Full wwPDB X-ray Structure Validation Report (i)

Aug 5, 2026 – 05:23 PM EDT

PDB ID	:	9CBO / pdb_00009cbo
Title	:	Methionine synthase from Thermus thermophilus HB8, Folate and Cobalamin Domains, Pre-Catalytic Conformation (Pre-Fol-on)
Authors	:	Yamada, K.; Mendoza, J.; Koutmos, M.
Deposited on	:	2024-06-19
Resolution	:	3.34 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 (i) 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 (i)) were used in the production of this report:

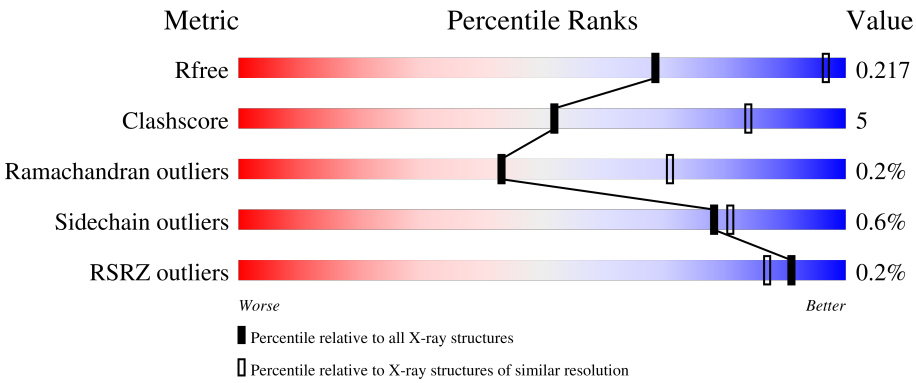
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.34 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	1434 (3.38-3.30)
Clashscore	190562	1479 (3.38-3.30)
Ramachandran outliers	187476	1456 (3.38-3.30)
Sidechain outliers	187428	1455 (3.38-3.30)
RSRZ outliers	180081	1434 (3.38-3.30)

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 <=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	511	86% 11% • •
1	B	511	89% 11%
1	C	511	89% 10% •
1	D	511	88% 11% • •
1	E	511	88% 11%

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Mol	Chain	Length	Quality of chain
1	F	511	91% 9%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	COB	A	901	X	-	-	-
2	COB	B	901	X	-	-	-
2	COB	C	901	X	-	-	-
2	COB	D	901	X	-	-	-
2	COB	E	901	X	-	-	-
2	COB	F	901	X	-	-	-
4	TRS	A	905	-	X	-	-

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 24580 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.

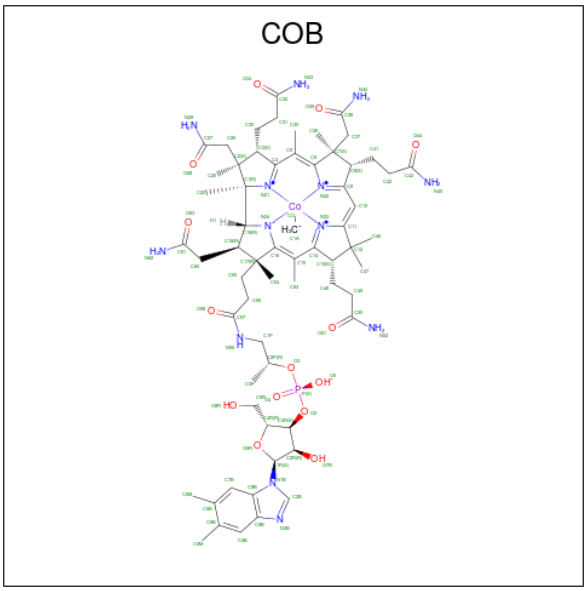
- Molecule 1 is a protein called Methionine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	510	3978	2536	696	731	15	0	0	0
1	B	510	3978	2536	696	731	15	0	0	0
1	C	510	3978	2536	696	731	15	0	0	0
1	D	507	3955	2523	692	725	15	0	0	0
1	E	509	3969	2531	694	729	15	0	0	0
1	F	510	3978	2536	696	731	15	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

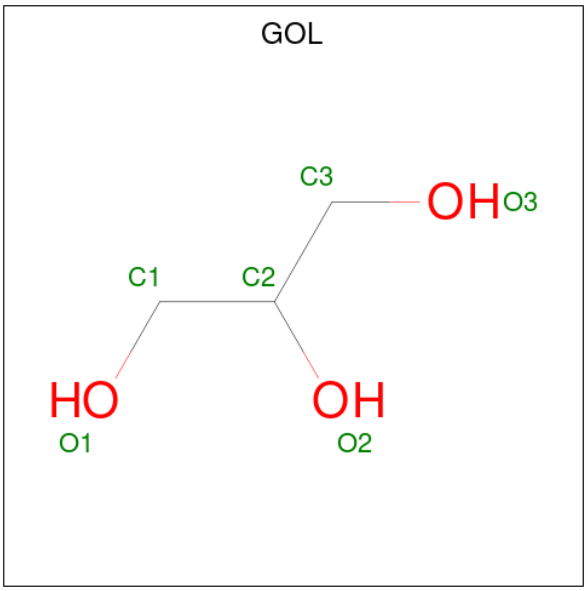
Chain	Residue	Modelled	Actual	Comment	Reference
A	759	ALA	ASP	engineered mutation	UNP Q5SKM5
B	759	ALA	ASP	engineered mutation	UNP Q5SKM5
C	759	ALA	ASP	engineered mutation	UNP Q5SKM5
D	759	ALA	ASP	engineered mutation	UNP Q5SKM5
E	759	ALA	ASP	engineered mutation	UNP Q5SKM5
F	759	ALA	ASP	engineered mutation	UNP Q5SKM5

- Molecule 2 is CO-METHYLCOBALAMIN (CCD ID: COB) (formula: $C_{63}H_{92}CoN_{13}O_{14}P$) (labeled as "Ligand of Interest" by depositor).



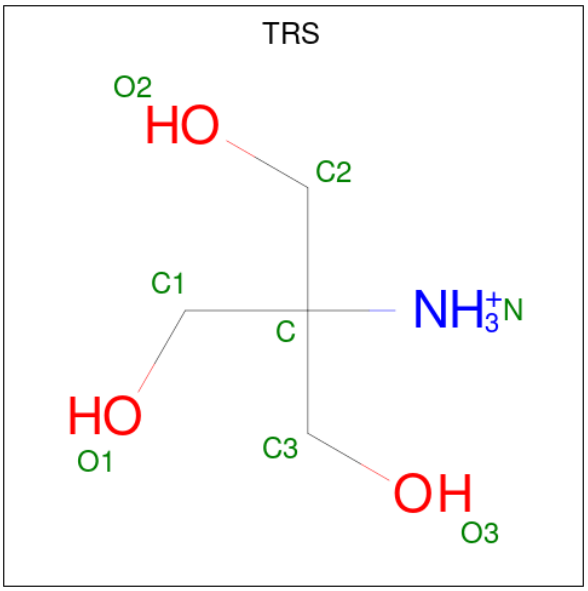
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	Co	N	O	P		
2	A	1	92	63	1	13	14	1	0	0
2	B	1	92	63	1	13	14	1	0	0
2	C	1	92	63	1	13	14	1	0	0
2	D	1	92	63	1	13	14	1	0	0
2	E	1	92	63	1	13	14	1	0	0
2	F	1	92	63	1	13	14	1	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	A	1	6	3	3	0	0
			Total	C	O		
3	B	1	6	3	3	0	0
			Total	C	O		
3	C	1	6	3	3	0	0
			Total	C	O		
3	C	1	6	3	3	0	0
			Total	C	O		
3	C	1	6	3	3	0	0
			Total	C	O		
3	C	1	6	3	3	0	0
			Total	C	O		
3	D	1	6	3	3	0	0
			Total	C	O		
3	E	1	6	3	3	0	0
			Total	C	O		
3	F	1	6	3	3	0	0

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	8	4	1	3	0	0
			Total	C	N	O		
4	A	1	8	4	1	3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	27	27	27	0	0
			Total	O		
5	B	18	18	18	0	0
			Total	O		
5	C	26	26	26	0	0
			Total	O		
5	D	17	17	17	0	0
			Total	O		
5	E	6	6	6	0	0
			Total	O		
5	F	16	16	16	0	0

3 Residue-property plots (i)

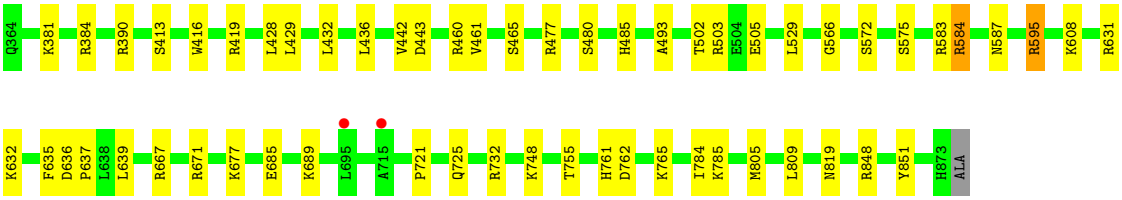
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

Chain A:

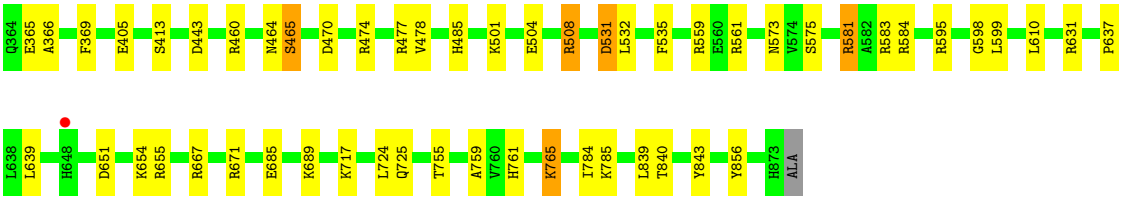
• Molecule 1: Methionine synthase

Chain B:



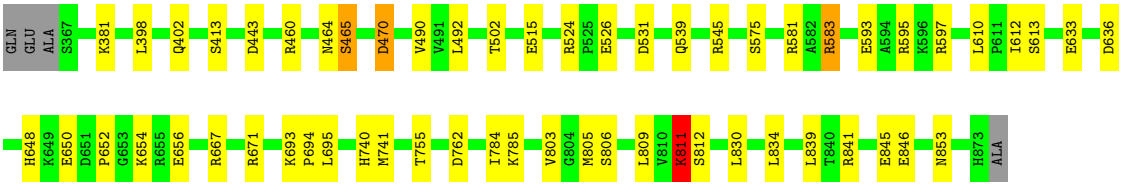
• Molecule 1: Methionine synthase

Chain C:



• Molecule 1: Methionine synthase

Chain D:

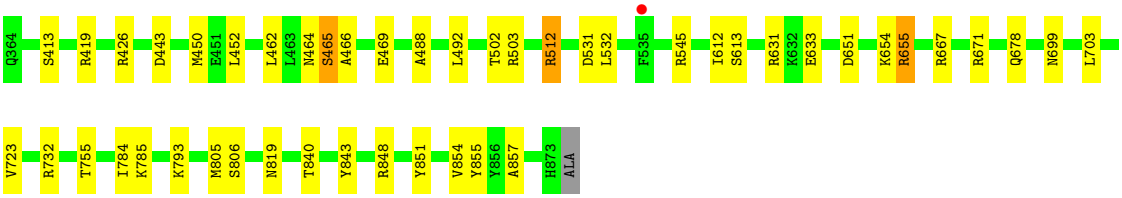


- Molecule 1: Methionine synthase

Chain E:

