2:B:77:LYS:HG3	2.19	0.42
1:C:78:ASN:O	5:C:1459:HOH:O	2.21	0.42
2:B:15:PHE:CZ	2:B:25:VAL:HG13	2.54	0.42
1:A:207:LEU:HD22	5:A:1287:HOH:O	2.18	0.42
1:A:12:LYS:CE	5:A:1378:HOH:O	2.44	0.42
1:A:179:LYS:HE3	5:A:1275:HOH:O	2.19	0.42
2:B:13:THR:HA	2:B:16:LYS:HG2	2.01	0.42
1:C:137:THR:CG2	5:C:1471:HOH:O	2.59	0.42
1:C:71:TYR:HB2	5:C:1456:HOH:O	2.19	0.42
1:A:31:ARG:CZ	2:B:91:LYS:HB3	2.50	0.42
1:A:124:ILE:HD13	1:A:266:PHE:CD2	2.53	0.42
1:A:11:GLU:O	1:A:12:LYS:C	2.64	0.41
1:A:97:LYS:NZ	5:A:1412:HOH:O	2.36	0.41
1:A:133:THR:HB	1:A:137:THR:HG21	2.02	0.41
1:A:79:ASP:C	1:A:80:PRO:C	2.87	0.41
2:B:54:THR:HG22	4:B:109:HEC:O2D	2.20	0.41
2:B:23:HIS:CD2	2:B:37:LEU:HD11	2.56	0.41
1:C:2:THR:HB	1:C:3:PRO:HD2	2.03	0.41
1:C:138:THR:HA	1:C:139:PRO:HD3	1.78	0.41
2:B:31:HIS:HB2	2:B:51:TYR:CD1	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:40:ILE:O	2:B:64:TRP:HB2	2.20	0.41
1:C:6:HIS:HD2	5:C:1443:HOH:O	1.99	0.41
1:A:19:PHE:CE1	1:A:102:ILE:HD12	2.56	0.41
2:B:104:LYS:HB2	2:B:104:LYS:HE2	1.38	0.41
1:C:264:LYS:HE2	5:C:1542:HOH:O	2.21	0.41
2:B:69:MET:HE1	4:B:109:HEC:CMA	2.51	0.40
2:B:69:MET:HE1	4:B:109:HEC:HMA2	2.03	0.40
1:A:35:GLU:HB2	5:A:1385:HOH:O	2.20	0.40
2:B:101:THR:CA	2:B:104:LYS:HD3	2.29	0.40

All (9) 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)
1:A:226:LYS:NZ	5:A:1233:HOH:O[1_445]	0.39	1.81
5:C:1321:HOH:O	5:C:1332:HOH:O[1_655]	0.68	1.52
1:A:226:LYS:CE	5:A:1233:HOH:O[1_445]	1.14	1.06
1:A:1:THR:OG1	5:C:1472:HOH:O[1_564]	1.86	0.34
1:A:17:GLU:OE1	2:B:49:GLU:OE1[1_665]	1.87	0.33
1:C:84:GLY:CA	1:C:278:LYS:NZ[1_455]	2.02	0.18
2:B:2:GLU:OE2	1:C:93:GLU:OE1[1_565]	2.05	0.15
1:C:291:GLU:OE1	5:B:1168:HOH:O[1_655]	2.06	0.14
2:B:108:GLU:OE2	1:C:256:ASP:OD1[1_455]	2.18	0.02

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	292/296 (99%)	281 (96%)	5 (2%)	6 (2%)	5 1
1	C	292/296 (99%)	284 (97%)	6 (2%)	2 (1%)	18 7
2	B	106/108 (98%)	97 (92%)	6 (6%)	3 (3%)	4 0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	690/700 (99%)	662 (96%)	17 (2%)	11 (2%)	7 1

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	C	279	ASP
1	C	280	ALA
1	A	2	THR
2	B	66	GLU
1	A	134	PRO
1	A	136	ASP
2	B	32	LYS
1	A	11	GLU
1	A	12	LYS
2	B	4	LYS

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	252/254 (99%)	232 (92%)	20 (8%)	11 2
1	C	252/254 (99%)	228 (90%)	24 (10%)	8 1
2	B	89/89 (100%)	69 (78%)	20 (22%)	1 0
All	All	593/597 (99%)	529 (89%)	64 (11%)	6 1

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	THR
1	A	2	THR
1	A	12	LYS
1	A	32	GLU
1	A	51	TRP

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Mol	Chain	Res	Type
1	A	53	ILE
1	A	86	GLN
1	A	93	GLU
1	A	95	ILE
1	A	102	ILE
1	A	168	VAL
1	A	196	ASN
1	A	200	ASN
1	A	215	LYS
1	A	219	ASN
1	A	245	LEU
1	A	269	LEU
1	A	270	LEU
1	A	278	LYS
1	A	294	LEU
2	B	3	PHE
2	B	4	LYS
2	B	16	LYS
2	B	21	GLN
2	B	24	THR
2	B	32	LYS
2	B	40	ILE
2	B	47	GLN
2	B	60	LYS
2	B	63	LEU
2	B	68	ASN
2	B	69	MET
2	B	84	LYS
2	B	94	LYS
2	B	96	ARG
2	B	102	TYR
2	B	103	LEU
2	B	104	LYS
2	B	105	LYS
2	B	107	SER
1	C	12	LYS
1	C	28	LEU
1	C	30	LEU
1	C	31	ARG
1	C	32	GLU
1	C	35	GLU
1	C	49	LEU

