



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 11:13 AM EDT

PDB ID : 1BMT / pdb_00001bmt
Title : HOW A PROTEIN BINDS B12: A 3.0 ANGSTROM X-RAY STRUCTURE OF THE B12-BINDING DOMAINS OF METHIONINE SYNTHASE
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Deposited on : 1994-09-02
Resolution : 3.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

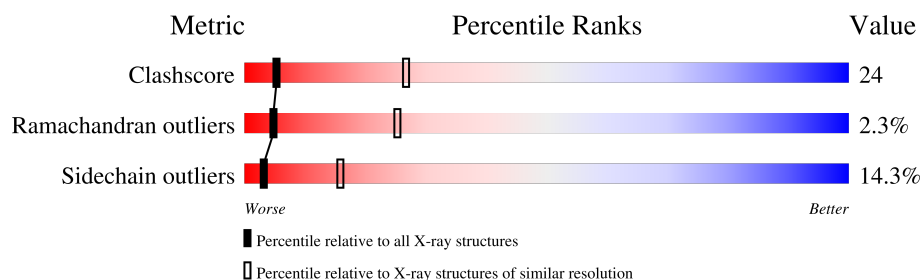
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	246	52% 38% 10%
1	B	246	48% 41% 10%

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	B	122	-	-	X	-

2 Entry composition [i](#)

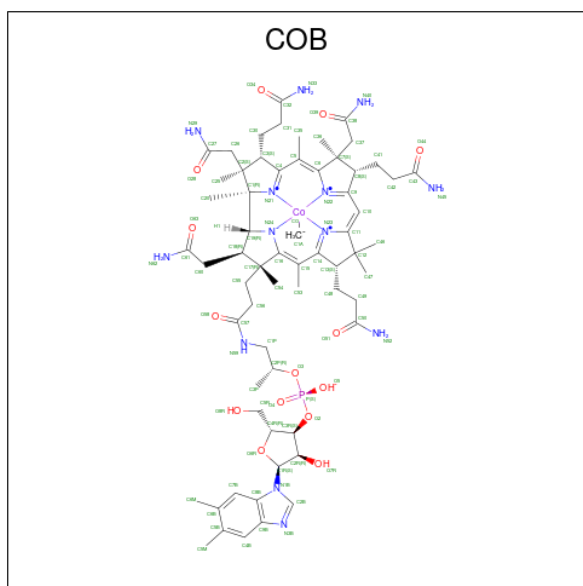
There are 3 unique types of molecules in this entry. The entry contains 4000 atoms, of which 0 are hydrogens and 0 are deuteriums.

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
1	A	246	Total	C	N	O	S	0	0	0
			1907	1198	332	368	9			
1	B	246	Total	C	N	O	S	0	0	0
			1907	1198	332	368	9			

- Molecule 2 is CO-METHYLCOBALAMIN (CCD ID: COB) (formula: $C_{63}H_{92}CoN_{13}O_{14}P$).



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

- Molecule 3 is water.

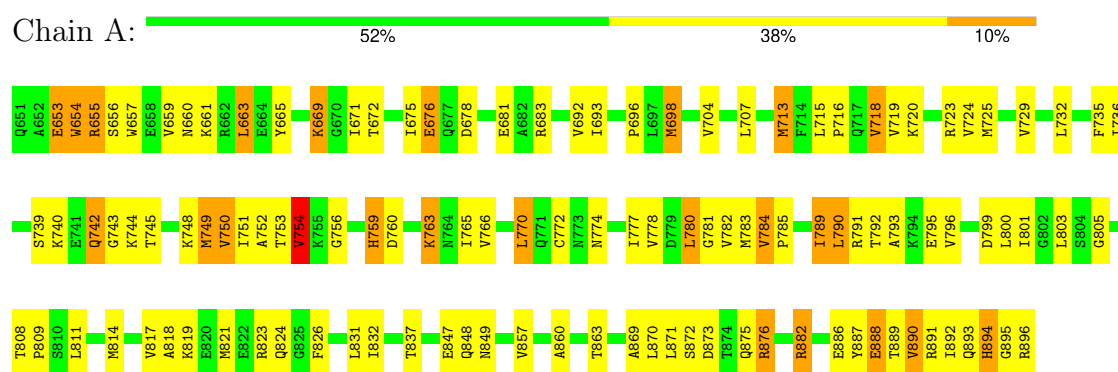
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	O 1	0	0
3	B	1	Total 1	O 1	0	0

3 Residue-property plots

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

Note EDS was not executed.

• Molecule 1: METHIONINE SYNTHASE



• Molecule 1: METHIONINE SYNTHASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	96.70Å 55.30Å 103.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 3.00	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.170 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4000	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: COB

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	0.72	0/1933	1.16	12/2611 (0.5%)
1	B	0.76	1/1933 (0.1%)	1.19	15/2611 (0.6%)
All	All	0.74	1/3866 (0.0%)	1.18	27/5222 (0.5%)

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	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	759	HIS	CD2-NE2	5.21	1.43	1.37

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	784	VAL	CA-C-N	9.54	129.63	119.90
1	B	784	VAL	C-N-CA	9.54	129.63	119.90
1	A	824	GLN	OE1-CD-NE2	-9.22	113.38	122.60
1	A	818	ALA	N-CA-C	-7.94	102.22	111.03
1	A	784	VAL	N-CA-C	7.91	115.44	108.63
1	B	733	GLU	CA-C-N	-6.47	111.76	119.84
1	B	733	GLU	C-N-CA	-6.47	111.76	119.84
1	A	847	GLU	N-CA-C	6.43	118.87	111.02
1	A	824	GLN	CG-CD-NE2	6.19	125.68	116.40
1	B	891	ARG	N-CA-C	-6.18	104.62	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	763	LYS	N-CA-C	-6.06	104.00	111.33
1	B	875	GLN	N-CA-C	5.99	117.89	111.36
1	B	845	LYS	N-CA-C	5.90	121.39	113.72
1	A	732	LEU	N-CA-C	5.85	117.74	111.36
1	A	675	ILE	N-CA-C	5.79	115.98	110.42
1	B	860	ALA	N-CA-C	5.72	117.97	111.11
1	B	738	ALA	N-CA-C	-5.62	105.92	112.89
1	B	746	ASN	N-CA-C	-5.60	106.61	113.50
1	A	882	ARG	N-CA-C	-5.56	105.12	111.07
1	B	718	VAL	N-CA-C	-5.52	104.15	111.05
1	B	759	HIS	CE1-NE2-CD2	-5.48	103.52	109.00
1	B	844	VAL	N-CA-C	5.29	115.50	110.53
1	B	858	GLN	N-CA-C	5.19	117.02	111.36
1	A	653	GLU	N-CA-C	5.17	117.80	110.50
1	B	818	ALA	N-CA-C	-5.13	105.60	111.14
1	A	759	HIS	CE1-NE2-CD2	-5.06	103.94	109.00
1	A	718	VAL	N-CA-C	-5.03	104.80	111.09