- Molecule 1: Methionine synthase

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	187.77Å 187.77Å 325.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	163.18 – 3.34 162.65 – 3.34	Depositor EDS
% Data completeness (in resolution range)	99.9 (163.18-3.34) 99.9 (162.65-3.34)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.47 (at 3.33Å)	Xtriage
Refinement program	REFMAC 5.8.0431	Depositor
R, R_{free}	0.176 , 0.219 0.178 , 0.217	Depositor DCC
R_{free} test set	4227 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å^2)	100.1	Xtriage
Anisotropy	0.339	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 127.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	24580	wwPDB-VP
Average B, all atoms (Å^2)	119.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.83% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: COB, GOL, TRS

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	0.93	10/4044 (0.2%)	1.14	18/5458 (0.3%)
1	B	0.82	4/4044 (0.1%)	1.02	1/5458 (0.0%)
1	C	0.83	2/4044 (0.0%)	1.06	4/5458 (0.1%)
1	D	0.81	1/4021 (0.0%)	1.05	10/5427 (0.2%)
1	E	0.77	1/4035 (0.0%)	0.97	4/5446 (0.1%)
1	F	0.78	0/4044	1.05	5/5458 (0.1%)
All	All	0.83	18/24232 (0.1%)	1.05	42/32705 (0.1%)

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	11
1	B	0	3
1	C	0	5
1	D	0	3
1	E	0	5
1	F	0	5
All	All	0	32

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	677	LYS	C-N	15.66	1.54	1.33
1	A	800	PRO	N-CD	-14.81	1.27	1.47
1	A	840	THR	C-N	9.60	1.46	1.33
1	A	841	ARG	C-N	-7.90	1.23	1.33
1	A	503	ARG	C-N	-7.20	1.24	1.33
1	A	831	PRO	C-N	-7.10	1.24	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	485	HIS	CG-CD2	-6.33	1.28	1.35
1	D	740	HIS	CG-CD2	-6.11	1.29	1.35
1	B	485	HIS	CG-CD2	-6.07	1.29	1.35
1	A	382	ARG	NE-CZ	-6.03	1.26	1.33
1	A	512	ARG	CZ-NH2	-5.73	1.26	1.33
1	B	503	ARG	NE-CZ	5.68	1.39	1.33
1	A	799	LYS	C-N	5.63	1.40	1.33
1	B	477	ARG	CZ-NH1	5.36	1.40	1.32
1	C	761	HIS	ND1-CE1	5.30	1.37	1.32
1	A	584	ARG	NE-CZ	-5.19	1.27	1.33
1	B	761	HIS	CE1-NE2	5.05	1.37	1.32
1	A	546	PRO	N-CA	-5.04	1.44	1.47

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	382	ARG	NE-CZ-NH1	-9.03	112.47	121.50
1	F	502	THR	OG1-CB-CG2	-8.42	92.45	109.30
1	D	502	THR	OG1-CB-CG2	-8.09	93.11	109.30
1	D	841	ARG	NE-CZ-NH1	-7.87	113.63	121.50
1	A	584	ARG	NE-CZ-NH1	-7.67	113.83	121.50
1	A	382	ARG	CG-CD-NE	-7.64	95.19	112.00
1	A	614	GLN	OE1-CD-NE2	7.24	129.84	122.60
1	C	508	ARG	NE-CZ-NH1	-7.05	114.45	121.50
1	A	508	ARG	NE-CZ-NH1	-6.84	114.66	121.50
1	A	631	ARG	NE-CZ-NH1	-6.65	114.85	121.50
1	E	677	LYS	CA-C-N	-6.27	112.17	120.95
1	E	677	LYS	C-N-CA	-6.27	112.17	120.95
1	F	654	LYS	N-CA-C	-5.99	104.83	111.36
1	A	531	ASP	CA-CB-CG	5.93	118.53	112.60
1	E	540	GLY	N-CA-C	5.92	120.97	114.40
1	A	584	ARG	CD-NE-CZ	5.75	132.45	124.40
1	A	549	LYS	CG-CD-CE	-5.74	98.09	111.30
1	A	561	ARG	NH1-CZ-NH2	5.65	126.65	119.30
1	C	573	ASN	OD1-CG-ND2	5.64	128.24	122.60
1	A	840	THR	CA-C-N	-5.62	112.75	120.28
1	A	840	THR	C-N-CA	-5.62	112.75	120.28
1	A	503	ARG	CA-C-N	5.58	128.22	120.29
1	A	503	ARG	C-N-CA	5.58	128.22	120.29
1	C	717	LYS	CG-CD-CE	5.58	124.13	111.30
1	D	845	GLU	CG-CD-OE1	5.50	131.06	118.40
1	F	793	LYS	CB-CG-CD	5.42	123.77	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	470	ASP	N-CA-C	-5.39	106.47	112.57
1	D	805	MET	CG-SD-CE	-5.37	89.09	100.90
1	A	580	PRO	CA-CB-CG	-5.28	94.47	104.50
1	D	785	LYS	CB-CA-C	5.26	119.48	111.80
1	F	512	ARG	NH1-CZ-NH2	5.25	126.13	119.30
1	D	805	MET	CB-CG-SD	5.24	128.41	112.70
1	F	855	TYR	N-CA-CB	-5.23	102.45	111.55
1	B	595	ARG	NE-CZ-NH2	5.22	123.90	119.20
1	D	811	LYS	CG-CD-CE	-5.21	99.33	111.30
1	A	595	ARG	NH1-CZ-NH2	-5.20	112.54	119.30
1	E	648	HIS	CA-CB-CG	5.17	118.97	113.80
1	C	531	ASP	CA-CB-CG	5.15	117.75	112.60
1	D	381	LYS	CG-CD-CE	5.11	123.06	111.30
1	D	853	ASN	OD1-CG-ND2	5.07	127.67	122.60
1	A	830	LEU	CA-C-N	5.05	124.96	119.76
1	A	830	LEU	C-N-CA	5.05	124.96	119.76

There are no chirality outliers.

All (32) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	374	ARG	Sidechain
1	A	382	ARG	Sidechain
1	A	384	ARG	Sidechain
1	A	390	ARG	Sidechain
1	A	460	ARG	Sidechain
1	A	503	ARG	Sidechain
1	A	508	ARG	Sidechain
1	A	581	ARG	Sidechain
1	A	584	ARG	Sidechain
1	A	595	ARG	Sidechain
1	A	631	ARG	Sidechain
1	B	390	ARG	Sidechain
1	B	584	ARG	Sidechain
1	B	595	ARG	Sidechain
1	C	460	ARG	Sidechain
1	C	508	ARG	Sidechain
1	C	561	ARG	Sidechain
1	C	581	ARG	Sidechain
1	C	655	ARG	Sidechain
1	D	460	ARG	Sidechain
1	D	583	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	D	595	ARG	Sidechain
1	E	503	ARG	Sidechain
1	E	545	ARG	Sidechain
1	E	581	ARG	Sidechain
1	E	595	ARG	Sidechain
1	E	676	ARG	Sidechain
1	F	419	ARG	Sidechain
1	F	503	ARG	Sidechain
1	F	512	ARG	Sidechain
1	F	655	ARG	Sidechain
1	F	732	ARG	Sidechain

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	3978	0	4096	47	0
1	B	3978	0	4096	27	0
1	C	3978	0	4096	34	0
1	D	3955	0	4077	35	0
1	E	3969	0	4088	29	0
1	F	3978	0	4096	28	0
2	A	92	0	87	12	0
2	B	92	0	87	10	0
2	C	92	0	87	11	0
2	D	92	0	87	12	0
2	E	92	0	87	13	0
2	F	92	0	87	12	0
3	A	18	0	24	1	0
3	B	6	0	8	0	0
3	C	24	0	32	2	0
3	D	6	0	8	0	0
3	E	6	0	8	0	0
3	F	6	0	8	0	0
4	A	16	0	24	0	0
5	A	27	0	0	3	0
5	B	18	0	0	0	0
5	C	26	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	17	0	0	0	0
5	E	6	0	0	0	0
5	F	16	0	0	0	0
All	All	24580	0	25183	251	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 5.

