



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:34 PM UTC

PDB ID : 1T4C / pdb_00001t4c
Title : Formyl-CoA Transferase in complex with Oxalyl-CoA
Authors : Ricagno, S.; Jonsson, S.; Richards, N.G.; Lindqvist, Y.
Deposited on : 2004-04-29
Resolution : 2.61 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

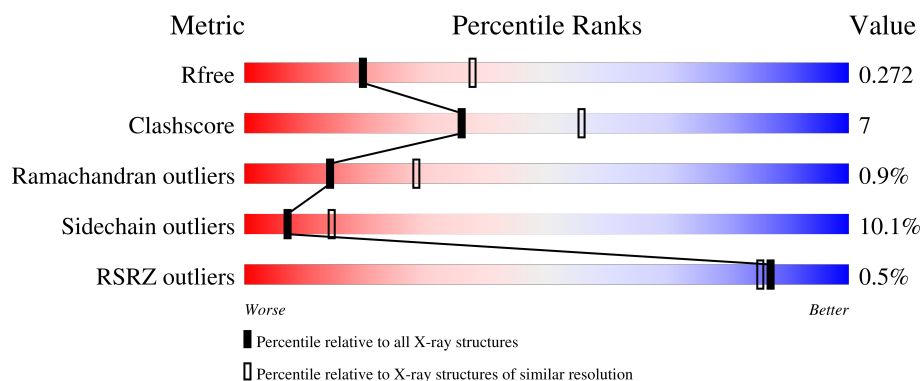
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 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	4951 (2.64-2.60)
Clashscore	190562	5303 (2.64-2.60)
Ramachandran outliers	187476	5217 (2.64-2.60)
Sidechain outliers	187428	5217 (2.64-2.60)
RSRZ outliers	180081	4950 (2.64-2.60)

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	427	79% 18% .
1	B	427	78% 19% .

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	OXD	B	1002	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6857 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.

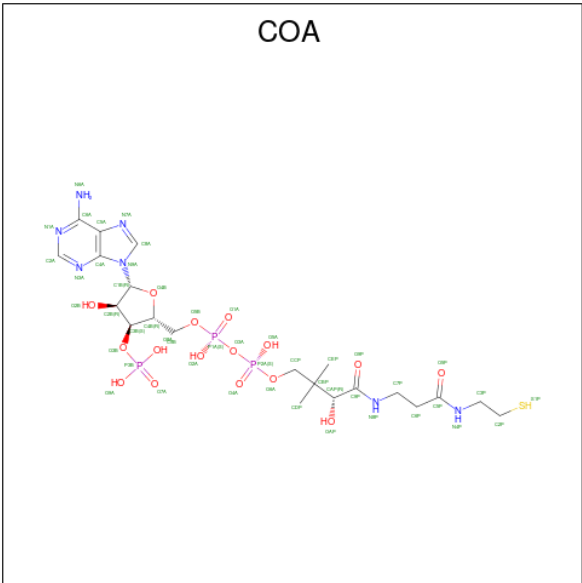
- Molecule 1 is a protein called Formyl-CoA:oxalate CoA-transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	427	3313	2096	568	626	23	0	0	0
1	B	427	3313	2096	568	626	23	4	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	186	ILE	MET	conflict	UNP O06644
B	186	ILE	MET	conflict	UNP O06644

- Molecule 2 is COENZYME A (CCD ID: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S).



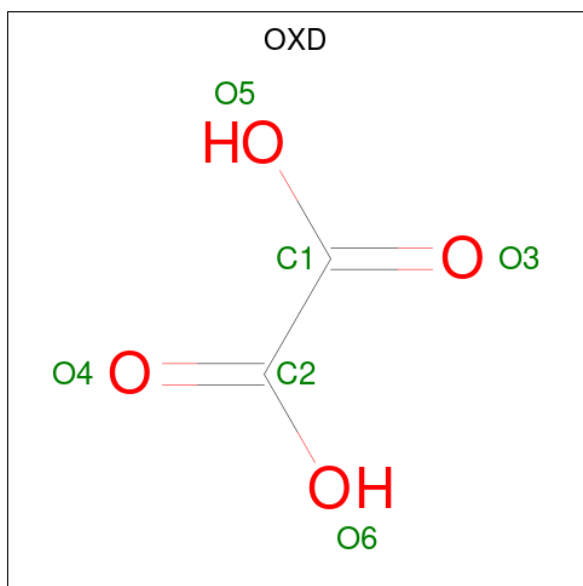
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
2	A	1	48	21	7	16	3	1	0	0

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 3 is OXALIC ACID (CCD ID: OXD) (formula: $C_2H_2O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			5	2	3		