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	823	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	1907	0	1939	86	0
1	B	1907	0	1939	93	0
2	A	92	0	78	14	0
2	B	92	0	80	22	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
All	All	4000	0	4036	195	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 24.

All (195) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:122:COB:C1	2:B:122:COB:C20	1.74	1.59
2:A:122:COB:C3R	2:A:122:COB:O2	1.65	1.42
1:A:698:MET:HE1	2:A:122:COB:H5R2	1.42	0.98
1:B:698:MET:HE1	2:B:122:COB:H5R2	1.45	0.97
2:B:122:COB:H362	2:B:122:COB:H351	1.59	0.84
1:B:791:ARG:O	1:B:794:LYS:HG2	1.81	0.80
1:A:752:ALA:HB1	1:A:784:VAL:HG11	1.63	0.79
1:A:748:LYS:HD3	1:A:796:VAL:HG12	1.63	0.79
1:B:752:ALA:HB1	1:B:784:VAL:CG1	2.13	0.78
2:B:122:COB:C20	2:B:122:COB:C2	2.61	0.78
1:B:748:LYS:HD3	1:B:796:VAL:HG12	1.67	0.76
1:A:808:THR:HG23	1:B:893:GLN:HG2	1.66	0.76
1:B:753:THR:HB	1:B:782:VAL:HA	1.68	0.76
1:A:869:ALA:HA	1:A:875:GLN:HE21	1.51	0.75
1:B:752:ALA:HB1	1:B:784:VAL:HG11	1.70	0.74
2:B:122:COB:C20	2:B:122:COB:N21	2.48	0.73
1:A:725:MET:O	1:A:729:VAL:HG23	1.88	0.73
1:B:749:MET:HE3	1:B:751:ILE:HD12	1.69	0.73
1:B:794:LYS:CG	1:B:795:GLU:N	2.52	0.72
2:A:122:COB:H552	2:A:122:COB:H531	1.71	0.72
1:B:808:THR:HB	1:B:809:PRO:HD3	1.71	0.71
2:A:122:COB:H362	2:A:122:COB:H351	1.72	0.71
1:B:660:ASN:OD1	1:B:696:PRO:CG	2.39	0.70
1:A:780:LEU:HD11	1:A:792:THR:HG21	1.74	0.70
2:A:122:COB:C3R	2:A:122:COB:P	2.81	0.68
1:A:720:LYS:HG2	1:A:723:ARG:HH12	1.58	0.68
1:A:707:LEU:HB3	1:A:713:MET:HG3	1.76	0.68
1:B:660:ASN:OD1	1:B:696:PRO:HG3	1.93	0.68
1:B:707:LEU:HD23	1:B:713:MET:HE3	1.75	0.68
1:A:736:ILE:O	1:A:739:SER:HB3	1.95	0.67
1:B:770:LEU:HB3	1:B:777:ILE:HD11	1.76	0.67
1:B:707:LEU:CD2	1:B:713:MET:HE3	2.26	0.66
1:A:752:ALA:HB1	1:A:784:VAL:CG1	2.27	0.65
1:B:794:LYS:HG2	1:B:795:GLU:H	1.61	0.65
1:A:754:VAL:HG23	1:A:803:LEU:HD22	1.79	0.65
1:A:718:VAL:HG13	2:A:122:COB:H462	1.80	0.64
1:A:753:THR:OG1	1:A:763:LYS:HD3	1.96	0.64
1:B:654:TRP:HA	1:B:657:TRP:CE2	2.32	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:693:ILE:HG21	1:A:765:ILE:HD13	1.81	0.63
1:B:754:VAL:HG23	1:B:803:LEU:HD22	1.79	0.63
1:B:869:ALA:HA	1:B:875:GLN:HE21	1.63	0.63
2:A:122:COB:O28	2:A:122:COB:H3	1.99	0.62
1:B:716:PRO:O	1:B:719:VAL:HG12	2.00	0.62
1:B:819:LYS:HG3	1:B:849:ASN:HB3	1.81	0.62
1:A:740:LYS:NZ	1:A:742:GLN:HB2	2.14	0.62
1:A:893:GLN:HG2	1:B:808:THR:HG23	1.83	0.61
1:A:793:ALA:CB	1:A:801:ILE:HD11	2.32	0.60
1:B:889:THR:O	1:B:892:ILE:HB	2.01	0.59
1:A:742:GLN:HB3	1:A:774:ASN:OD1	2.03	0.59
1:A:748:LYS:HD3	1:A:796:VAL:CG1	2.31	0.59
1:B:749:MET:CE	1:B:751:ILE:HD12	2.33	0.58
1:A:832:ILE:HG22	1:A:837:THR:HG21	1.86	0.58
1:A:654:TRP:HA	1:A:657:TRP:CE2	2.39	0.58
1:B:889:THR:HA	1:B:892:ILE:HD12	1.85	0.58
2:B:122:COB:C20	2:B:122:COB:C25	2.82	0.57
1:B:811:LEU:HD12	1:B:841:HIS:CE1	2.38	0.57
1:A:887:TYR:HA	1:A:890:VAL:CG1	2.34	0.57
1:A:750:VAL:CG1	1:A:801:ILE:HG12	2.35	0.57
1:B:794:LYS:NZ	1:B:795:GLU:HG3	2.19	0.57
1:A:750:VAL:HG13	1:A:801:ILE:HG12	1.87	0.56