All (251) 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:799:LYS:N	1:A:800:PRO:HD3	1.85	0.91
1:A:457:LEU:HD22	1:A:460:ARG:HD2	1.54	0.88
1:F:462:LEU:HD23	1:F:488:ALA:HB3	1.55	0.86
1:F:462:LEU:CD2	1:F:488:ALA:HB3	2.08	0.84
1:A:752:VAL:HG23	1:A:800:PRO:CG	2.08	0.83
2:C:901:COB:H552	2:C:901:COB:H531	1.60	0.82
2:C:901:COB:H252	2:C:901:COB:H601	1.60	0.82
1:E:639:LEU:HD11	1:E:725:GLN:HG3	1.60	0.82
2:E:901:COB:H362	2:E:901:COB:H351	1.60	0.81
2:B:901:COB:H362	2:B:901:COB:H351	1.63	0.80
1:A:752:VAL:HG23	1:A:800:PRO:HG2	1.63	0.79
1:B:639:LEU:HD11	1:B:725:GLN:HG3	1.64	0.77
2:A:901:COB:H552	2:A:901:COB:H531	1.66	0.76
1:C:631:ARG:HG2	1:C:637:PRO:HD3	1.67	0.76
1:C:840:THR:HG23	1:C:843:TYR:H	1.49	0.76
2:D:901:COB:H552	2:D:901:COB:H531	1.67	0.75
1:A:651:ASP:HB3	1:A:654:LYS:HB3	1.69	0.74
1:A:752:VAL:CG2	1:A:800:PRO:HG2	2.18	0.73
2:D:901:COB:H252	2:D:901:COB:H601	1.71	0.70
1:C:631:ARG:HG2	1:C:637:PRO:CD	2.22	0.70
1:C:584:ARG:HH22	1:D:545:ARG:NH2	1.88	0.70
2:E:901:COB:H531	2:E:901:COB:H552	1.74	0.69
1:B:809:LEU:HD22	2:B:901:COB:H302	1.75	0.69
2:D:901:COB:H362	2:D:901:COB:H351	1.75	0.68
1:D:492:LEU:HD23	1:D:531:ASP:HB3	1.75	0.68
2:B:901:COB:H531	2:B:901:COB:H552	1.75	0.67
1:C:639:LEU:HD11	1:C:725:GLN:HG3	1.76	0.67
1:A:669:LYS:HD3	5:A:1011:HOH:O	1.95	0.66
2:F:901:COB:H203	2:F:901:COB:H301	1.76	0.66
2:C:901:COB:H362	2:C:901:COB:H351	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:584:ARG:HD3	5:A:1019:HOH:O	1.98	0.63
2:F:901:COB:H362	2:F:901:COB:H351	1.81	0.63
2:E:901:COB:H601	2:E:901:COB:H252	1.81	0.62
1:A:584:ARG:HG3	1:A:608:LYS:O	1.99	0.62
2:A:901:COB:H601	2:A:901:COB:H262	1.81	0.61
1:F:464:ASN:O	1:F:465:SER:HB3	2.00	0.61
1:E:650:GLU:HB2	1:E:655:ARG:HD2	1.82	0.61
2:A:901:COB:H353	2:A:901:COB:H311	1.82	0.60
1:A:757:LYS:HG2	5:A:1014:HOH:O	2.02	0.60
1:F:545:ARG:HG3	1:F:631:ARG:HH21	1.65	0.60
1:E:502:THR:OG1	1:E:505:GLU:HG3	2.01	0.60
1:C:755:THR:HB	1:C:784:ILE:HA	1.82	0.60
2:B:901:COB:H262	2:B:901:COB:H601	1.83	0.59
1:A:611:PRO:HG2	1:A:614:GLN:HG3	1.84	0.59
2:F:901:COB:H252	2:F:901:COB:H601	1.84	0.59
1:C:504:GLU:HG2	1:D:846:GLU:HG3	1.85	0.59
1:D:762:ASP:H	2:D:901:COB:H522	1.49	0.59
2:B:901:COB:H362	2:B:901:COB:C35	2.32	0.58
1:F:465:SER:OG	1:F:466:ALA:N	2.35	0.58
1:A:416:TRP:HB3	1:A:419:ARG:HG3	1.86	0.57
1:F:651:ASP:O	1:F:655:ARG:HG3	2.04	0.57
1:B:677:LYS:HB3	1:B:732:ARG:NH1	2.20	0.57
1:F:755:THR:HB	1:F:784:ILE:HA	1.85	0.57
1:A:632:LYS:HB3	1:A:635:PHE:HB3	1.87	0.56
1:E:428:LEU:HG	1:E:432:LEU:HD23	1.86	0.56
1:A:457:LEU:CD2	1:A:460:ARG:HD2	2.33	0.56
1:D:755:THR:HB	1:D:784:ILE:HA	1.87	0.55
2:E:901:COB:H362	2:E:901:COB:C35	2.34	0.55
1:B:755:THR:HB	1:B:784:ILE:HA	1.88	0.55
1:A:803:VAL:HG23	1:A:830:LEU:HD11	1.87	0.55
1:F:443:ASP:HA	1:F:464:ASN:HB3	1.89	0.55
1:E:693:LYS:HD2	1:E:694:PRO:HD2	1.88	0.54
1:A:703:LEU:HA	1:A:730:MET:HE1	1.90	0.54
3:C:903:GOL:H32	1:D:539:GLN:HA	1.90	0.54
1:D:693:LYS:HE3	1:D:695:LEU:HB2	1.90	0.54
2:F:901:COB:H492	2:F:901:COB:H533	1.89	0.54
1:A:464:ASN:O	1:A:465:SER:HB3	2.07	0.54
1:E:506:LYS:HB3	1:E:532:LEU:HD23	1.90	0.54
1:E:575:SER:OG	1:E:583:ARG:HD2	2.08	0.54
1:A:581:ARG:HD2	1:D:470:ASP:OD2	2.08	0.53
1:E:755:THR:HB	1:E:784:ILE:HA	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:762:ASP:H	2:E:901:COB:H522	1.55	0.53
1:A:575:SER:OG	1:A:583:ARG:HD2	2.08	0.53
3:C:903:GOL:C3	1:D:539:GLN:HA	2.39	0.53
1:F:462:LEU:CD2	1:F:488:ALA:CB	2.83	0.53
1:A:755:THR:HB	1:A:784:ILE:HA	1.91	0.53
1:B:631:ARG:HG2	1:B:637:PRO:CD	2.39	0.53
1:C:413:SER:HB2	1:C:443:ASP:HB3	1.91	0.53
1:E:413:SER:HB2	1:E:443:ASP:HB3	1.91	0.52
1:E:567:PHE:H	1:E:600:THR:HG22	1.73	0.52
1:C:575:SER:OG	1:C:583:ARG:HD2	2.09	0.52
1:A:383:PHE:O	1:A:384:ARG:C	2.52	0.52
1:D:413:SER:HB2	1:D:443:ASP:HB3	1.90	0.52
1:C:470:ASP:HB2	1:C:474:ARG:HG3	1.92	0.52
2:D:901:COB:H203	2:D:901:COB:H301	1.91	0.52
1:E:464:ASN:O	1:E:490:VAL:HB	2.10	0.52
2:C:901:COB:H351	2:C:901:COB:H372	1.92	0.52
1:F:462:LEU:HD22	1:F:488:ALA:HB3	1.90	0.52
1:D:575:SER:OG	1:D:583:ARG:HD2	2.09	0.52
1:D:650:GLU:HG3	1:D:654:LYS:HG2	1.92	0.52
1:E:368:LEU:HD21	1:E:591:LEU:HD11	1.92	0.51
1:B:632:LYS:HB3	1:B:635:PHE:HB3	1.92	0.51
1:C:559:ARG:HD3	1:C:598:GLY:HA3	1.93	0.51
1:E:538:THR:HB	1:E:578:LEU:CD1	2.40	0.51
1:B:428:LEU:HG	1:B:432:LEU:HD23	1.93	0.51
1:A:798:HIS:C	1:A:800:PRO:HD3	2.36	0.51
1:D:612:ILE:HD12	1:D:613:SER:N	2.26	0.51
1:E:538:THR:HB	1:E:578:LEU:HD13	1.93	0.51
1:F:612:ILE:HD12	1:F:613:SER:N	2.25	0.51
1:F:413:SER:HB2	1:F:443:ASP:HB3	1.93	0.50
2:B:901:COB:H601	2:B:901:COB:H252	1.93	0.50
1:B:575:SER:OG	1:B:583:ARG:HD2	2.11	0.50
1:D:811:LYS:HG3	1:D:812:SER:N	2.27	0.50
1:A:413:SER:HB2	1:A:443:ASP:HB3	1.93	0.50
1:C:584:ARG:NH2	1:D:545:ARG:NH2	2.59	0.50
1:D:694:PRO:HB2	1:D:741:MET:HG2	1.94	0.49
1:A:470:ASP:HB2	1:A:474:ARG:HG3	1.94	0.49
1:B:381:LYS:HG2	1:B:384:ARG:NH2	2.27	0.49
2:A:901:COB:H601	2:A:901:COB:H252	1.93	0.49
1:E:447:PRO:HG2	1:E:477:ARG:HD3	1.95	0.49
1:B:416:TRP:HB3	1:B:419:ARG:HG3	1.95	0.49
1:D:464:ASN:HA	1:D:490:VAL:HB	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:723:VAL:HG12	2:F:901:COB:H532	1.95	0.49
1:A:632:LYS:HG2	1:A:633:GLU:H	1.78	0.48
1:A:762:ASP:HB2	1:A:765:LYS:HE3	1.93	0.48
1:F:667:ARG:O	1:F:671:ARG:HG3	2.14	0.48
1:F:851:TYR:O	1:F:854:VAL:HG23	2.13	0.48
2:A:901:COB:H541	2:A:901:COB:H602	1.76	0.48
1:C:584:ARG:HH22	1:D:545:ARG:HH22	1.61	0.48
1:D:524:ARG:HB3	1:D:526:GLU:HG2	1.96	0.48
1:D:667:ARG:O	1:D:671:ARG:HG3	2.14	0.48
1:F:426:ARG:HG3	1:F:452:LEU:HD21	1.95	0.48
1:F:805:MET:HE1	1:F:819:ASN:HB2	1.96	0.48
1:C:443:ASP:HA	1:C:464:ASN:HB3	1.94	0.48
1:A:374:ARG:NH1	1:A:382:ARG:NH1	2.62	0.47
2:D:901:COB:H362	2:D:901:COB:C35	2.44	0.47
1:C:685:GLU:O	1:C:689:LYS:HG2	2.14	0.47
1:B:413:SER:HB2	1:B:443:ASP:HB3	1.96	0.47
1:D:593:GLU:O	1:D:597:ARG:HG2	2.15	0.47
2:E:901:COB:H601	2:E:901:COB:H262	1.95	0.47
1:D:652:PRO:O	1:D:656:GLU:HG2	2.14	0.47
1:F:671:ARG:NH2	1:F:678:GLN:O	2.41	0.47
1:C:581:ARG:HG3	1:C:584:ARG:NH2	2.30	0.47
1:C:667:ARG:O	1:C:671:ARG:HG3	2.15	0.47
2:C:901:COB:H621	2:C:901:COB:H18	1.49	0.46
2:E:901:COB:H301	2:E:901:COB:H203	1.97	0.46
2:C:901:COB:H552	2:C:901:COB:C53	2.41	0.46
1:C:755:THR:HG22	1:C:759:ALA:HB3	1.97	0.46
1:D:803:VAL:HG23	1:D:830:LEU:HD11	1.98	0.46
1:E:466:ALA:HB2	1:E:489:LEU:HD11	1.98	0.46
2:D:901:COB:H601	2:D:901:COB:H262	1.97	0.46
2:C:901:COB:H301	2:C:901:COB:H203	1.97	0.46
1:A:752:VAL:HG23	1:A:800:PRO:HG3	1.95	0.46
1:C:464:ASN:O	1:C:465:SER:HB2	2.16	0.46
1:C:465:SER:O	1:C:478:VAL:HG11	2.16	0.46
1:A:460:ARG:O	1:A:460:ARG:HG3	2.16	0.45
1:A:465:SER:OG	1:A:466:ALA:N	2.47	0.45
1:A:581:ARG:NH1	1:A:584:ARG:NH2	2.63	0.45
2:A:901:COB:H362	2:A:901:COB:C35	2.46	0.45
1:C:765:LYS:HB3	1:C:765:LYS:HE3	1.63	0.45
1:E:667:ARG:O	1:E:671:ARG:HG3	2.16	0.45
1:A:752:VAL:HG21	1:A:800:PRO:HG2	1.98	0.45
1:E:477:ARG:O	1:E:481:LEU:HG	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:531:ASP:O	1:F:532:LEU:C	2.58	0.45
1:F:450:MET:HE1	1:F:465:SER:OG	2.15	0.45
1:E:632:LYS:O	1:E:635:PHE:HB3	2.17	0.45
1:B:502:THR:OG1	1:B:505:GLU:HG3	2.17	0.45
2:B:901:COB:H8	2:B:901:COB:H401	1.81	0.45
1:C:477:ARG:HA	1:C:477:ARG:HD2	1.79	0.45
1:C:651:ASP:HB3	1:C:654:LYS:HB2	1.99	0.45
2:A:901:COB:HM51	2:A:901:COB:HM63	1.70	0.45
1:B:631:ARG:HG2	1:B:637:PRO:HD3	1.99	0.45
1:F:857:ALA:CB	2:F:901:COB:HM62	2.46	0.45
1:C:784:ILE:HG13	1:C:785:LYS:HG3	1.98	0.45
1:C:365:GLU:HG2	1:C:366:ALA:H	1.82	0.44
1:B:572:SER:HA	1:B:587:ASN:ND2	2.32	0.44
1:B:762:ASP:HB2	1:B:765:LYS:HE3	1.99	0.44
1:D:464:ASN:O	1:D:465:SER:HB2	2.17	0.44
2:E:901:COB:H463	2:E:901:COB:H481	1.84	0.44
1:F:806:SER:HB2	2:F:901:COB:H4B	1.99	0.44
1:B:436:LEU:O	1:B:460:ARG:NH2	2.47	0.44
2:C:901:COB:N29	2:C:901:COB:H3	2.31	0.44
1:B:636:ASP:OD2	1:B:639:LEU:HG	2.18	0.44
1:B:667:ARG:O	1:B:671:ARG:HG3	2.18	0.44
1:B:721:PRO:O	1:B:725:GLN:HG2	2.18	0.44
2:E:901:COB:H541	2:E:901:COB:H602	1.83	0.44
1:F:465:SER:O	1:F:492:LEU:HD12	2.18	0.44
1:A:534:THR:H	3:A:903:GOL:H11	1.83	0.44
1:D:806:SER:HB2	2:D:901:COB:H4B	2.00	0.44
1:F:462:LEU:HD23	1:F:488:ALA:CB	2.39	0.44
1:A:443:ASP:HA	1:A:464:ASN:HB3	1.99	0.44
2:F:901:COB:H362	2:F:901:COB:C35	2.48	0.44
2:F:901:COB:H2B	2:F:901:COB:O7R	2.18	0.44
1:A:390:ARG:HA	1:A:390:ARG:HD3	1.60	0.44
1:D:809:LEU:HD12	1:D:811:LYS:HE3	2.00	0.44
1:E:529:LEU:HD23	1:E:566:GLY:HA3	2.00	0.44
1:B:584:ARG:HG3	1:B:608:LYS:O	2.17	0.43
2:A:901:COB:H362	2:A:901:COB:H351	1.99	0.43
2:B:901:COB:H541	2:B:901:COB:H602	1.82	0.43
1:A:763:ILE:HG22	2:A:901:COB:H3P3	2.01	0.43
2:B:901:COB:H481	2:B:901:COB:H463	1.72	0.43
2:F:901:COB:N29	2:F:901:COB:H3	2.33	0.43
1:B:848:ARG:HA	1:B:851:TYR:O	2.19	0.43
1:C:365:GLU:HB2	1:C:369:PHE:CD1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:584:ARG:NH2	1:D:545:ARG:HH22	2.15	0.43
2:C:901:COB:H463	2:C:901:COB:H481	1.74	0.43
2:D:901:COB:H552	2:D:901:COB:C53	2.41	0.43
1:F:848:ARG:HA	1:F:851:TYR:O	2.18	0.43
1:C:405:GLU:HG2	1:D:633:GLU:HG2	2.01	0.43
1:E:806:SER:HB2	2:E:901:COB:H4B	2.00	0.43
2:D:901:COB:O28	2:D:901:COB:H3	2.19	0.43
1:F:784:ILE:HG13	1:F:785:LYS:HG3	2.00	0.43
2:F:901:COB:C55	2:F:901:COB:H531	2.49	0.43
2:B:901:COB:H552	2:B:901:COB:C53	2.46	0.43
1:A:632:LYS:HE3	1:A:635:PHE:HB2	2.00	0.42
1:A:667:ARG:O	1:A:671:ARG:HG3	2.19	0.42
1:E:784:ILE:HG13	1:E:785:LYS:HG3	2.01	0.42
1:B:442:VAL:HG23	1:B:461:VAL:HG13	2.00	0.42
1:C:839:LEU:HD23	1:C:839:LEU:HA	1.86	0.42
1:D:398:LEU:O	1:D:402:GLN:HG3	2.20	0.42
2:D:901:COB:H2B	2:D:901:COB:O7R	2.20	0.42
2:F:901:COB:H541	2:F:901:COB:H602	1.74	0.42
1:C:531:ASP:O	1:C:532:LEU:C	2.63	0.42
1:C:839:LEU:HB3	1:C:856:TYR:HE1	1.84	0.42
1:D:526:GLU:HG2	1:D:526:GLU:H	1.60	0.42
1:B:784:ILE:HG13	1:B:785:LYS:HG3	2.01	0.42
1:C:610:LEU:HD21	1:D:636:ASP:HB2	2.01	0.42
1:C:501:LYS:HD3	1:C:535:PHE:CE2	2.55	0.41
1:E:469:GLU:HG3	1:E:470:ASP:OD1	2.19	0.41
1:E:595:ARG:HA	1:E:599:LEU:HB3	2.01	0.41
1:E:721:PRO:O	1:E:725:GLN:HG2	2.20	0.41
1:A:784:ILE:HG13	1:A:785:LYS:HG3	2.01	0.41
1:D:581:ARG:NH2	1:D:648:HIS:O	2.53	0.41
1:A:685:GLU:OE2	1:A:689:LYS:HE2	2.20	0.41
1:A:809:LEU:HD11	2:A:901:COB:H363	2.02	0.41
1:B:429:LEU:HA	1:B:432:LEU:HG	2.02	0.41
2:C:901:COB:H362	2:C:901:COB:C35	2.49	0.41
1:E:848:ARG:HA	1:E:851:TYR:O	2.21	0.41
1:A:787:PRO:HD2	1:A:790:GLU:HG3	2.02	0.41
1:D:464:ASN:O	1:D:465:SER:CB	2.68	0.41
1:A:477:ARG:HA	1:A:477:ARG:HD2	1.95	0.41
1:B:685:GLU:O	1:B:689:LYS:HG2	2.21	0.41
1:D:612:ILE:HD12	1:D:613:SER:H	1.86	0.41
2:D:901:COB:H411	2:D:901:COB:H363	1.73	0.41
2:E:901:COB:HM63	2:E:901:COB:HM51	1.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:461:VAL:HG12	1:A:463:LEU:HD12	2.02	0.41
1:A:677:LYS:HE3	1:A:677:LYS:HB2	1.86	0.41
1:B:805:MET:HE1	1:B:819:ASN:HB2	2.02	0.41
2:A:901:COB:H363	2:A:901:COB:H411	1.82	0.40
1:B:529:LEU:HD23	1:B:566:GLY:HA3	2.03	0.40
1:F:840:THR:O	1:F:843:TYR:N	2.50	0.40
1:A:398:LEU:O	1:A:402:GLN:HG3	2.22	0.40
2:C:901:COB:HM63	2:C:901:COB:HM51	1.82	0.40
1:D:834:LEU:HB3	1:D:839:LEU:HD11	2.03	0.40
2:E:901:COB:H3	2:E:901:COB:N29	2.35	0.40
1:F:699:ASN:HA	1:F:703:LEU:HD12	2.04	0.40
1:A:710:GLY:HA3	2:A:901:COB:H602	2.04	0.40
1:E:559:ARG:HD3	1:E:598:GLY:HA3	2.02	0.40
2:E:901:COB:H552	2:E:901:COB:C53	2.49	0.40
1:C:595:ARG:HA	1:C:599:LEU:HB3	2.03	0.40
1:E:474:ARG:O	1:E:477:ARG:HG2	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	508/511 (99%)	495 (97%)	12 (2%)	1 (0%)	43	71
1	B	508/511 (99%)	491 (97%)	15 (3%)	2 (0%)	30	59
1	C	508/511 (99%)	490 (96%)	17 (3%)	1 (0%)	43	71
1	D	505/511 (99%)	493 (98%)	11 (2%)	1 (0%)	43	71
1	E	507/511 (99%)	494 (97%)	13 (3%)	0	100	100
1	F	508/511 (99%)	494 (97%)	13 (3%)	1 (0%)	43	71
All	All	3044/3066 (99%)	2957 (97%)	81 (3%)	6 (0%)	43	71