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Mol	Chain	Res	Type
1	C	51	TRP
1	C	72	ARG
1	C	74	LYS
1	C	135	GLU
1	C	144	LEU
1	C	168	VAL
1	C	195	ASN
1	C	196	ASN
1	C	200	ASN
1	C	214	GLU
1	C	226	LYS
1	C	245	LEU
1	C	260	LYS
1	C	269	LEU
1	C	282	SER
1	C	289	LEU
1	C	294	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	HIS
1	A	20	GLN
1	A	24	ASN
1	A	38	ASN
1	A	86	GLN
1	A	120	GLN
1	A	159	GLN
1	A	181	HIS
1	A	195	ASN
1	A	196	ASN
1	A	200	ASN
1	A	208	ASN
1	A	222	GLN
1	A	253	ASN
1	A	255	GLN
1	A	292	GLN
2	B	21	GLN
2	B	31	HIS
2	B	36	ASN
2	B	38	HIS
2	B	57	ASN

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Mol	Chain	Res	Type
2	B	61	ASN
2	B	68	ASN
1	C	6	HIS
1	C	20	GLN
1	C	24	ASN
1	C	62	ASN
1	C	82	ASN
1	C	159	GLN
1	C	195	ASN
1	C	200	ASN
1	C	208	ASN
1	C	222	GLN
1	C	292	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](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	HEC	B	109	2	46,50,50	1.29	3 (6%)	58,82,82	0.97	1 (1%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ZNH	C	295	1	50,50,50	1.52	10 (20%)	65,82,82	1.44	9 (13%)
3	ZNH	A	295	1	50,50,50	1.09	2 (4%)	65,82,82	1.69	12 (18%)

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	HEC	B	109	2	-	5/14/54/54	-
3	ZNH	C	295	1	1/1/3/9	2/14/54/54	-
3	ZNH	A	295	1	1/1/3/9	4/14/54/54	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	295	ZNH	C1A-C2A	-3.70	1.38	1.45
3	C	295	ZNH	CMC-C2C	3.48	1.57	1.50
4	B	109	HEC	CBB-CAB	-3.19	1.37	1.49
3	C	295	ZNH	C4A-C3A	-2.90	1.39	1.45
4	B	109	HEC	CBC-CAC	-2.90	1.38	1.49
3	C	295	ZNH	ZN-NB	2.82	2.12	2.03
3	C	295	ZNH	C1D-ND	-2.70	1.33	1.38
3	C	295	ZNH	C1B-NB	-2.40	1.32	1.37
3	C	295	ZNH	C1B-C2B	2.27	1.49	1.44
3	C	295	ZNH	CAC-C3C	-2.18	1.41	1.47
3	A	295	ZNH	CMC-C2C	2.17	1.55	1.50
3	C	295	ZNH	CAB-C3B	-2.14	1.41	1.47
4	B	109	HEC	CAC-C3C	2.09	1.41	1.35
3	A	295	ZNH	CBC-CAC	2.06	1.40	1.30
3	C	295	ZNH	ZN-NA	2.01	2.10	2.03

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	295	ZNH	C4B-C3B-C2B	-4.86	102.82	107.28
3	A	295	ZNH	CAD-C3D-C4D	4.85	133.16	124.70
3	A	295	ZNH	CAD-C3D-C2D	-4.34	119.74	127.87
3	C	295	ZNH	C4B-C3B-C2B	-4.33	103.30	107.28
3	A	295	ZNH	CBD-CAD-C3D	4.19	124.11	112.53
3	C	295	ZNH	CHC-C4B-NB	-3.71	119.97	124.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	295	ZNH	CMC-C2C-C1C	-3.06	120.76	125.42
3	C	295	ZNH	C2C-C1C-NC	2.88	112.64	109.02
3	A	295	ZNH	C4C-C3C-C2C	-2.82	103.80	107.30
3	C	295	ZNH	C4C-C3C-C2C	-2.76	103.87	107.30
3	A	295	ZNH	C3B-C2B-C1B	2.75	108.48	106.41
3	A	295	ZNH	CHA-C4D-C3D	2.71	128.73	124.77
3	C	295	ZNH	C1C-CHC-C4B	2.71	132.61	126.25
3	C	295	ZNH	CBA-CAA-C2A	2.68	119.94	112.53
3	A	295	ZNH	CBA-CAA-C2A	2.34	118.99	112.53
3	A	295	ZNH	C2C-C1C-NC	2.31	111.92	109.02
4	B	109	HEC	CHC-C4B-NB	-2.28	121.97	124.45
3	C	295	ZNH	CHD-C4C-NC	-2.28	120.31	124.24
3	C	295	ZNH	C1A-CHA-C4D	2.17	130.64	126.02
3	C	295	ZNH	C4D-C3D-C2D	-2.17	103.73	106.89
3	A	295	ZNH	CHC-C4B-NB	-2.04	122.11	124.75
3	A	295	ZNH	CMC-C2C-C3C	2.00	131.26	126.55

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	295	ZNH	NA
3	C	295	ZNH	NA

All (11) torsion outliers are listed below:

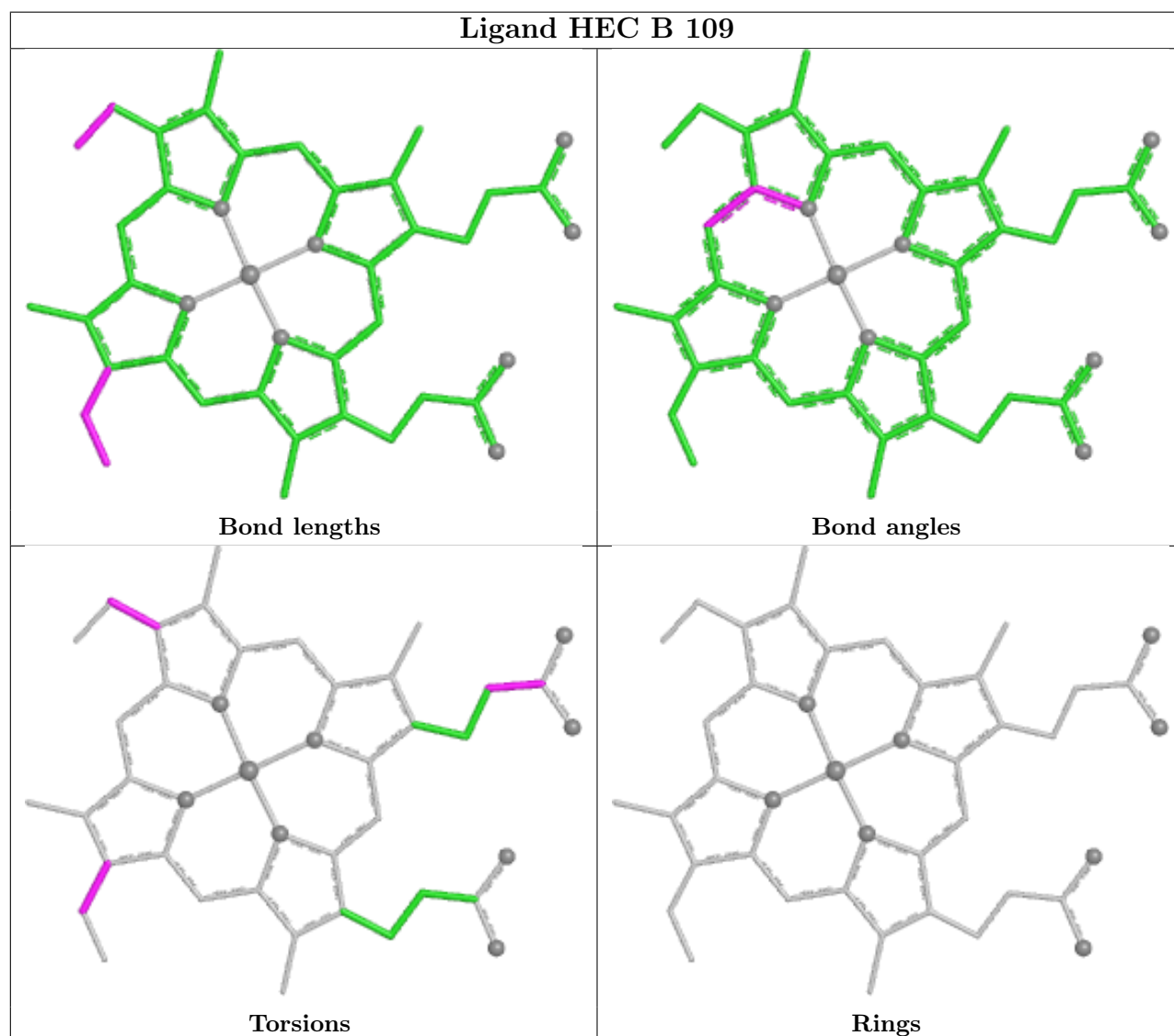
Mol	Chain	Res	Type	Atoms
4	B	109	HEC	C4C-C3C-CAC-CBC
3	A	295	ZNH	C4D-C3D-CAD-CBD
3	A	295	ZNH	C2D-C3D-CAD-CBD
4	B	109	HEC	C2C-C3C-CAC-CBC
3	C	295	ZNH	CAA-CBA-CGA-O2A
3	A	295	ZNH	CAA-CBA-CGA-O2A
3	A	295	ZNH	CAA-CBA-CGA-O1A
3	C	295	ZNH	CAA-CBA-CGA-O1A
4	B	109	HEC	C4B-C3B-CAB-CBB
4	B	109	HEC	CAA-CBA-CGA-O2A
4	B	109	HEC	CAA-CBA-CGA-O1A