2:B:122:COB:C20	2:B:122:COB:C19	2.74	0.56
1:A:857:VAL:HG11	1:A:863:THR:HA	1.87	0.56
2:B:122:COB:N40	2:B:122:COB:H8	2.21	0.56
1:A:693:ILE:CG2	1:A:765:ILE:HD13	2.35	0.56
1:B:832:ILE:HG22	1:B:837:THR:HG21	1.87	0.56
1:B:794:LYS:HG3	1:B:795:GLU:N	2.21	0.55
1:B:780:LEU:HD11	1:B:792:THR:HG21	1.88	0.55
1:A:756:GLY:HA2	1:A:783:MET:HE3	1.89	0.55
1:B:718:VAL:HG13	2:B:122:COB:H462	1.87	0.54
1:B:853:PRO:HG3	1:B:880:VAL:HG21	1.90	0.54
1:A:663:LEU:CD1	1:A:678:ASP:HB3	2.38	0.54
1:A:748:LYS:CD	1:A:796:VAL:HG12	2.36	0.54
1:B:698:MET:SD	2:B:122:COB:H3P2	2.48	0.54
1:A:760:ASP:HB2	1:A:763:LYS:HE2	1.88	0.54
1:A:888:GLU:O	1:A:892:ILE:HG13	2.08	0.54
1:B:744:LYS:HD3	1:B:745:THR:HG22	1.88	0.54
1:A:753:THR:HB	1:A:782:VAL:HA	1.90	0.54
1:A:765:ILE:HD12	2:A:122:COB:H3P1	1.89	0.54
1:A:720:LYS:HG2	1:A:723:ARG:NH1	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:770:LEU:HB3	1:A:777:ILE:HD11	1.90	0.53
1:A:872:SER:O	1:A:876:ARG:HB2	2.08	0.53
1:A:808:THR:HB	1:A:809:PRO:HD3	1.91	0.53
1:B:786:ALA:O	1:B:790:LEU:HD22	2.09	0.53
1:B:860:ALA:H	2:B:122:COB:H5R1	1.75	0.52
1:A:791:ARG:O	1:A:795:GLU:HG3	2.09	0.52
1:A:869:ALA:HA	1:A:875:GLN:NE2	2.23	0.52
1:A:793:ALA:HB2	1:A:801:ILE:HD11	1.91	0.52
1:B:663:LEU:CD1	1:B:678:ASP:HB3	2.40	0.52
1:B:813:GLU:O	1:B:817:VAL:HG13	2.09	0.52
1:A:790:LEU:HG	1:A:826:PHE:CE1	2.45	0.51
1:A:716:PRO:C	1:A:719:VAL:HG12	2.35	0.51
1:B:857:VAL:HB	2:B:122:COB:HM61	1.92	0.51
2:B:122:COB:H552	2:B:122:COB:H531	1.92	0.51
2:B:122:COB:C20	2:B:122:COB:H253	2.40	0.51
2:A:122:COB:O2	2:A:122:COB:C4R	2.53	0.51
2:B:122:COB:H262	2:B:122:COB:H601	1.92	0.51
1:B:654:TRP:HZ3	1:B:662:ARG:HB3	1.76	0.51
1:B:758:VAL:HG12	2:B:122:COB:H421	1.93	0.51
1:A:886:GLU:O	1:A:890:VAL:HG12	2.12	0.50
1:A:740:LYS:HZ1	1:A:742:GLN:HB2	1.75	0.50
1:B:794:LYS:HZ2	1:B:795:GLU:HG3	1.77	0.50
1:A:805:GLY:HA3	1:A:814:MET:HE3	1.94	0.49
1:B:704:VAL:CG1	1:B:718:VAL:HG23	2.42	0.49
1:B:771:GLN:C	1:B:773:ASN:H	2.20	0.49
1:A:716:PRO:O	1:A:719:VAL:HG12	2.12	0.49
1:B:741:GLU:C	1:B:742:GLN:HG3	2.38	0.49
1:A:744:LYS:HG3	1:A:745:THR:H	1.77	0.49
1:A:873:ASP:HA	1:A:876:ARG:NH1	2.27	0.49
1:A:665:TYR:HE1	1:A:671:ILE:HD11	1.77	0.49
1:B:672:THR:HG22	1:B:724:VAL:HG22	1.94	0.48
1:A:657:TRP:HB3	1:A:661:LYS:HD3	1.94	0.48
1:A:752:ALA:CB	1:A:784:VAL:HG11	2.39	0.48
1:B:873:ASP:HA	1:B:876:ARG:NH1	2.28	0.48
1:A:749:MET:HE3	1:A:751:ILE:HD12	1.96	0.48
1:A:654:TRP:N	1:A:654:TRP:CD1	2.82	0.48
1:A:672:THR:HG22	1:A:724:VAL:HG22	1.96	0.47
1:B:660:ASN:OD1	1:B:696:PRO:HG2	2.14	0.47
1:B:842:THR:HG21	1:B:856:TYR:HB2	1.96	0.47
1:A:882:ARG:HE	1:B:710:GLU:CD	2.22	0.47
1:A:763:LYS:O	1:A:766:VAL:HG12	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:652:ALA:O	1:B:655:ARG:HB2	2.15	0.47
1:B:654:TRP:C	1:B:656:SER:H	2.22	0.47
1:B:831:LEU:HD12	1:B:855:VAL:HB	1.95	0.47
1:B:654:TRP:CZ3	1:B:662:ARG:HB3	2.49	0.47
1:A:790:LEU:HD11	1:A:821:MET:HG2	1.96	0.47
1:A:860:ALA:H	2:A:122:COB:H5R1	1.80	0.47
1:A:893:GLN:C	1:A:895:GLY:H	2.23	0.46
1:B:784:VAL:HA	1:B:785:PRO:HD3	1.73	0.46
1:A:759:HIS:CE1	2:A:122:COB:H412	2.51	0.46
1:A:676:GLU:H	1:A:676:GLU:HG2	1.55	0.46