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	465	SER
1	B	465	SER
1	C	465	SER
1	D	465	SER
1	F	465	SER
1	B	493	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	414/414 (100%)	411 (99%)	3 (1%)	76	79
1	B	414/414 (100%)	412 (100%)	2 (0%)	81	82
1	C	414/414 (100%)	412 (100%)	2 (0%)	81	82
1	D	412/414 (100%)	409 (99%)	3 (1%)	76	79
1	E	413/414 (100%)	410 (99%)	3 (1%)	76	79
1	F	414/414 (100%)	412 (100%)	2 (0%)	81	82
All	All	2481/2484 (100%)	2466 (99%)	15 (1%)	78	81

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	580	PRO
1	A	581	ARG
1	A	841	ARG
1	B	480	SER
1	B	748	LYS
1	C	724	LEU
1	C	765	LYS
1	D	515	GLU
1	D	610	LEU
1	D	811	LYS
1	E	455	LYS

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Mol	Chain	Res	Type
1	E	676	ARG
1	E	809	LEU
1	F	469	GLU
1	F	633	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	467	ASN
1	B	464	ASN
1	B	587	ASN
1	B	719	GLN
1	B	853	ASN
1	C	819	ASN
1	C	853	ASN
1	E	464	ASN
1	E	853	ASN
1	F	364	GLN
1	F	464	ASN

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)

19 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	$\#\|Z\| > 2$
3	GOL	D	902	-	5,5,5	0.93	0	5,5,5	0.85	0
2	COB	F	901	1	93,102,102	1.17	2 (2%)	149,170,170	1.42	14 (9%)
3	GOL	C	905	-	5,5,5	0.99	0	5,5,5	1.15	0
3	GOL	A	904	-	5,5,5	1.01	0	5,5,5	0.71	0
3	GOL	C	904	-	5,5,5	1.18	0	5,5,5	0.88	0
3	GOL	C	903	-	5,5,5	0.30	0	5,5,5	0.60	0
2	COB	C	901	1	93,102,102	1.40	4 (4%)	149,170,170	1.49	17 (11%)
3	GOL	F	902	-	5,5,5	0.38	0	5,5,5	0.64	0
4	TRS	A	905	-	7,7,7	1.71	1 (14%)	9,9,9	1.45	2 (22%)
3	GOL	A	903	-	5,5,5	1.02	0	5,5,5	1.09	0
4	TRS	A	906	-	7,7,7	1.77	2 (28%)	9,9,9	1.19	1 (11%)
3	GOL	A	902	-	5,5,5	0.38	0	5,5,5	0.36	0
2	COB	B	901	1	93,102,102	1.23	2 (2%)	149,170,170	1.44	18 (12%)
2	COB	D	901	1	93,102,102	1.33	4 (4%)	149,170,170	1.55	22 (14%)
2	COB	A	901	1	93,102,102	1.27	3 (3%)	149,170,170	1.55	19 (12%)
3	GOL	C	902	-	5,5,5	0.14	0	5,5,5	0.50	0
3	GOL	E	902	-	5,5,5	0.62	0	5,5,5	1.15	1 (20%)
2	COB	E	901	1	93,102,102	1.28	4 (4%)	149,170,170	1.45	17 (11%)
3	GOL	B	902	-	5,5,5	0.21	0	5,5,5	0.23	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	D	902	-	-	3/4/4/4	-
2	COB	F	901	1	1/1/36/38	12/56/231/231	0/3/11/11
3	GOL	C	905	-	-	4/4/4/4	-
3	GOL	A	904	-	-	0/4/4/4	-
3	GOL	C	904	-	-	3/4/4/4	-
3	GOL	C	903	-	-	0/4/4/4	-
2	COB	C	901	1	1/1/36/38	16/56/231/231	0/3/11/11
3	GOL	F	902	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	TRS	A	905	-	-	9/9/9/9	-
3	GOL	A	903	-	-	0/4/4/4	-
4	TRS	A	906	-	-	3/9/9/9	-
3	GOL	A	902	-	-	1/4/4/4	-
2	COB	B	901	1	1/1/36/38	20/56/231/231	0/3/11/11
2	COB	D	901	1	1/1/36/38	14/56/231/231	0/3/11/11
2	COB	A	901	1	1/1/36/38	19/56/231/231	0/3/11/11
3	GOL	C	902	-	-	2/4/4/4	-
3	GOL	E	902	-	-	2/4/4/4	-
2	COB	E	901	1	1/1/36/38	19/56/231/231	0/3/11/11
3	GOL	B	902	-	-	4/4/4/4	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	901	COB	C19-N24	-10.28	1.27	1.49
2	C	901	COB	C19-N24	-10.25	1.27	1.49
2	A	901	COB	C19-N24	-10.14	1.28	1.49
2	B	901	COB	C19-N24	-10.05	1.28	1.49
2	E	901	COB	C19-N24	-9.95	1.28	1.49
2	F	901	COB	C19-N24	-9.69	1.29	1.49
2	A	901	COB	C54-C17	-3.40	1.49	1.54
2	D	901	COB	C18-C19	-2.75	1.48	1.54
4	A	906	TRS	C-N	2.70	1.58	1.49
2	C	901	COB	C18-C19	-2.69	1.49	1.54
4	A	905	TRS	C-N	2.67	1.58	1.49
2	E	901	COB	C18-C19	-2.59	1.49	1.54
2	D	901	COB	C14-C15	-2.52	1.32	1.43
2	B	901	COB	C18-C19	-2.51	1.49	1.54
2	E	901	COB	C36-C7	-2.30	1.50	1.54
2	A	901	COB	C18-C19	-2.20	1.50	1.54
2	F	901	COB	C18-C19	-2.18	1.50	1.54
4	A	906	TRS	C2-C	2.15	1.58	1.53
2	C	901	COB	C14-N23	2.09	1.41	1.33
2	C	901	COB	C36-C7	-2.08	1.51	1.54
2	E	901	COB	C54-C17	-2.07	1.51	1.54
2	D	901	COB	C4-C5	-2.05	1.34	1.43