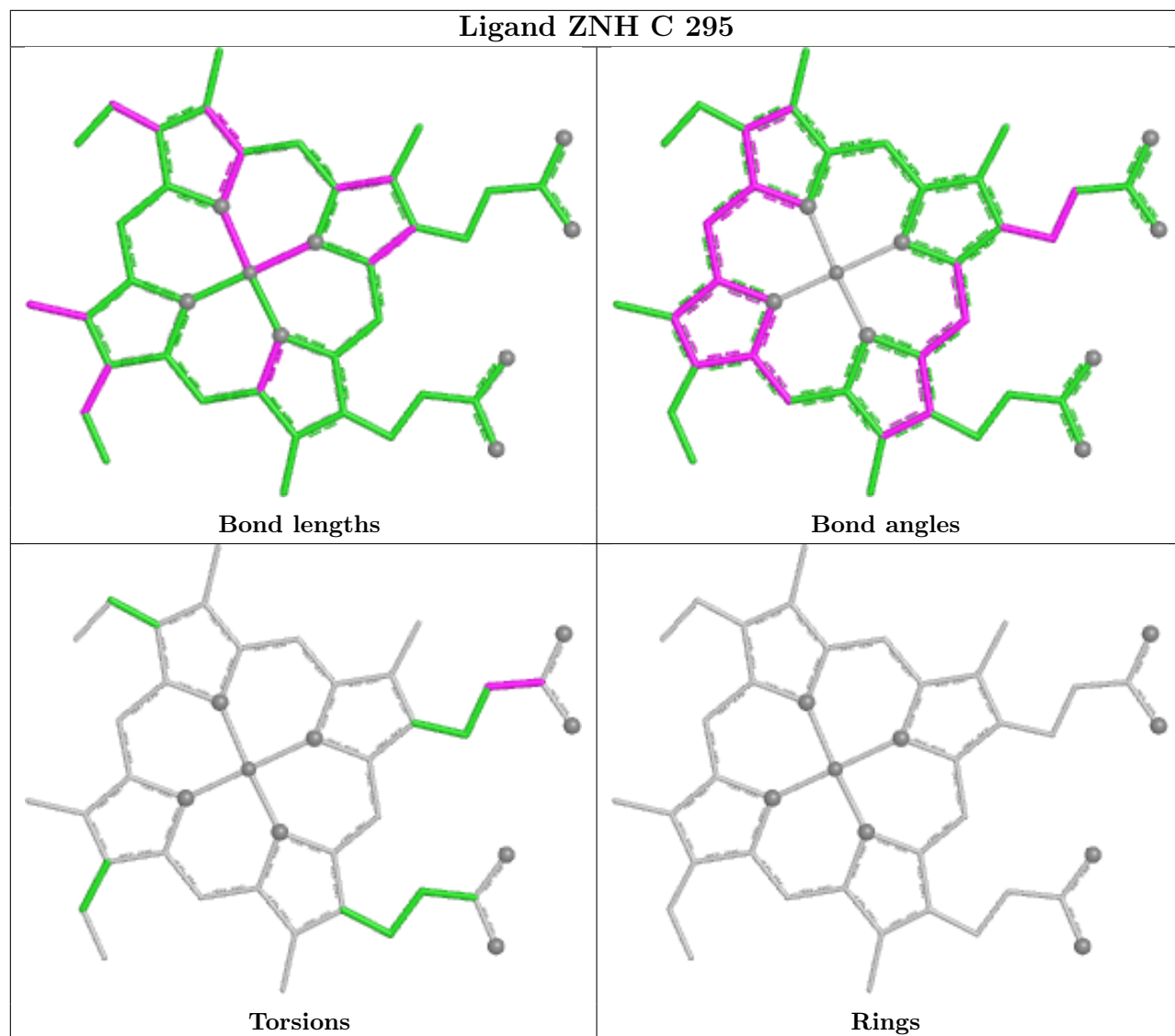
There are no ring outliers.