1:A:749:MET:CE	1:A:751:ILE:HD12	2.46	0.45
1:A:805:GLY:HA3	1:A:814:MET:CE	2.46	0.45
1:B:896:ARG:O	1:B:896:ARG:HD3	2.16	0.45
1:A:665:TYR:CE1	1:A:669:LYS:HG3	2.50	0.45
1:A:759:HIS:NE2	2:A:122:COB:H202	2.31	0.45
2:A:122:COB:H601	2:A:122:COB:H262	1.97	0.45
1:B:831:LEU:CD1	1:B:855:VAL:HB	2.47	0.45
1:B:859:ASN:HB2	2:B:122:COB:H5R1	1.99	0.45
1:B:692:VAL:CG2	1:B:732:LEU:HD11	2.46	0.45
1:B:866:VAL:HG12	1:B:870:LEU:HD22	1.97	0.45
1:A:659:VAL:HG21	1:A:681:GLU:HB3	1.98	0.45
1:A:660:ASN:OD1	1:A:696:PRO:CG	2.65	0.45
1:B:731:TYR:O	1:B:734:PRO:HD2	2.16	0.44
1:B:805:GLY:HA3	1:B:814:MET:HE3	2.00	0.44
1:B:759:HIS:HE1	2:B:122:COB:H412	1.82	0.44
1:B:753:THR:HG22	1:B:757:ASP:HB3	2.00	0.44
1:A:660:ASN:OD1	1:A:696:PRO:HG3	2.18	0.44
1:A:872:SER:HB3	1:A:875:GLN:HB2	1.98	0.44
1:B:752:ALA:HB1	1:B:784:VAL:HG13	1.98	0.44
1:B:754:VAL:HG13	1:B:755:LYS:N	2.32	0.44
1:B:720:LYS:O	1:B:723:ARG:HG2	2.18	0.43
2:B:122:COB:O28	2:B:122:COB:H3	2.17	0.43
1:A:715:LEU:O	1:A:718:VAL:HG12	2.19	0.43
2:B:122:COB:H252	2:B:122:COB:N62	2.34	0.43
1:A:754:VAL:CG2	1:A:803:LEU:HD22	2.48	0.43
1:A:707:LEU:CD2	1:A:713:MET:HE3	2.49	0.43
1:A:784:VAL:HG13	1:A:789:ILE:HD11	2.01	0.43
1:B:755:LYS:HB3	1:B:813:GLU:HG3	2.01	0.43
1:B:753:THR:HG21	1:B:759:HIS:O	2.19	0.42
1:B:830:LEU:O	1:B:854:THR:HA	2.19	0.42
1:B:806:LEU:HD23	2:B:122:COB:H201	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:801:ILE:HD12	1:A:821:MET:HE3	2.01	0.42
1:B:763:LYS:HZ2	1:B:781:GLY:HA2	1.84	0.42
1:B:888:GLU:O	1:B:892:ILE:HG13	2.19	0.42
1:A:718:VAL:CG1	2:A:122:COB:H462	2.49	0.42
1:B:763:LYS:HE2	1:B:763:LYS:HB3	1.88	0.42
1:B:796:VAL:O	1:B:797:ASN:HB2	2.19	0.42
1:B:816:ASN:C	1:B:816:ASN:ND2	2.77	0.42
1:B:692:VAL:HG21	1:B:732:LEU:HD11	2.01	0.42
1:B:712:LYS:HE2	1:B:712:LYS:HB3	1.89	0.42
1:B:750:VAL:HA	1:B:778:VAL:O	2.20	0.41
1:A:894:HIS:C	1:B:845:LYS:HD2	2.45	0.41
1:B:707:LEU:HD22	1:B:713:MET:HE3	2.01	0.41
1:B:660:ASN:CG	1:B:696:PRO:HG3	2.44	0.41
1:B:831:LEU:HD12	1:B:831:LEU:HA	1.92	0.41
1:A:704:VAL:CG1	1:A:718:VAL:HG23	2.50	0.41
1:B:715:LEU:O	1:B:718:VAL:HG12	2.20	0.41
1:A:819:LYS:HG3	1:A:849:ASN:HB3	2.02	0.41
1:B:872:SER:HB3	1:B:875:GLN:HB2	2.03	0.41
1:A:683:ARG:NH2	1:A:735:PHE:CE1	2.88	0.41
1:B:771:GLN:C	1:B:773:ASN:N	2.77	0.41
2:B:122:COB:H541	2:B:122:COB:H602	1.98	0.41
1:B:790:LEU:HD13	1:B:790:LEU:N	2.35	0.41
1:A:889:THR:HA	1:A:892:ILE:HD12	2.02	0.40
1:A:891:ARG:HB3	1:A:896:ARG:HD2	2.03	0.40
1:B:807:ILE:HD13	1:B:807:ILE:HG21	1.82	0.40
1:B:823:ARG:C	1:B:825:GLY:H	2.29	0.40
1:A:707:LEU:HD22	1:A:713:MET:HE3	2.03	0.40
1:A:784:VAL:HA	1:A:785:PRO:HD3	1.82	0.40
1:B:657:TRP:HB3	1:B:661:LYS:HD3	2.02	0.40
1:B:705:GLY:O	1:B:708:PHE:HB3	2.21	0.40
1:A:763:LYS:NZ	1:A:781:GLY:HA2	2.36	0.40
1:A:796:VAL:HG12	1:A:796:VAL:O	2.21	0.40
1:B:794:LYS:HZ3	1:B:795:GLU:CG	2.34	0.40
1:B:803:LEU:HD11	1:B:817:VAL:HG22	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	244/246 (99%)	225 (92%)	11 (4%)	8 (3%)	3	17
1	B	244/246 (99%)	218 (89%)	23 (9%)	3 (1%)	10	40
All	All	488/492 (99%)	443 (91%)	34 (7%)	11 (2%)	5	25