All (111) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	COB	C1-C19-N24	8.58	119.98	106.29
2	F	901	COB	C1-C19-N24	8.50	119.86	106.29
2	B	901	COB	C1-C19-N24	8.29	119.53	106.29
2	D	901	COB	C1-C19-N24	8.05	119.14	106.29
2	E	901	COB	C1-C19-N24	7.99	119.03	106.29
2	C	901	COB	C1-C19-N24	7.60	118.41	106.29
2	E	901	COB	C18-C19-N24	6.60	112.42	101.92
2	F	901	COB	C18-C19-N24	6.59	112.41	101.92
2	C	901	COB	C18-C19-N24	6.51	112.28	101.92
2	B	901	COB	C18-C19-N24	6.18	111.76	101.92
2	A	901	COB	C18-C19-N24	5.78	111.11	101.92
2	D	901	COB	C18-C19-N24	5.58	110.80	101.92
2	F	901	COB	C26-C2-C1	4.77	117.37	110.00
2	A	901	COB	C9B-C8B-N1B	4.62	107.35	105.30
2	A	901	COB	C1-C19-C18	4.56	128.60	121.75
2	D	901	COB	C1-C19-C18	4.55	128.59	121.75
2	C	901	COB	C1-C19-C18	4.13	127.96	121.75
2	B	901	COB	C1-C19-C18	4.09	127.91	121.75
2	E	901	COB	C9B-C8B-N1B	3.80	106.99	105.30
2	C	901	COB	C41-C8-C9	3.78	117.78	111.19
2	C	901	COB	C2-C3-C4	-3.73	97.45	101.64
2	C	901	COB	C26-C2-C1	3.70	115.72	110.00
2	F	901	COB	C1-C19-C18	3.69	127.30	121.75
2	E	901	COB	C1-C19-C18	3.62	127.19	121.75
2	A	901	COB	C19-C1-N21	3.58	106.20	101.69
2	E	901	COB	C54-C17-C18	-3.48	107.99	112.99
2	B	901	COB	C35-C5-C6	-3.31	117.08	122.41
2	D	901	COB	C2-C3-C4	-3.20	98.05	101.64
2	B	901	COB	C9B-C8B-N1B	3.08	106.67	105.30
2	C	901	COB	O39-C38-C37	-3.05	112.55	121.98
2	B	901	COB	C19-C1-N21	3.02	105.50	101.69
2	B	901	COB	C1-N21-C4	3.01	114.37	109.48
2	F	901	COB	C1-N21-C4	3.00	114.36	109.48
2	F	901	COB	C19-C1-N21	2.99	105.46	101.69
2	E	901	COB	C26-C2-C1	2.98	114.61	110.00
2	E	901	COB	C2-C3-C4	-2.96	98.31	101.64
2	E	901	COB	O34-C32-N33	-2.93	114.70	122.53
2	D	901	COB	O39-C38-C37	2.92	131.04	121.98
2	A	901	COB	C26-C2-C1	2.89	114.48	110.00
4	A	905	TRS	O1-C1-C	2.89	118.92	110.88
2	A	901	COB	C5M-C5B-C4B	2.87	124.63	119.57
2	C	901	COB	C10-C9-N22	-2.86	122.47	125.74
2	A	901	COB	C35-C5-C6	-2.85	117.82	122.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	901	COB	C9-C10-C11	-2.85	121.65	125.84
2	C	901	COB	C54-C17-C18	-2.82	108.94	112.99
2	D	901	COB	C1-N21-C4	2.79	114.01	109.48
2	A	901	COB	C54-C17-C18	-2.71	109.10	112.99
2	A	901	COB	C10-C9-N22	-2.69	122.67	125.74
2	D	901	COB	O51-C50-C49	-2.67	112.97	121.04
2	D	901	COB	C19-C1-N21	2.66	105.04	101.69
2	A	901	COB	C1-N21-C4	2.63	113.76	109.48
2	D	901	COB	C26-C2-C1	2.63	114.07	110.00
2	C	901	COB	C19-C1-N21	2.61	104.98	101.69
2	B	901	COB	C26-C2-C1	2.61	114.04	110.00
2	D	901	COB	C1-C2-C3	2.60	104.87	101.60
2	E	901	COB	C1-N21-C4	2.59	113.69	109.48
2	A	901	COB	C5M-C5B-C6B	-2.58	115.48	120.76
2	D	901	COB	C42-C43-N45	2.57	124.71	116.49
2	E	901	COB	O51-C50-C49	-2.56	113.30	121.04
2	D	901	COB	C54-C17-C18	-2.52	109.38	112.99
2	E	901	COB	C49-C50-N52	2.48	124.45	116.49
2	A	901	COB	C47-C12-C46	-2.46	105.33	109.41
2	F	901	COB	C13-C12-C11	-2.45	97.58	100.86
2	E	901	COB	C37-C7-C6	2.45	114.69	107.11
2	C	901	COB	O39-C38-N40	2.45	129.07	122.53
2	D	901	COB	C36-C7-C37	-2.44	106.59	110.74
2	B	901	COB	C3-C4-C5	2.42	127.88	123.82
2	D	901	COB	O3-C2P-C1P	2.41	111.71	106.94
2	E	901	COB	C19-C1-N21	2.39	104.69	101.69
2	B	901	COB	O39-C38-C37	-2.37	114.63	121.98
2	F	901	COB	C54-C17-C18	-2.36	109.60	112.99
2	B	901	COB	C26-C2-C3	-2.34	103.33	107.42
2	F	901	COB	C25-C2-C1	-2.33	110.27	113.75
2	F	901	COB	C3-C4-C5	2.32	127.70	123.82
2	F	901	COB	C48-C13-C14	2.31	115.09	109.66
2	B	901	COB	C42-C43-N45	2.31	123.90	116.49
2	C	901	COB	C12-C13-C14	-2.30	98.86	101.86
2	A	901	COB	C8B-C9B-N3B	-2.30	107.51	110.00
2	A	901	COB	C3-C4-C5	2.30	127.66	123.82
2	C	901	COB	C31-C32-N33	-2.29	109.16	116.49
2	A	901	COB	C5-C4-N21	-2.28	120.52	124.19
2	D	901	COB	O58-C57-N59	2.27	127.48	123.03
2	D	901	COB	C49-C50-N52	2.25	123.71	116.49
2	E	901	COB	C41-C8-C9	2.25	115.11	111.19
2	B	901	COB	C1-C2-C3	2.25	104.43	101.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	901	COB	C2-C3-C4	-2.23	99.13	101.64
2	F	901	COB	C3-C4-N21	-2.23	109.21	111.98
2	A	901	COB	C2-C1-C19	2.23	122.62	118.77
2	A	901	COB	O58-C57-C56	-2.22	118.00	122.02
2	D	901	COB	C20-C1-C2	-2.21	109.64	113.28
2	B	901	COB	C2-C1-C19	2.18	122.53	118.77
2	C	901	COB	C20-C1-C2	-2.18	109.69	113.28
2	C	901	COB	O34-C32-C31	2.18	127.61	121.04
2	D	901	COB	O44-C43-C42	-2.18	114.47	121.04
2	B	901	COB	C2R-C1R-N1B	2.16	118.67	113.30
2	E	901	COB	C7B-C8B-C9B	-2.15	119.89	122.47
2	F	901	COB	C9B-C8B-N1B	2.15	106.26	105.30
2	B	901	COB	C7-C37-C38	2.12	120.55	114.28
2	D	901	COB	C25-C2-C1	-2.10	110.61	113.75
2	C	901	COB	C1-C2-C3	2.09	104.23	101.60
4	A	905	TRS	C2-C-C1	-2.09	105.08	110.66
2	B	901	COB	C49-C50-N52	-2.09	109.80	116.49
2	D	901	COB	C37-C7-C6	2.09	113.57	107.11
2	B	901	COB	C5-C4-N21	-2.07	120.86	124.19
2	E	901	COB	C2-C1-C19	2.06	122.32	118.77
2	C	901	COB	C2-C1-C19	2.05	122.30	118.77
2	E	901	COB	C31-C32-N33	2.05	123.06	116.49
3	E	902	GOL	O2-C2-C1	-2.04	100.73	109.18
4	A	906	TRS	O2-C2-C	2.02	116.52	110.88
2	D	901	COB	O58-C57-C56	-2.02	118.36	122.02
2	D	901	COB	C26-C2-C3	-2.01	103.91	107.42

All (6) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	901	COB	C19
2	B	901	COB	C19
2	C	901	COB	C19
2	D	901	COB	C19
2	E	901	COB	C19
2	F	901	COB	C19

All (133) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901	COB	C38-C37-C7-C6
2	A	901	COB	C38-C37-C7-C8

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Mol	Chain	Res	Type	Atoms
2	B	901	COB	C2-C3-C30-C31
2	B	901	COB	C4-C3-C30-C31
2	B	901	COB	C14-C13-C48-C49
2	C	901	COB	C18-C60-C61-O63
2	C	901	COB	C18-C60-C61-N62
2	F	901	COB	C16-C17-C55-C56
2	F	901	COB	C54-C17-C55-C56
2	F	901	COB	C18-C17-C55-C56
2	F	901	COB	C2R-C1R-N1B-C2B
3	C	904	GOL	O1-C1-C2-C3
3	C	904	GOL	C1-C2-C3-O3
3	C	905	GOL	O1-C1-C2-C3
3	D	902	GOL	O1-C1-C2-C3
4	A	905	TRS	N-C-C1-O1
4	A	905	TRS	C1-C-C2-O2
4	A	905	TRS	C3-C-C2-O2
4	A	905	TRS	N-C-C2-O2
2	A	901	COB	C4-C3-C30-C31
2	C	901	COB	C42-C41-C8-C9
2	D	901	COB	C42-C41-C8-C9
2	D	901	COB	C14-C13-C48-C49
2	E	901	COB	C42-C41-C8-C9
2	F	901	COB	C42-C41-C8-C9
2	B	901	COB	C7-C37-C38-O39
2	B	901	COB	C7-C37-C38-N40
3	C	904	GOL	O2-C2-C3-O3
2	B	901	COB	C12-C13-C48-C49
2	D	901	COB	C12-C13-C48-C49
2	A	901	COB	C2R-C1R-N1B-C2B
2	A	901	COB	C18-C17-C55-C56
2	B	901	COB	C18-C17-C55-C56
2	D	901	COB	C18-C17-C55-C56
2	D	901	COB	C42-C41-C8-C7
2	A	901	COB	C16-C17-C55-C56
2	B	901	COB	C16-C17-C55-C56
2	D	901	COB	C16-C17-C55-C56
2	E	901	COB	C14-C13-C48-C49
3	B	902	GOL	O1-C1-C2-C3
3	B	902	GOL	C1-C2-C3-O3
3	C	902	GOL	C1-C2-C3-O3
3	C	905	GOL	C1-C2-C3-O3
3	F	902	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
2	A	901	COB	C2R-C1R-N1B-C8B
2	F	901	COB	C2R-C1R-N1B-C8B
3	C	905	GOL	O2-C2-C3-O3
3	D	902	GOL	O1-C1-C2-O2
2	F	901	COB	C41-C42-C43-N45
2	D	901	COB	C2R-C1R-N1B-C2B
4	A	905	TRS	C3-C-C1-O1
4	A	905	TRS	C1-C-C3-O3
4	A	906	TRS	C2-C-C1-O1
2	E	901	COB	C41-C42-C43-O44
2	F	901	COB	C41-C42-C43-O44
2	B	901	COB	O6R-C4R-C5R-O8R
2	A	901	COB	C7-C37-C38-N40
2	A	901	COB	C2-C3-C30-C31
2	C	901	COB	C12-C13-C48-C49
2	E	901	COB	C12-C13-C48-C49
3	B	902	GOL	O2-C2-C3-O3
2	E	901	COB	C16-C17-C55-C56
2	A	901	COB	C41-C42-C43-N45
2	E	901	COB	C41-C42-C43-N45
2	B	901	COB	C3R-C4R-C5R-O8R
2	C	901	COB	C14-C13-C48-C49
3	C	905	GOL	O1-C1-C2-O2
3	E	902	GOL	O1-C1-C2-O2
2	B	901	COB	C41-C42-C43-O44
2	A	901	COB	C42-C41-C8-C7
2	B	901	COB	C2R-C1R-N1B-C2B
4	A	905	TRS	C2-C-C1-O1
4	A	905	TRS	N-C-C3-O3
4	A	906	TRS	N-C-C1-O1
2	B	901	COB	C41-C42-C43-N45
2	E	901	COB	C18-C17-C55-C56
2	C	901	COB	C2-C3-C30-C31
3	C	902	GOL	O2-C2-C3-O3
3	A	902	GOL	O1-C1-C2-C3
2	A	901	COB	O6R-C1R-N1B-C8B
2	F	901	COB	O6R-C1R-N1B-C8B
2	A	901	COB	C41-C42-C43-O44
2	C	901	COB	C2R-C1R-N1B-C2B
2	E	901	COB	C2R-C1R-N1B-C2B
2	D	901	COB	O6R-C4R-C5R-O8R
3	B	902	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	A	901	COB	C7-C37-C38-O39
4	A	906	TRS	C3-C-C1-O1
2	E	901	COB	C30-C31-C32-O34
2	E	901	COB	C18-C60-C61-O63
2	B	901	COB	C42-C41-C8-C7
2	C	901	COB	C7-C37-C38-O39
3	F	902	GOL	O1-C1-C2-O2
3	D	902	GOL	C1-C2-C3-O3
3	E	902	GOL	O1-C1-C2-C3
2	A	901	COB	C17-C18-C60-C61
2	B	901	COB	C17-C18-C60-C61
2	F	901	COB	C17-C18-C60-C61
2	B	901	COB	C2R-C1R-N1B-C8B
2	B	901	COB	O6R-C1R-N1B-C8B
2	C	901	COB	C2R-C1R-N1B-C8B
2	C	901	COB	O6R-C1R-N1B-C8B
2	D	901	COB	C2R-C1R-N1B-C8B
2	D	901	COB	O6R-C1R-N1B-C8B
2	E	901	COB	C2R-C1R-N1B-C8B
2	E	901	COB	O6R-C1R-N1B-C8B
2	C	901	COB	C41-C42-C43-N45
2	C	901	COB	C4-C3-C30-C31
2	C	901	COB	N59-C1P-C2P-O3
2	E	901	COB	C2-C3-C30-C31
2	A	901	COB	C42-C41-C8-C9
2	D	901	COB	C3R-C4R-C5R-O8R
4	A	905	TRS	C2-C-C3-O3
2	A	901	COB	C19-C18-C60-C61
2	B	901	COB	C19-C18-C60-C61
2	F	901	COB	C19-C18-C60-C61
2	E	901	COB	O6R-C1R-N1B-C2B
2	D	901	COB	C41-C42-C43-N45
2	C	901	COB	N59-C1P-C2P-C3P
2	C	901	COB	C41-C42-C43-O44
2	E	901	COB	C30-C31-C32-N33
2	A	901	COB	O6R-C4R-C5R-O8R
2	E	901	COB	O6R-C4R-C5R-O8R
2	A	901	COB	C38-C37-C7-C36
2	C	901	COB	O6R-C1R-N1B-C2B
2	E	901	COB	C17-C18-C60-C61
2	B	901	COB	O6R-C1R-N1B-C2B
2	D	901	COB	C41-C42-C43-O44