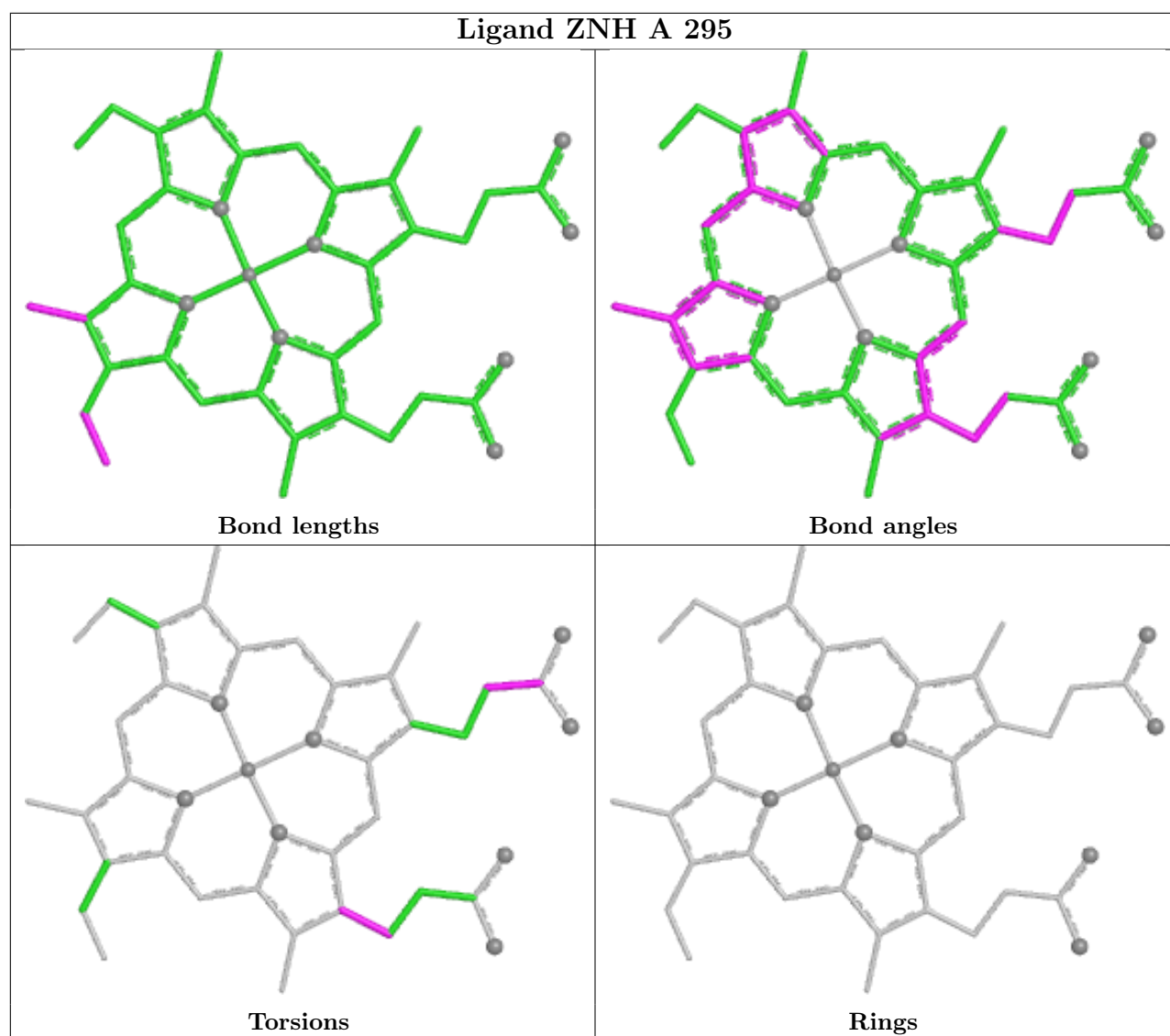
1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	109	HEC	5	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	7
1	C	2
2	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	11:GLU	C	12:LYS	N	1.74
1	A	79:ASP	C	80:PRO	N	1.74
1	A	2:THR	C	3:PRO	N	1.69
1	C	79:ASP	C	80:PRO	N	1.66
1	C	277:PRO	C	278:LYS	N	1.65
1	A	134:PRO	C	135:GLU	N	1.63
1	B	9:LYS	C	10:LYS	N	1.20
1	A	73:PHE	C	74:LYS	N	1.18
1	A	135:GLU	C	136:ASP	N	1.14
1	A	1:THR	C	2:THR	N	0.65

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	294/296 (99%)	1.05	42 (14%) 6 5	20, 24, 36, 50	0
1	C	294/296 (99%)	1.17	52 (17%) 4 3	18, 24, 39, 49	0
2	B	108/108 (100%)	2.98	87 (80%) 0 0	18, 34, 49, 50	0
All	All	696/700 (99%)	1.40	181 (26%) 1 1	18, 25, 43, 50	0

All (181) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	THR	7.6
2	B	55	ASP	6.8
2	B	3	PHE	6.7
1	C	138	THR	6.2
2	B	58	ILE	6.1
2	B	30	PRO	5.9
1	A	1	THR	5.9
1	C	77	PHE	5.6
1	C	1	THR	5.2
2	B	56	ALA	5.2
2	B	48	ALA	5.1
1	C	79	ASP	5.1
1	C	81	SER	5.0
1	C	294	LEU	4.8
2	B	101	THR	4.8
2	B	50	GLY	4.6
2	B	22	CYS	4.6
2	B	106	ALA	4.6
2	B	61	ASN	4.5
1	C	35	GLU	4.4
1	C	80	PRO	4.3
2	B	64	TRP	4.3
2	B	57	ASN	4.3