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	740	LYS
1	A	848	GLN
1	A	876	ARG
1	A	655	ARG
1	A	742	GLN
1	A	743	GLY
1	A	754	VAL
1	B	655	ARG
1	B	876	ARG
1	A	654	TRP
1	A	894	HIS

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	207/207 (100%)	180 (87%)	27 (13%)	4	19
1	B	207/207 (100%)	175 (84%)	32 (16%)	2	13

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	414/414 (100%)	355 (86%)	59 (14%)	3 16

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	653	GLU
1	A	655	ARG
1	A	656	SER
1	A	663	LEU
1	A	669	LYS
1	A	676	GLU
1	A	692	VAL
1	A	698	MET
1	A	713	MET
1	A	749	MET
1	A	750	VAL
1	A	754	VAL
1	A	770	LEU
1	A	772	CYS
1	A	778	VAL
1	A	780	LEU
1	A	789	ILE
1	A	790	LEU
1	A	799	ASP
1	A	800	LEU
1	A	811	LEU
1	A	817	VAL
1	A	831	LEU
1	A	870	LEU
1	A	871	LEU
1	A	888	GLU
1	A	890	VAL
1	B	654	TRP
1	B	655	ARG
1	B	663	LEU
1	B	669	LYS
1	B	675	ILE
1	B	676	GLU
1	B	689	PRO
1	B	692	VAL
1	B	697	LEU
1	B	701	MET

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Mol	Chain	Res	Type
1	B	715	LEU
1	B	719	VAL
1	B	737	GLU
1	B	742	GLN
1	B	750	VAL
1	B	754	VAL
1	B	770	LEU
1	B	772	CYS
1	B	778	VAL
1	B	780	LEU
1	B	784	VAL
1	B	789	ILE
1	B	790	LEU
1	B	800	LEU
1	B	806	LEU
1	B	817	VAL
1	B	831	LEU
1	B	870	LEU
1	B	871	LEU
1	B	888	GLU
1	B	890	VAL
1	B	896	ARG

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	685	GLN
1	A	771	GLN
1	A	841	HIS
1	A	875	GLN
1	B	685	GLN
1	B	771	GLN
1	B	774	ASN
1	B	816	ASN
1	B	841	HIS
1	B	848	GLN
1	B	875	GLN

5.3.3 RNA ⓘ

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](#)

2 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	$\# Z > 2$
2	COB	B	122	1	93,102,102	3.10	41 (44%)	149,170,170	2.46	50 (33%)
2	COB	A	122	1	93,102,102	2.95	42 (45%)	149,170,170	2.53	53 (35%)

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
2	COB	B	122	1	-	11/56/231/231	0/3/11/11
2	COB	A	122	1	-	16/56/231/231	0/3/11/11

All (83) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	122	COB	C1-N21	-11.52	1.32	1.50
2	B	122	COB	C20-C1	10.86	1.74	1.53
2	A	122	COB	O6R-C1R	9.21	1.63	1.42
2	A	122	COB	C53-C15	8.90	1.69	1.50
2	B	122	COB	C4B-C9B	8.58	1.53	1.40
2	A	122	COB	C1-N21	-7.66	1.38	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	122	COB	C7B-C6B	7.05	1.49	1.39
2	B	122	COB	C5M-C5B	-6.23	1.39	1.51
2	A	122	COB	O2-C3R	6.22	1.65	1.44
2	B	122	COB	C37-C38	-5.92	1.32	1.51
2	B	122	COB	C4B-C5B	5.74	1.47	1.39
2	A	122	COB	C5M-C5B	-5.62	1.40	1.51
2	A	122	COB	C4B-C5B	5.61	1.47	1.39
2	B	122	COB	O2-C3R	5.29	1.62	1.44
2	B	122	COB	C6M-C6B	5.12	1.60	1.51