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Mol	Chain	Res	Type	Atoms
2	E	901	COB	C4-C3-C30-C31
2	B	901	COB	C18-C60-C61-O63
2	E	901	COB	C18-C60-C61-N62
2	D	901	COB	O6R-C1R-N1B-C2B
2	F	901	COB	C3-C30-C31-C32

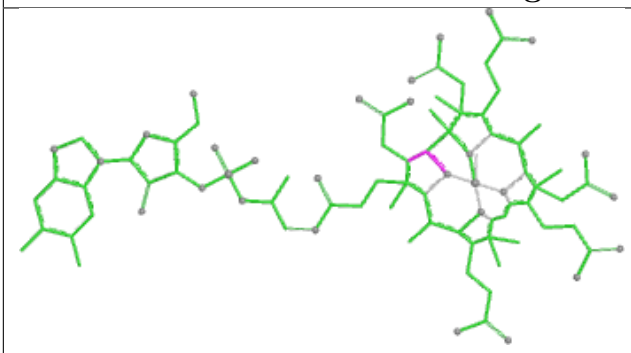
There are no ring outliers.

8 monomers are involved in 73 short contacts:

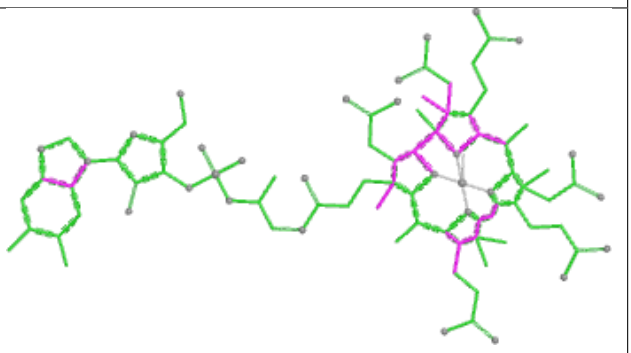
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	901	COB	12	0
3	C	903	GOL	2	0
2	C	901	COB	11	0
3	A	903	GOL	1	0
2	B	901	COB	10	0
2	D	901	COB	12	0
2	A	901	COB	12	0
2	E	901	COB	13	0

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 then 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.