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Mol	Chain	Res	Type	RSRZ
2	B	59	LYS	4.3
1	C	135	GLU	4.3
2	B	34	GLY	4.2
2	B	54	THR	4.2
1	C	4	LEU	4.2
2	B	51	TYR	4.2
2	B	83	THR	4.1
1	C	71	TYR	4.1
1	C	72	ARG	4.0
2	B	33	VAL	4.0
2	B	99	LEU	3.9
1	A	76	GLU	3.9
1	C	76	GLU	3.9
1	C	73	PHE	3.9
2	B	63	LEU	3.9
2	B	39	GLY	3.9
1	C	39	TYR	3.9
1	C	134	PRO	3.8
2	B	89	GLY	3.7
2	B	53	TYR	3.7
1	C	82	ASN	3.7
1	A	3	PRO	3.7
1	A	4	LEU	3.7
1	C	210	ASP	3.7
2	B	62	VAL	3.7
1	A	78	ASN	3.7
2	B	49	GLU	3.6
2	B	24	THR	3.6
2	B	29	GLY	3.6
1	A	148	ASP	3.6
2	B	41	PHE	3.6
2	B	107	SER	3.6
1	A	73	PHE	3.5
2	B	31	HIS	3.5
2	B	37	LEU	3.5
1	C	40	ILE	3.4
1	A	80	PRO	3.4
2	B	74	THR	3.3
1	C	3	PRO	3.3
1	A	139	PRO	3.3
1	C	139	PRO	3.3
1	C	2	THR	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	136	ASP	3.2
1	A	89	PHE	3.2
2	B	87	PHE	3.2
2	B	84	LYS	3.2
2	B	81	PRO	3.2
2	B	27	LYS	3.2
1	A	71	TYR	3.1
2	B	2	GLU	3.1
2	B	5	ALA	3.1
1	C	89	PHE	3.1
1	C	168	VAL	3.1
1	C	34	ASP	3.1
2	B	9	LYS	3.1
2	B	1	THR	3.1
1	C	75	LYS	3.0
2	B	73	LEU	3.0
1	A	74	LYS	3.0
1	C	293	GLY	3.0
1	A	280	ALA	3.0
2	B	86	ALA	3.0
2	B	52	SER	3.0
2	B	42	GLY	3.0
1	C	141	ASN	3.0
2	B	8	ALA	3.0
2	B	90	LEU	3.0
2	B	72	TYR	2.9
2	B	32	LYS	2.9
2	B	17	THR	2.9
1	C	280	ALA	2.9
2	B	102	TYR	2.9
1	C	195	ASN	2.9
1	C	219	ASN	2.9
1	A	77	PHE	2.8
2	B	60	LYS	2.8
1	A	39	TYR	2.8
1	C	279	ASP	2.8
2	B	79	TYR	2.8
1	C	148	ASP	2.8
2	B	108	GLU	2.7
1	C	278	LYS	2.7
1	A	294	LEU	2.7
2	B	20	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	138	THR	2.7
2	B	88	GLY	2.7
1	C	97	LYS	2.7
2	B	76	PRO	2.7
2	B	95	ASP	2.7
1	C	5	VAL	2.7
1	A	219	ASN	2.7
1	A	281	PRO	2.7
2	B	6	GLY	2.7
1	C	246	SER	2.6
2	B	77	LYS	2.6
2	B	44	HIS	2.6
2	B	103	LEU	2.6
1	A	16	TYR	2.6
2	B	38	HIS	2.6
1	A	70	THR	2.6
1	C	137	THR	2.6
1	A	140	ASP	2.6
1	C	16	TYR	2.6
2	B	4	LYS	2.6
1	C	245	LEU	2.5
2	B	25	VAL	2.5
2	B	100	ILE	2.5
2	B	46	GLY	2.5
1	C	70	THR	2.5
1	A	278	LYS	2.5
2	B	45	SER	2.5
2	B	43	ARG	2.5
1	A	5	VAL	2.5
1	A	12	LYS	2.5
1	C	133	THR	2.5
2	B	69	MET	2.5
2	B	71	GLU	2.4
2	B	11	GLY	2.4
2	B	28	GLY	2.4
2	B	40	ILE	2.4
1	A	79	ASP	2.4
2	B	65	ASP	2.4
1	A	285	ILE	2.4
1	A	210	ASP	2.4
2	B	93	GLU	2.4
1	A	263	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	142	GLY	2.3
2	B	97	ASN	2.3
2	B	35	PRO	2.3
1	A	137	THR	2.3
2	B	36	ASN	2.3
2	B	85	MET	2.3
2	B	23	HIS	2.2
1	A	17	GLU	2.2
1	C	196	ASN	2.2
1	C	193	ALA	2.2
2	B	82	GLY	2.2
2	B	47	GLN	2.2
2	B	80	ILE	2.2
1	C	144	LEU	2.1
1	C	74	LYS	2.1
1	C	208	ASN	2.1
1	C	227	SER	2.1
1	A	259	PHE	2.1
1	C	194	ALA	2.1
1	C	218	ALA	2.1
2	B	68	ASN	2.1
1	C	223	TRP	2.1
2	B	66	GLU	2.1
1	C	248	VAL	2.1
2	B	7	SER	2.1
1	A	82	ASN	2.0
1	A	141	ASN	2.0
1	A	28	LEU	2.0
1	A	144	LEU	2.0
1	A	97	LYS	2.0
1	A	215	LYS	2.0
1	A	279	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