2	A	122	COB	C4B-C9B	4.97	1.47	1.40
2	B	122	COB	O6R-C1R	4.83	1.53	1.42
2	A	122	COB	C20-C1	4.80	1.62	1.53
2	A	122	COB	C7B-C8B	4.63	1.47	1.39
2	A	122	COB	C49-C50	-4.30	1.34	1.51
2	B	122	COB	C1P-C2P	-4.25	1.40	1.51
2	B	122	COB	C7B-C8B	4.21	1.46	1.39
2	A	122	COB	C6M-C6B	4.12	1.58	1.51
2	B	122	COB	C1-C19	-4.09	1.43	1.54
2	B	122	COB	C60-C18	-3.94	1.45	1.54
2	A	122	COB	C41-C8	-3.92	1.45	1.54
2	A	122	COB	C50-N52	-3.87	1.20	1.32
2	A	122	COB	C35-C5	3.76	1.58	1.50
2	A	122	COB	O8R-C5R	-3.70	1.26	1.42
2	B	122	COB	C56-C57	-3.69	1.44	1.51
2	B	122	COB	C2-C3	-3.68	1.46	1.57
2	A	122	COB	O6R-C4R	-3.67	1.36	1.45
2	B	122	COB	C42-C43	3.56	1.65	1.51
2	A	122	COB	C8B-C9B	3.55	1.46	1.40
2	B	122	COB	C35-C5	3.42	1.57	1.50
2	B	122	COB	C19-N24	-3.37	1.42	1.49
2	A	122	COB	C26-C27	-3.37	1.40	1.51
2	B	122	COB	C13-C14	3.36	1.60	1.51
2	B	122	COB	C41-C8	-3.29	1.46	1.54
2	A	122	COB	C25-C2	3.23	1.60	1.54
2	B	122	COB	C50-N52	-3.18	1.22	1.32
2	A	122	COB	C32-N33	-3.18	1.22	1.32
2	A	122	COB	C60-C18	-3.16	1.47	1.54
2	B	122	COB	O8R-C5R	-3.13	1.29	1.42
2	A	122	COB	C5R-C4R	-3.13	1.41	1.51
2	B	122	COB	C5R-C4R	-3.07	1.41	1.51
2	A	122	COB	C56-C57	3.03	1.57	1.51
2	A	122	COB	C48-C13	-2.99	1.47	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	122	COB	C43-N45	-2.99	1.23	1.32
2	A	122	COB	C47-C12	2.99	1.60	1.54
2	B	122	COB	C43-N45	-2.98	1.23	1.32
2	B	122	COB	P-O5	-2.93	1.41	1.55
2	B	122	COB	C49-C50	-2.93	1.39	1.51
2	A	122	COB	C37-C38	2.92	1.61	1.51
2	A	122	COB	C1-C19	-2.91	1.46	1.54
2	B	122	COB	C7B-C6B	2.88	1.43	1.39
2	B	122	COB	O6R-C4R	-2.87	1.38	1.45
2	B	122	COB	C17-C18	-2.85	1.49	1.54
2	B	122	COB	C41-C42	-2.84	1.44	1.52
2	A	122	COB	C6B-C5B	2.81	1.47	1.40
2	B	122	COB	C53-C15	2.80	1.56	1.50
2	B	122	COB	C38-N40	-2.79	1.23	1.32
2	A	122	COB	C27-N29	-2.78	1.23	1.32
2	A	122	COB	C61-N62	-2.67	1.24	1.32
2	B	122	COB	C12-C13	-2.67	1.47	1.55
2	A	122	COB	C2-C3	-2.67	1.49	1.57
2	A	122	COB	C8-C9	-2.66	1.45	1.51
2	A	122	COB	C18-C19	-2.64	1.49	1.54
2	A	122	COB	C41-C42	-2.62	1.45	1.52
2	B	122	COB	C30-C31	-2.61	1.45	1.52
2	A	122	COB	C12-C13	-2.56	1.47	1.55
2	B	122	COB	C30-C3	-2.51	1.48	1.54
2	B	122	COB	C61-N62	-2.49	1.24	1.32
2	A	122	COB	C30-C3	-2.44	1.48	1.54
2	A	122	COB	C19-N24	-2.42	1.44	1.49
2	B	122	COB	C2B-N1B	-2.38	1.33	1.37
2	B	122	COB	C8-C9	-2.37	1.46	1.51
2	A	122	COB	C4-C5	-2.31	1.33	1.43
2	A	122	COB	C17-C18	-2.20	1.50	1.54
2	B	122	COB	C26-C2	-2.19	1.50	1.55
2	B	122	COB	C7-C6	2.08	1.59	1.54
2	B	122	COB	C25-C2	2.08	1.58	1.54
2	A	122	COB	C55-C56	-2.07	1.48	1.53