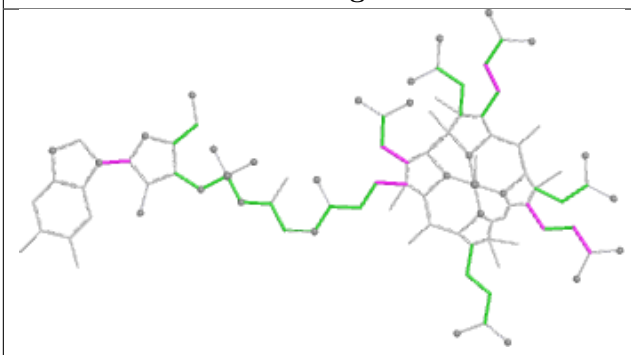
Ligand COB F 901



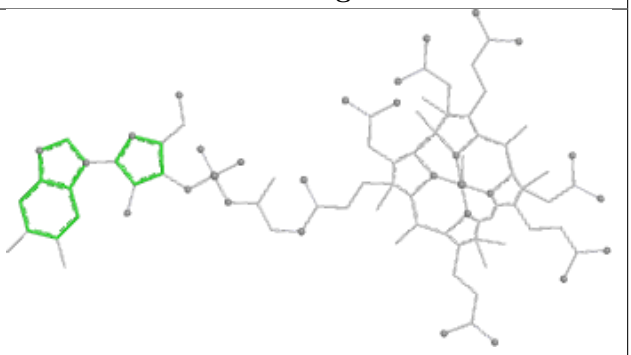
Bond lengths



Bond angles

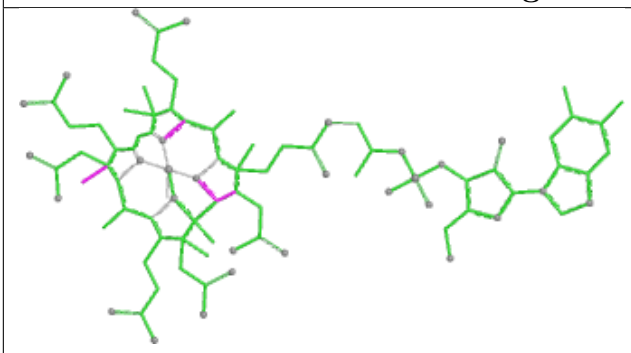


Torsions

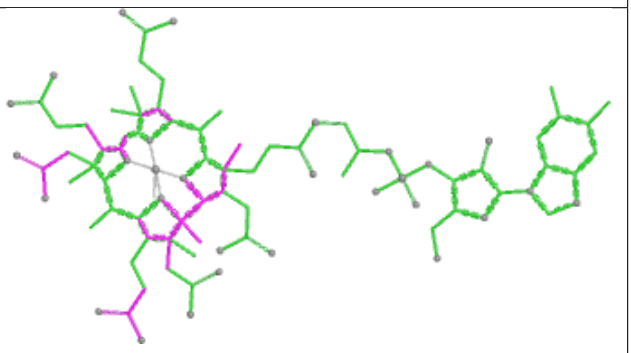


Rings

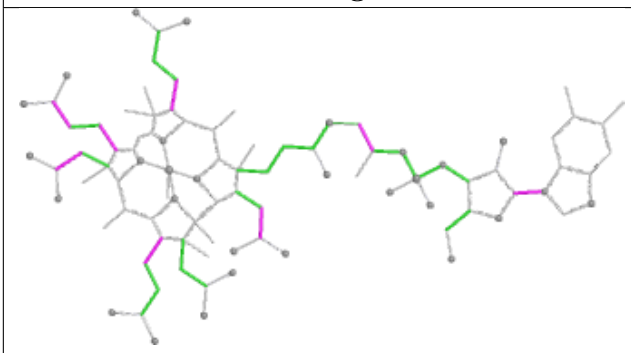
Ligand COB C 901



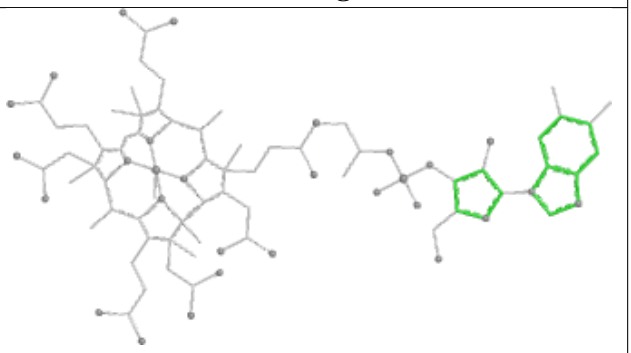
Bond lengths



Bond angles

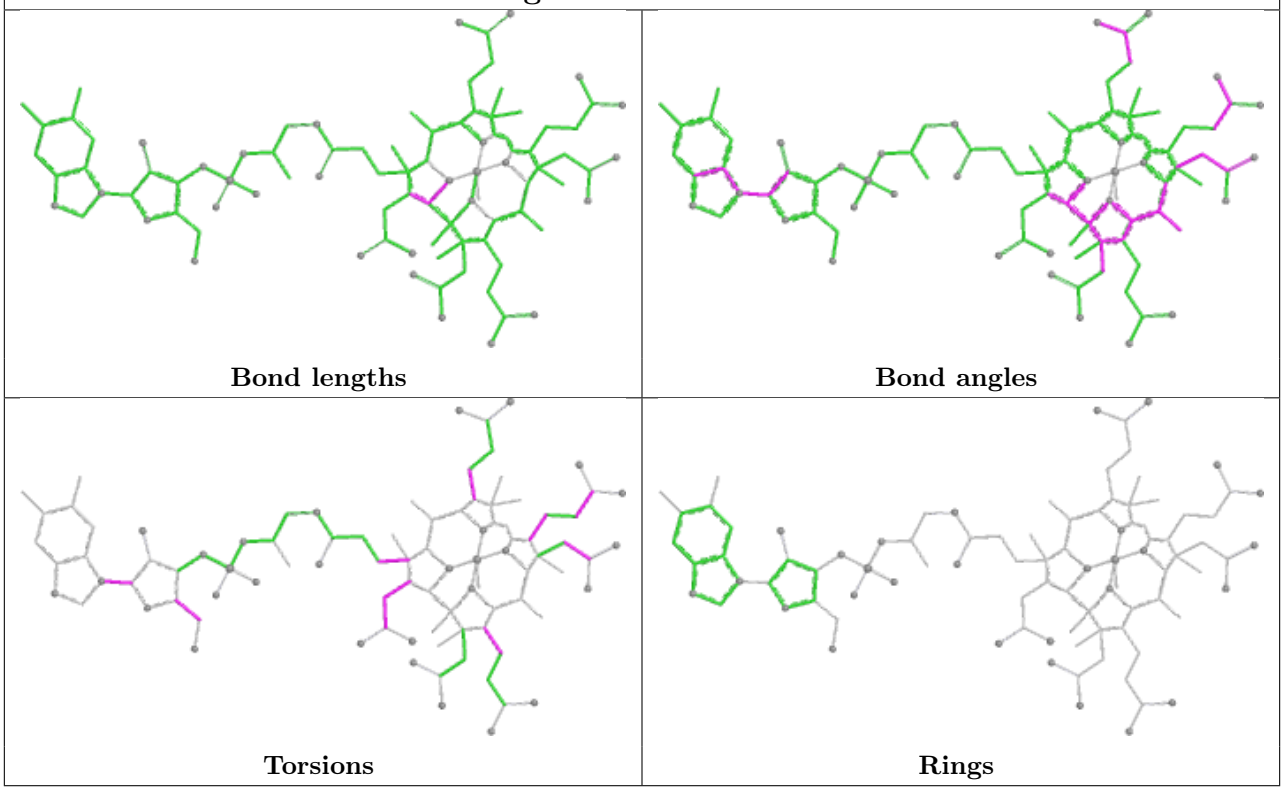


Torsions

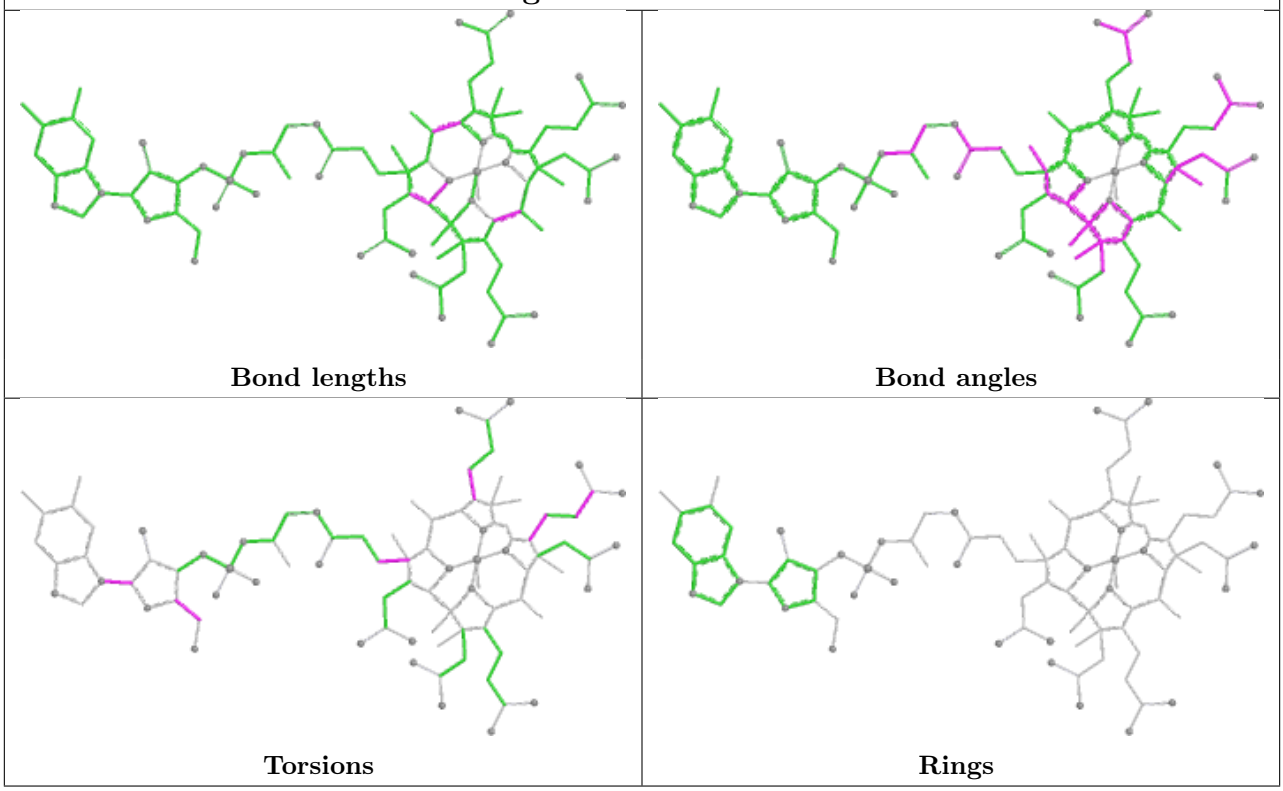


Rings

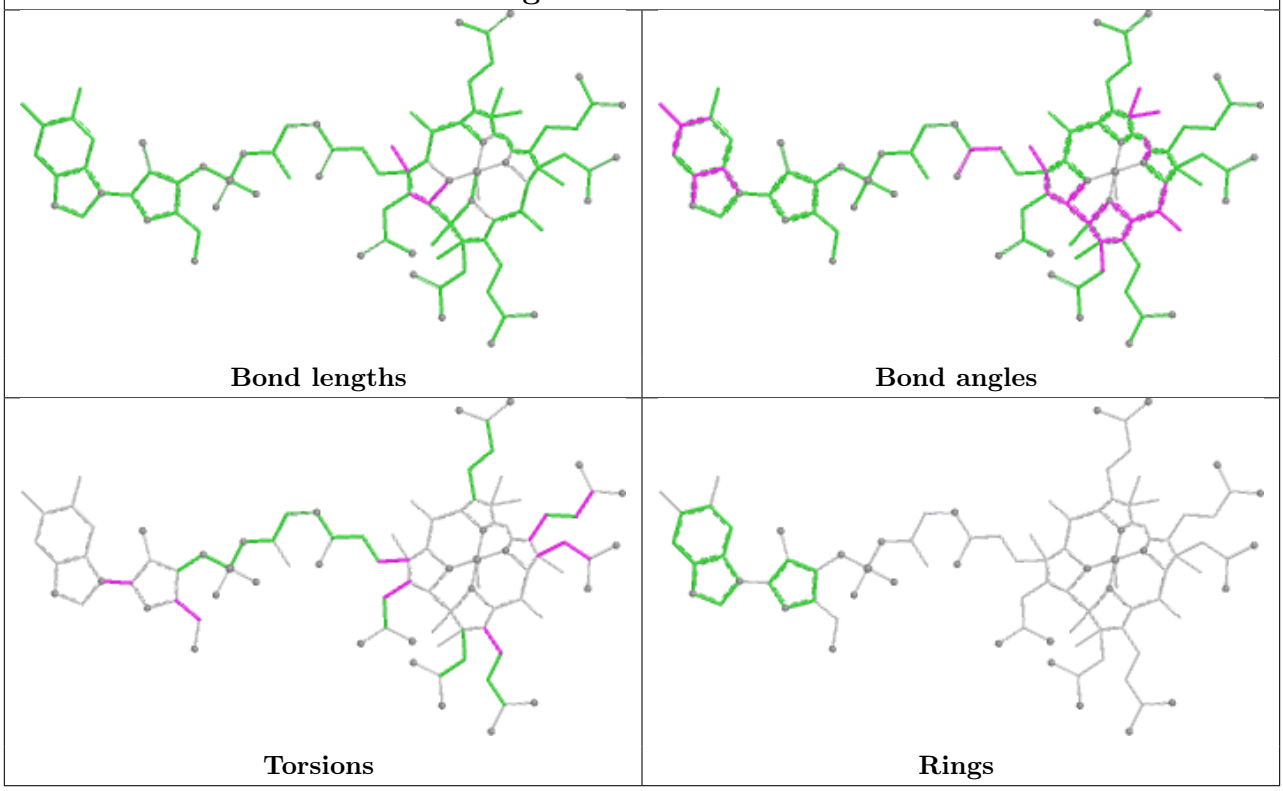
Ligand COB B 901



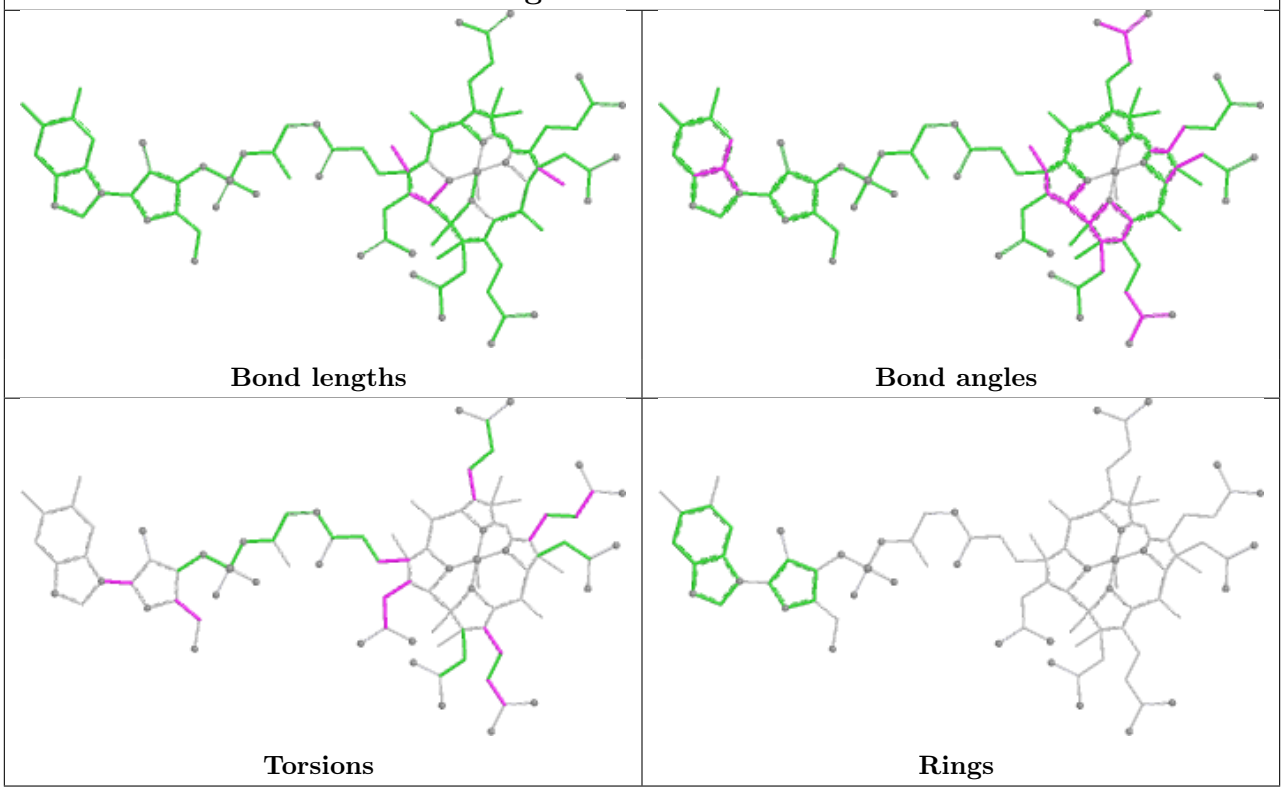
Ligand COB D 901



Ligand COB A 901



Ligand COB E 901



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 (i)

6.1 Protein, DNA and RNA chains (i)

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	510/511 (99%)	-0.44	0	100	100	67, 107, 186, 248	0
1	B	510/511 (99%)	-0.56	2 (0%)	88	80	72, 106, 170, 207	0
1	C	510/511 (99%)	-0.60	1 (0%)	91	87	67, 101, 153, 225	0
1	D	507/511 (99%)	-0.64	0	100	100	69, 108, 157, 191	0
1	E	509/511 (99%)	-0.43	2 (0%)	88	80	90, 147, 199, 233	0
1	F	510/511 (99%)	-0.55	1 (0%)	91	87	81, 118, 159, 254	0
All	All	3056/3066 (99%)	-0.54	6 (0%)	91	87	67, 113, 180, 254	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	647	ALA	3.2
1	C	648	HIS	3.0
1	B	695	LEU	2.9
1	B	715	ALA	2.6
1	E	830	LEU	2.3
1	F	535	PHE	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)

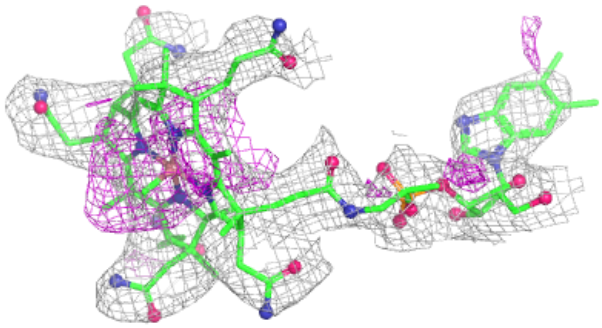
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