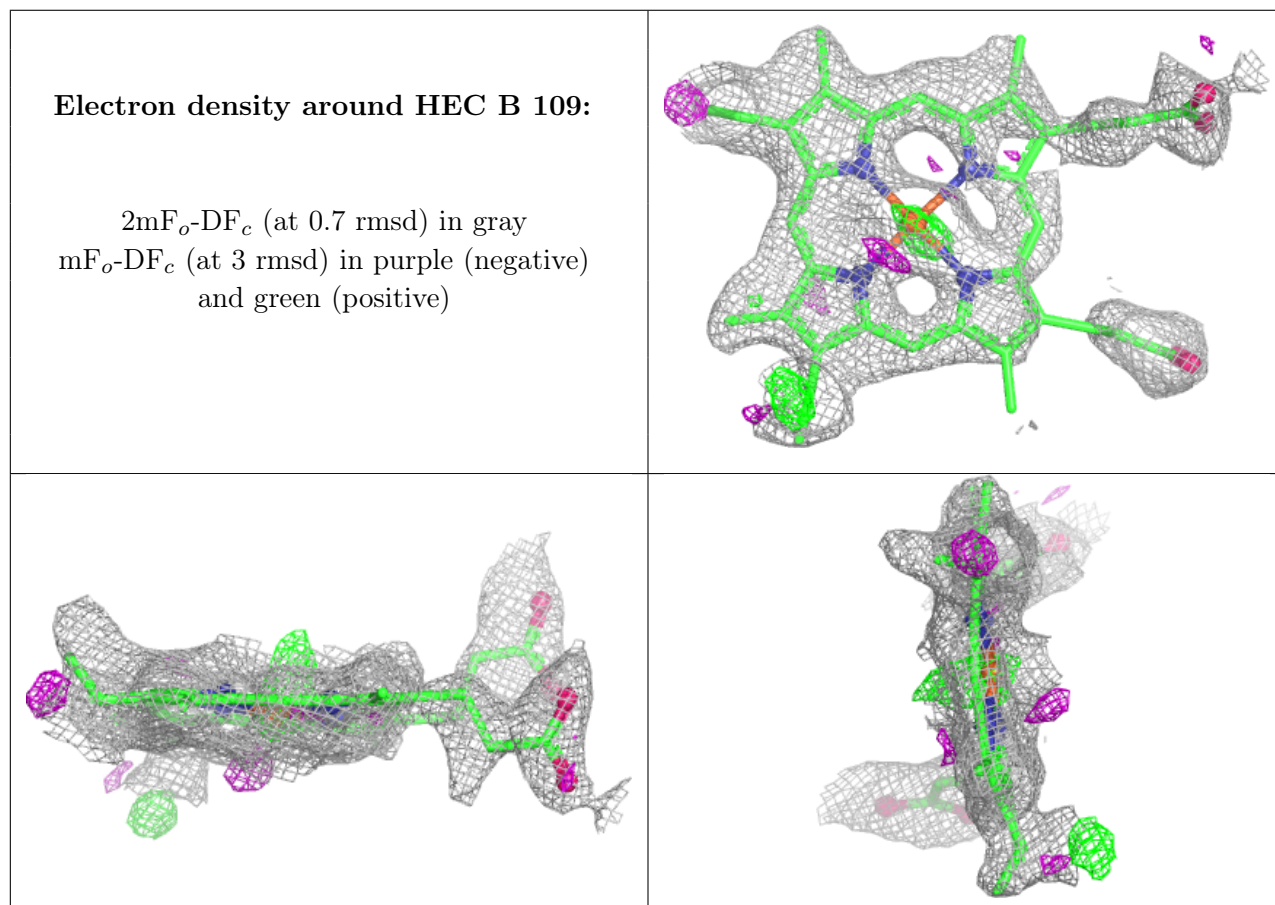
There are no oligosaccharides in this entry.

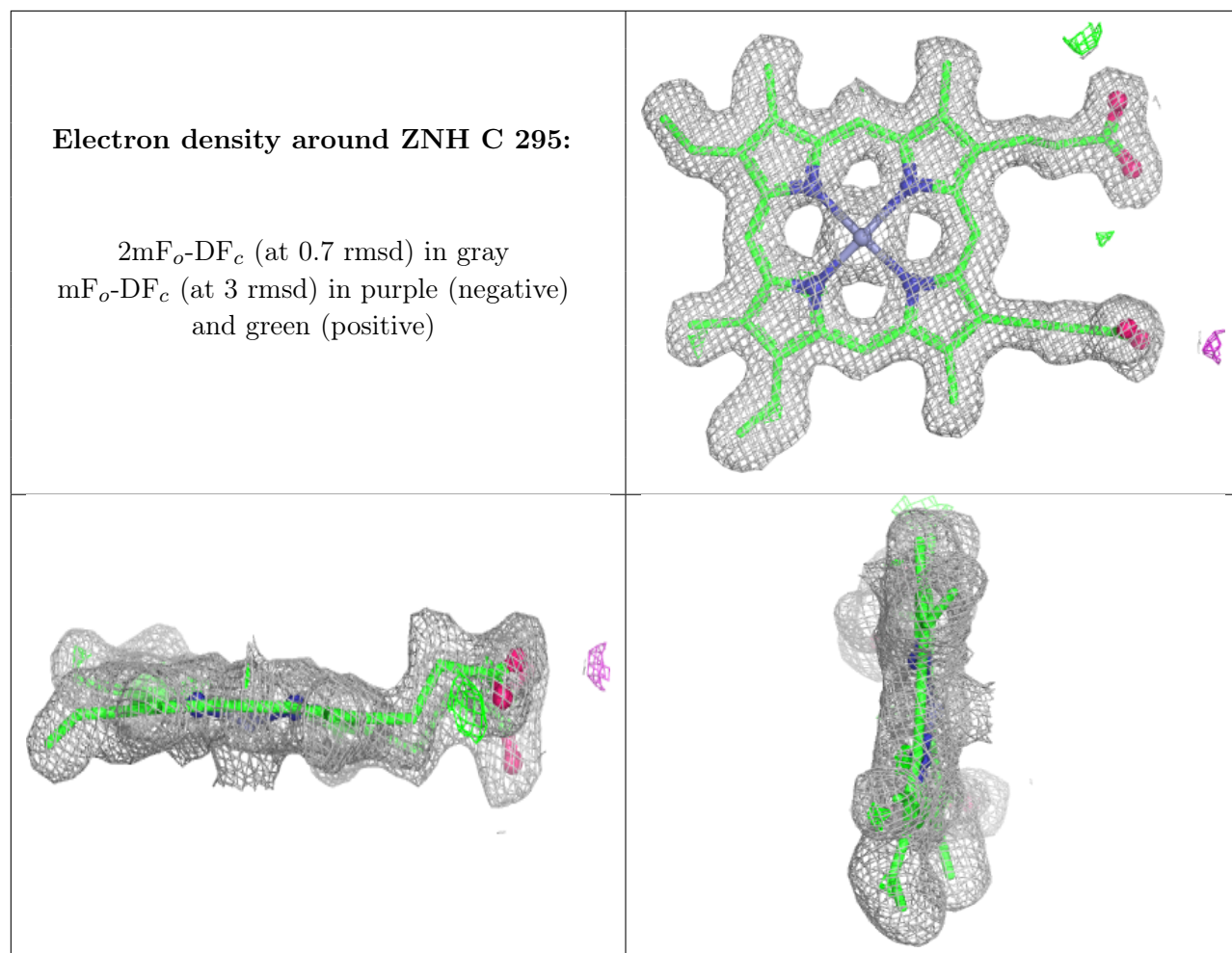
6.4 Ligands [i](#)