All (103) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	122	COB	O58-C57-C56	-8.63	106.39	122.02
2	A	122	COB	C1-N21-C4	8.57	123.41	109.48
2	A	122	COB	C19-N24-C16	7.78	124.41	111.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	122	COB	O6R-C1R-N1B	7.39	122.28	108.09
2	B	122	COB	C56-C55-C17	6.65	128.40	115.58
2	B	122	COB	C56-C57-N59	6.51	128.21	116.34
2	B	122	COB	C20-C1-C2	-6.51	102.54	113.28
2	A	122	COB	C1-C19-C18	6.23	131.11	121.75
2	A	122	COB	C54-C17-C55	-6.07	99.17	109.27
2	A	122	COB	C17-C16-N24	-5.61	101.49	110.91
2	A	122	COB	C1-C19-N24	5.42	114.93	106.29
2	B	122	COB	O3-C2P-C1P	-5.41	96.23	106.94
2	A	122	COB	O58-C57-C56	-5.34	112.35	122.02
2	A	122	COB	C2-C1-C19	5.12	127.59	118.77
2	B	122	COB	C19-N24-C16	5.10	120.12	111.96
2	B	122	COB	O39-C38-C37	-5.04	106.36	121.98
2	A	122	COB	O39-C38-C37	-4.98	106.56	121.98
2	A	122	COB	O3-C2P-C1P	-4.91	97.21	106.94
2	A	122	COB	C18-C19-N24	-4.91	94.12	101.92
2	A	122	COB	C7-C6-N22	-4.87	99.07	107.94
2	A	122	COB	C15-C16-N24	4.72	127.12	122.38
2	B	122	COB	C17-C18-C19	4.66	108.37	102.70
2	A	122	COB	C35-C5-C4	-4.52	107.62	116.79
2	A	122	COB	C19-C1-N21	4.51	107.37	101.69
2	A	122	COB	C12-C11-N23	-4.42	106.77	111.49
2	B	122	COB	C8-C7-C6	4.42	108.42	100.92
2	B	122	COB	C20-C1-C19	4.37	118.89	110.23
2	A	122	COB	C17-C18-C19	4.31	107.94	102.70
2	A	122	COB	C3-C4-N21	-4.29	106.65	111.98
2	B	122	COB	C17-C16-N24	-4.24	103.79	110.91
2	A	122	COB	C54-C17-C18	4.23	119.06	112.99
2	B	122	COB	O28-C27-C26	-4.21	108.94	121.98
2	A	122	COB	C20-C1-C2	-4.14	106.45	113.28
2	B	122	COB	C7-C8-C9	-4.10	95.69	100.89
2	A	122	COB	O39-C38-N40	4.10	133.48	122.53
2	B	122	COB	C37-C38-N40	4.05	129.05	116.49
2	B	122	COB	O6R-C4R-C3R	3.89	113.13	104.92
2	A	122	COB	O7R-C2R-C3R	-3.89	100.31	111.19
2	B	122	COB	O6R-C4R-C5R	3.87	117.40	109.22
2	B	122	COB	C31-C30-C3	-3.85	103.73	114.65
2	A	122	COB	C2R-C1R-N1B	3.84	122.85	113.30
2	B	122	COB	C1-N21-C4	3.83	115.70	109.48
2	B	122	COB	C7-C6-N22	-3.77	101.08	107.94
2	B	122	COB	C18-C19-N24	-3.70	96.03	101.92
2	A	122	COB	C5-C6-N22	3.68	129.45	123.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	122	COB	C35-C5-C6	3.63	128.26	122.41
2	B	122	COB	O28-C27-N29	3.62	132.19	122.53
2	A	122	COB	C25-C2-C3	3.47	121.72	112.91
2	B	122	COB	O44-C43-N45	-3.42	113.40	122.53
2	A	122	COB	C25-C2-C1	-3.37	108.72	113.75
2	A	122	COB	O6R-C1R-C2R	-3.31	99.53	106.62
2	B	122	COB	O8R-C5R-C4R	-3.30	100.11	111.33
2	B	122	COB	C9-C10-C11	-3.29	120.99	125.84
2	B	122	COB	C60-C61-N62	-3.26	108.38	116.19
2	B	122	COB	C49-C48-C13	-3.25	105.43	114.65
2	A	122	COB	C10-C9-N22	3.25	129.45	125.74
2	A	122	COB	C8-C7-C6	3.15	106.26	100.92
2	B	122	COB	C35-C5-C6	-3.15	117.34	122.41
2	B	122	COB	O51-C50-C49	-3.12	111.61	121.04
2	A	122	COB	C17-C16-C15	3.12	131.39	126.70
2	B	122	COB	C54-C17-C18	3.11	117.45	112.99
2	A	122	COB	O2-P-O3	-3.11	94.19	102.87
2	B	122	COB	C17-C16-C15	3.09	131.34	126.70
2	B	122	COB	C8B-N1B-C2B	3.02	109.00	106.26
2	B	122	COB	C53-C15-C16	2.95	126.75	120.42
2	A	122	COB	C20-C1-C19	-2.94	104.40	110.23
2	A	122	COB	O63-C61-C60	-2.93	114.75	120.87
2	A	122	COB	C1-C2-C3	2.92	105.27	101.60
2	B	122	COB	C1-C19-N24	2.91	110.93	106.29
2	B	122	COB	C7-C6-C5	2.87	132.56	128.07
2	A	122	COB	C25-C2-C26	-2.77	104.24	109.74
2	A	122	COB	C56-C57-N59	2.77	121.38	116.34
2	B	122	COB	N1B-C2B-N3B	-2.75	110.04	113.94
2	A	122	COB	C36-C7-C8	2.71	117.05	112.05
2	A	122	COB	C3P-C2P-C1P	-2.70	106.21	111.42
2	A	122	COB	C5-C4-N21	2.63	128.42	124.19
2	B	122	COB	C2-C1-C19	2.58	123.23	118.77
2	A	122	COB	C2-C1-N21	-2.55	98.25	101.78
2	A	122	COB	O8R-C5R-C4R	-2.51	102.78	111.33
2	B	122	COB	C6M-C6B-C5B	2.51	125.88	120.76
2	B	122	COB	C12-C11-C10	-2.46	121.33	123.54
2	A	122	COB	C2P-C1P-N59	2.46	116.53	112.92
2	A	122	COB	C7-C8-C9	-2.41	97.83	100.89
2	A	122	COB	C3R-C2R-C1R	2.40	105.17	99.89
2	B	122	COB	C15-C16-N24	2.39	124.78	122.38
2	A	122	COB	C41-C8-C7	2.38	120.72	114.19
2	B	122	COB	C42-C43-N45	2.31	123.90	116.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	122	COB	O51-C50-N52	2.31	128.69	122.53
2	B	122	COB	C9B-C4B-C5B	-2.28	116.61	120.83
2	B	122	COB	C53-C15-C14	-2.26	112.20	116.79
2	B	122	COB	C2-C26-C27	2.24	121.43	115.19
2	A	122	COB	C8B-N1B-C1R	-2.21	121.67	125.41
2	A	122	COB	C8-C9-N22	-2.19	106.44	110.77
2	A	122	COB	C4R-O6R-C1R	2.17	114.26	109.47
2	A	122	COB	O28-C27-N29	2.17	128.32	122.53
2	A	122	COB	C7-C6-C5	2.16	131.46	128.07
2	B	122	COB	C7-C37-C38	2.16	120.66	114.28
2	A	122	COB	C42-C41-C8	-2.15	108.56	114.65
2	B	122	COB	C12-C11-N23	2.12	113.76	111.49
2	B	122	COB	O7R-C2R-C3R	-2.10	105.32	111.19
2	B	122	COB	C4R-O6R-C1R	-2.05	104.94	109.47
2	B	122	COB	C41-C42-C43	-2.05	105.59	112.55
2	A	122	COB	C9B-C8B-N1B	-2.01	104.41	105.30