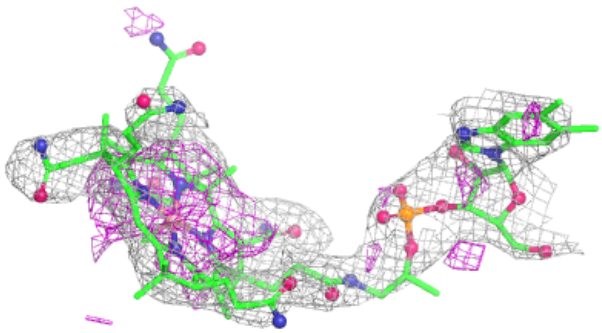
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GOL	F	902	6/6	0.81	0.14	156,168,191,199	0
3	GOL	A	904	6/6	0.82	0.12	82,101,110,131	0
3	GOL	D	902	6/6	0.83	0.14	101,105,116,120	0
3	GOL	C	905	6/6	0.85	0.16	69,93,121,136	0
3	GOL	C	902	6/6	0.86	0.16	149,177,195,208	0
3	GOL	A	903	6/6	0.88	0.16	79,100,123,136	0
4	TRS	A	905	8/8	0.88	0.13	100,137,157,171	0
3	GOL	C	904	6/6	0.89	0.16	91,106,114,124	0
3	GOL	E	902	6/6	0.89	0.23	122,130,137,148	0
4	TRS	A	906	8/8	0.91	0.13	78,99,139,151	0
3	GOL	C	903	6/6	0.93	0.14	110,135,143,145	0
3	GOL	A	902	6/6	0.94	0.14	104,110,128,134	0
2	COB	A	901	92/92	0.94	0.11	83,134,176,220	0
3	GOL	B	902	6/6	0.96	0.09	75,111,133,150	0
2	COB	F	901	92/92	0.97	0.09	66,93,117,153	0
2	COB	B	901	92/92	0.97	0.08	78,113,140,168	0
2	COB	E	901	92/92	0.98	0.07	73,103,138,165	0
2	COB	C	901	92/92	0.98	0.07	53,75,105,135	0
2	COB	D	901	92/92	0.98	0.07	60,82,115,133	0

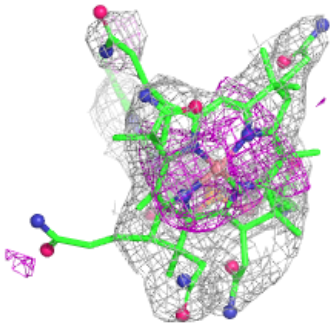
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around COB A 901:

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

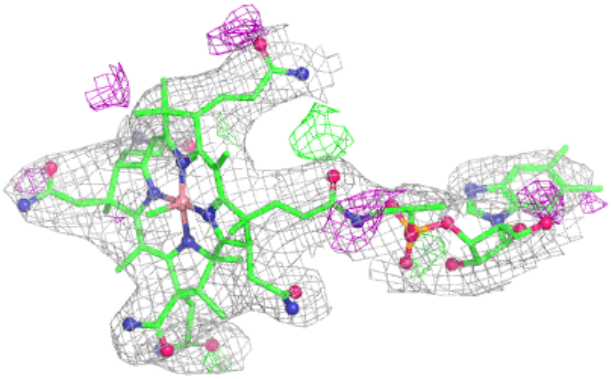


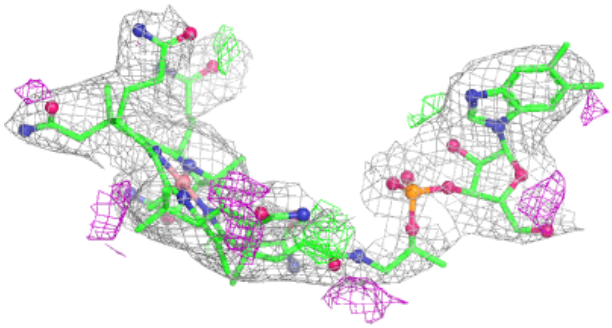


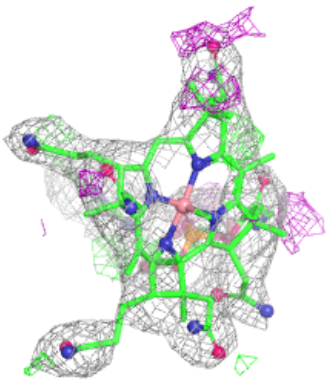


Electron density around COB F 901:

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

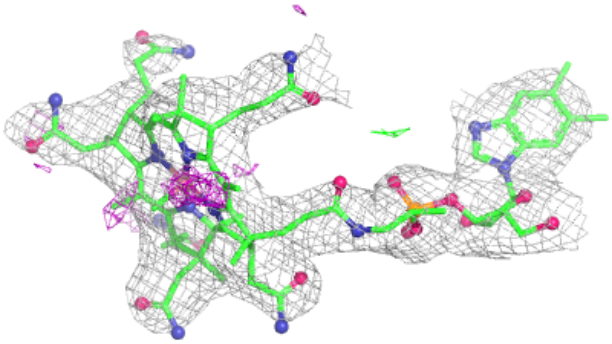


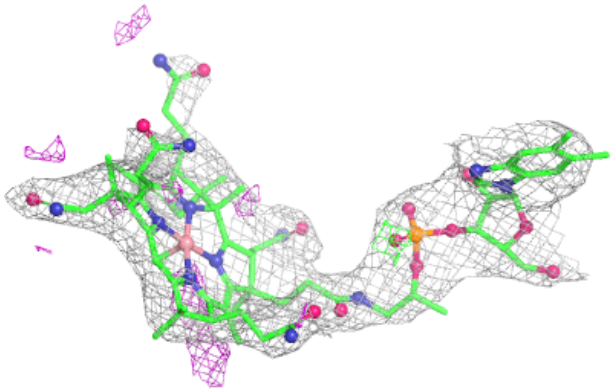


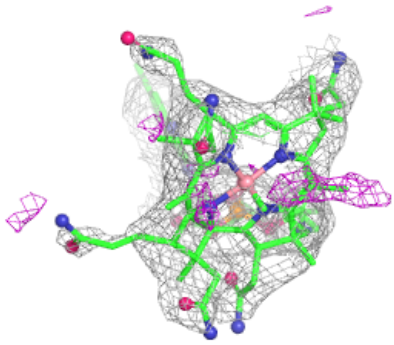


Electron density around COB B 901:

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

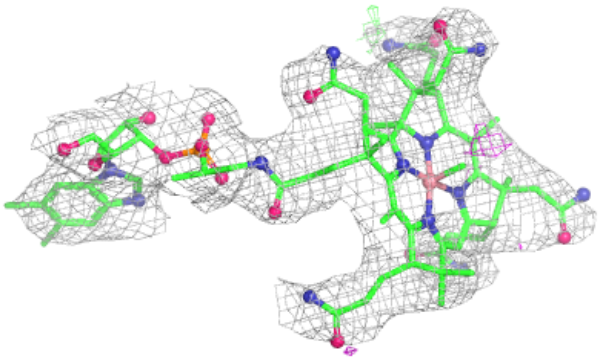


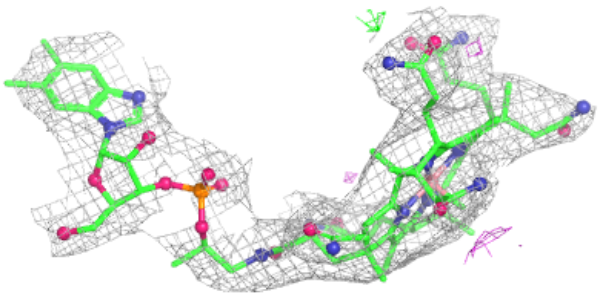


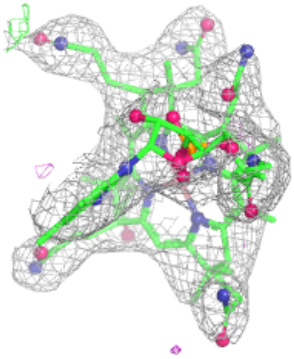


Electron density around COB E 901:

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







Electron density around COB C 901:

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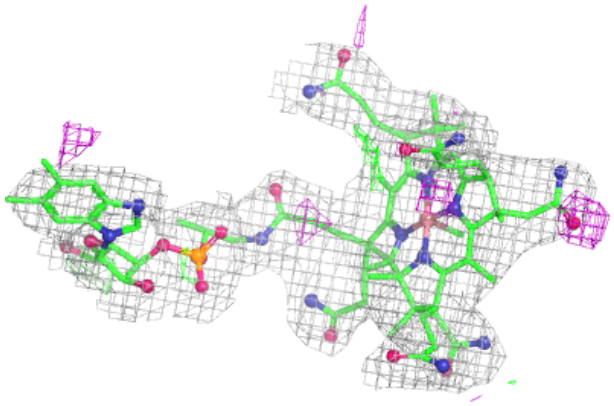
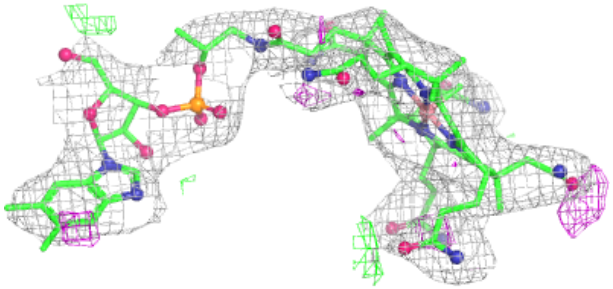
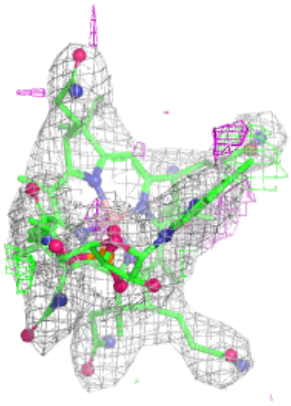
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

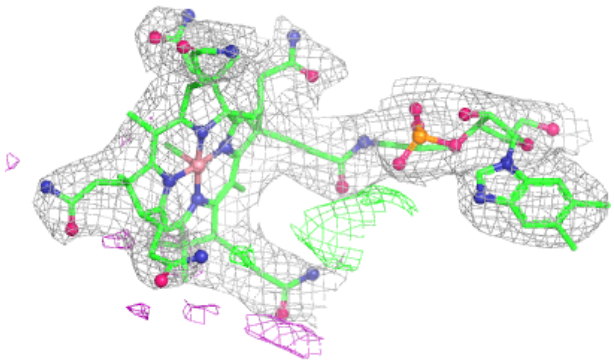


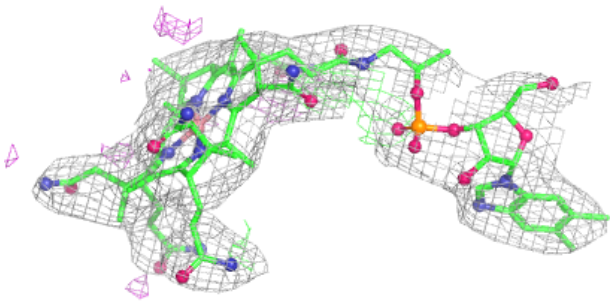


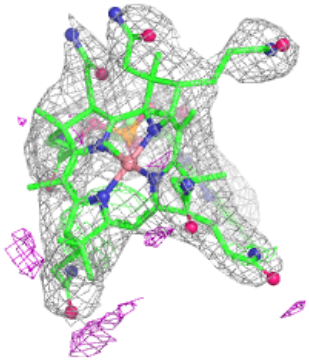


Electron density around COB D 901:

$2mF_o$-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 (i)

There are no such residues in this entry.