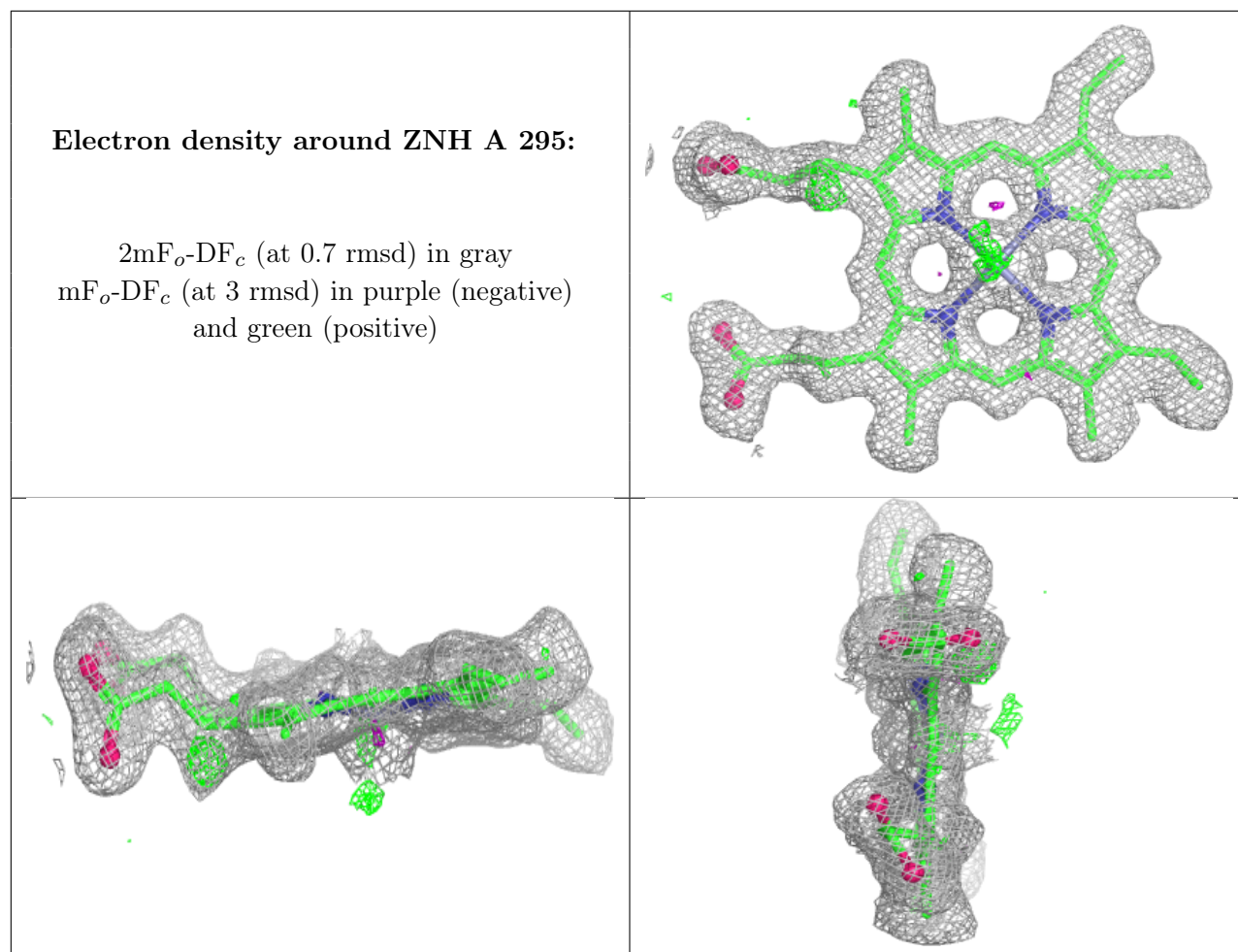
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	HEC	B	109	43/43	0.83	0.13	17,20,21,23	0
3	ZNH	C	295	43/43	0.97	0.09	19,21,23,26	0
3	ZNH	A	295	43/43	0.98	0.08	19,22,26,28	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.







6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.