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	122	COB	C25-C2-C26-C27
2	A	122	COB	C42-C41-C8-C9
2	B	122	COB	C7-C37-C38-N40
2	A	122	COB	C42-C41-C8-C7
2	B	122	COB	C7-C37-C38-O39
2	A	122	COB	C2R-C1R-N1B-C8B
2	A	122	COB	O6R-C1R-N1B-C8B
2	B	122	COB	C2R-C1R-N1B-C8B
2	B	122	COB	C3R-C4R-C5R-O8R
2	A	122	COB	C3P-C2P-O3-P
2	A	122	COB	C7-C37-C38-N40
2	B	122	COB	O58-C57-N59-C1P
2	A	122	COB	C7-C37-C38-O39
2	A	122	COB	C3-C2-C26-C27
2	B	122	COB	C42-C41-C8-C9
2	A	122	COB	C2R-C1R-N1B-C2B
2	A	122	COB	C17-C18-C60-C61
2	A	122	COB	C19-C18-C60-C61
2	B	122	COB	O6R-C1R-N1B-C8B
2	B	122	COB	C19-C18-C60-C61
2	A	122	COB	N59-C1P-C2P-O3

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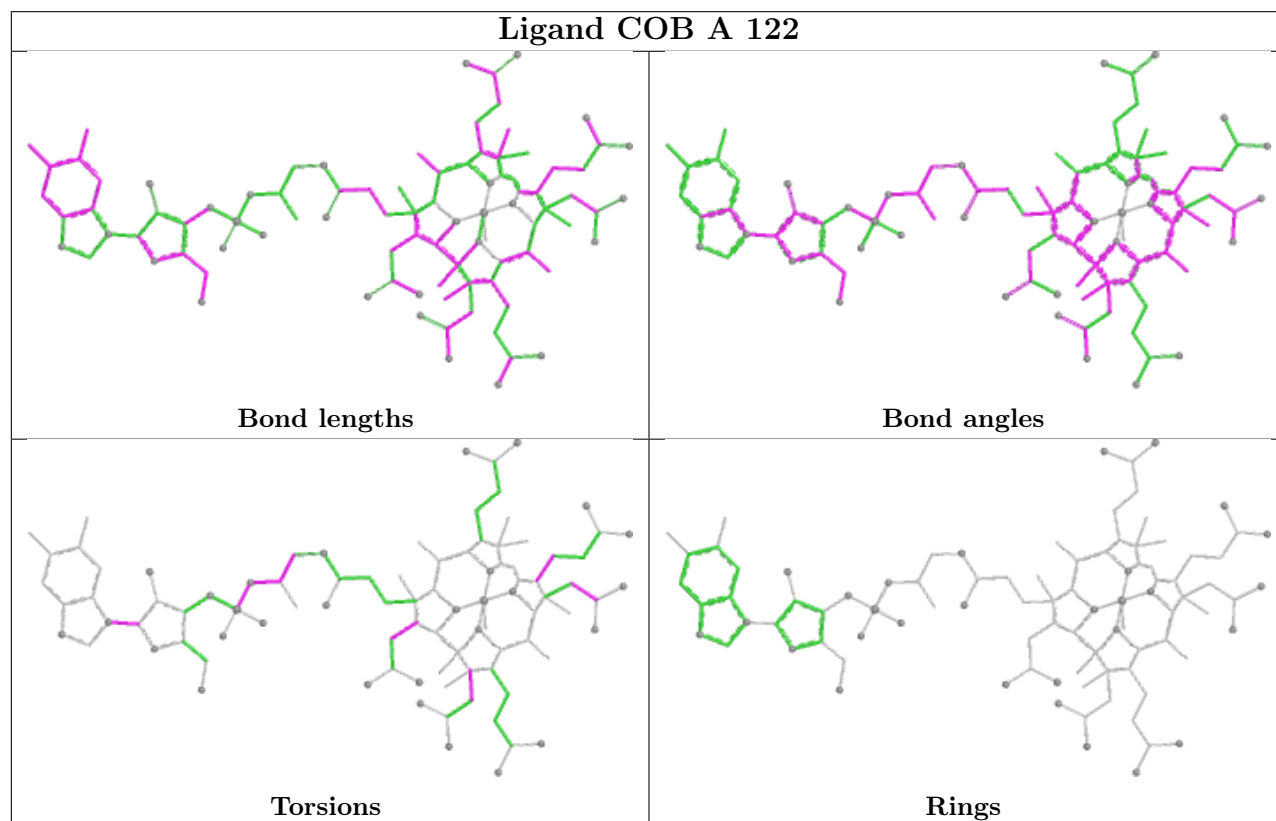
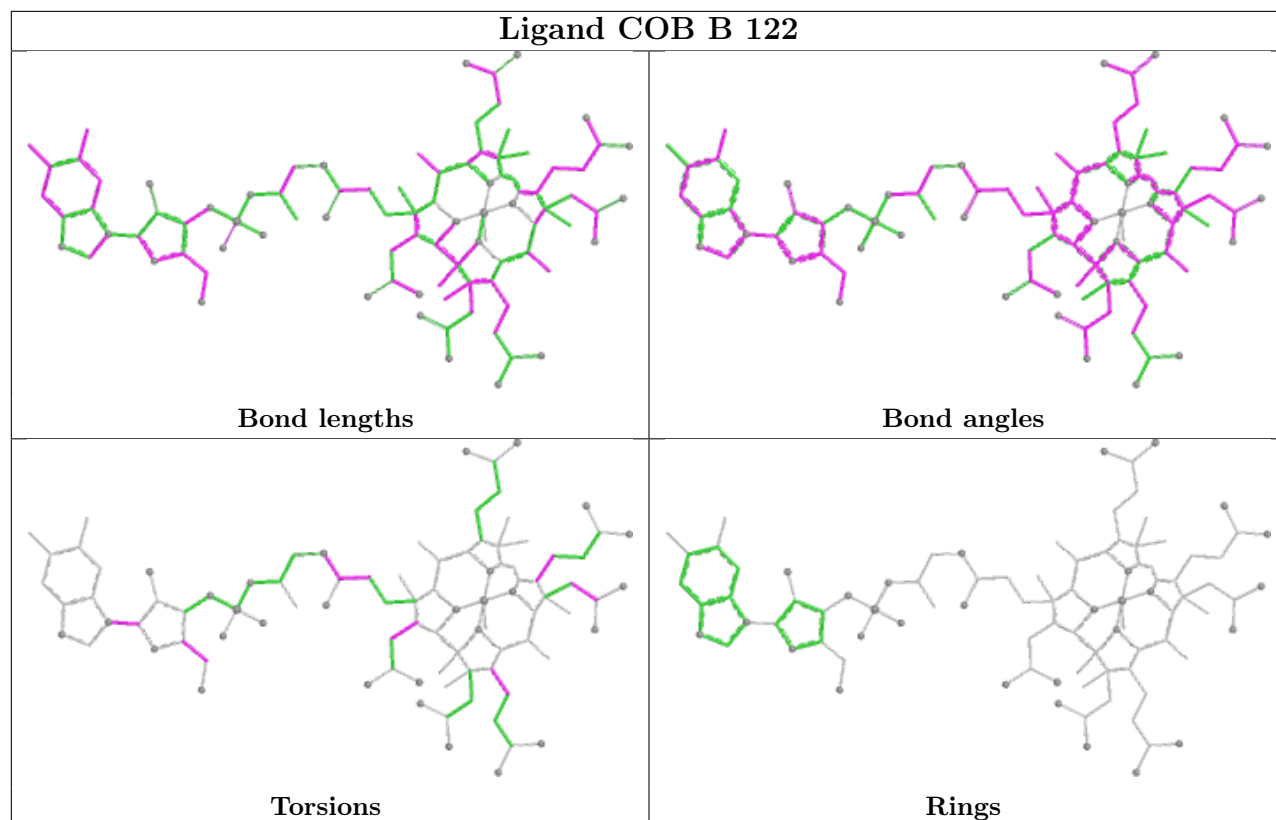
Mol	Chain	Res	Type	Atoms
2	B	122	COB	C4-C3-C30-C31
2	A	122	COB	C1-C2-C26-C27
2	B	122	COB	C17-C18-C60-C61
2	A	122	COB	O6R-C1R-N1B-C2B
2	A	122	COB	C2P-O3-P-O5
2	B	122	COB	C55-C56-C57-O58

There are no ring outliers.

2 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	122	COB	22	0
2	A	122	COB	14	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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.