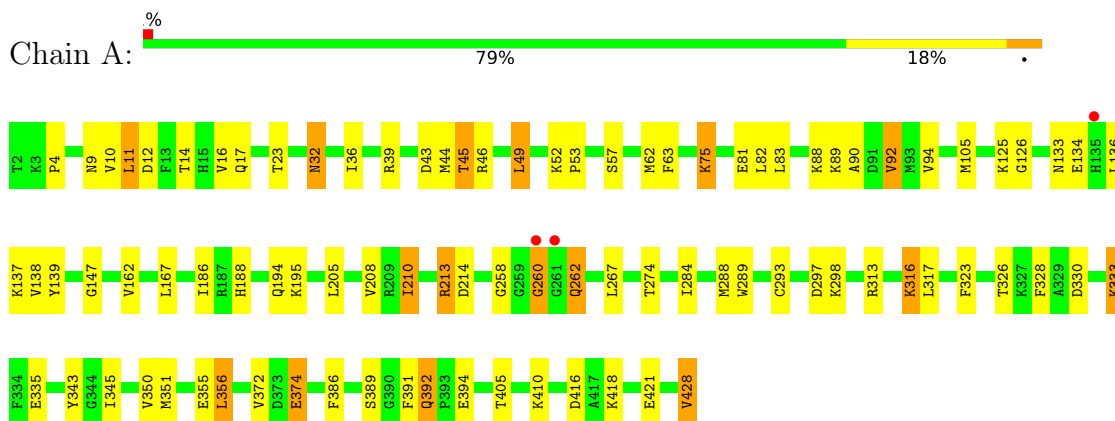
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	62	Total	O	0	0
			62	62		
4	B	68	Total	O	0	0
			68	68		

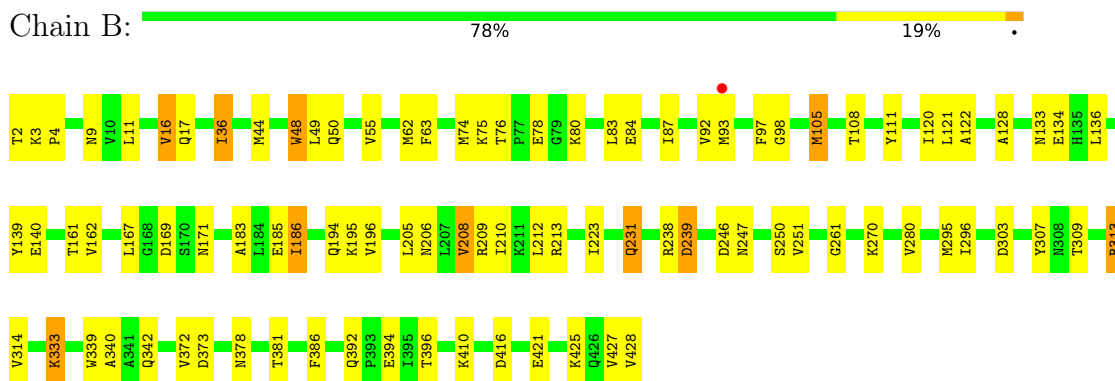
3 Residue-property plots

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

- Molecule 1: Formyl-CoA:oxalate CoA-transferase



- Molecule 1: Formyl-CoA:oxalate CoA-transferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	100.16Å 100.16Å 196.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.61 20.00 – 2.61	Depositor EDS
% Data completeness (in resolution range)	96.6 (20.00-2.61) 96.4 (20.00-2.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.60Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.200 , 0.273 0.198 , 0.272	Depositor DCC
R_{free} test set	1514 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	51.5	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 26.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6857	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.13% 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

5.1 Standard geometry

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

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	1.15	5/3387 (0.1%)	1.12	2/4580 (0.0%)
1	B	1.15	2/3387 (0.1%)	1.16	7/4580 (0.2%)
All	All	1.15	7/6774 (0.1%)	1.14	9/9160 (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	B	0	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	169	ASP	CG-OD2	10.92	1.46	1.25
1	A	372	VAL	CA-CB	6.79	1.62	1.54
1	A	260	GLY	C-O	6.59	1.32	1.23
1	B	128	ALA	CA-CB	-6.45	1.42	1.53
1	A	284	ILE	CA-C	5.96	1.57	1.52
1	A	90	ALA	CA-CB	-5.56	1.44	1.53
1	A	53	PRO	CA-C	5.03	1.59	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	48	TRP	CB-CA-C	7.92	123.31	110.16
1	B	239	ASP	N-CA-C	5.54	120.35	111.81
1	A	391	PHE	N-CA-C	5.46	116.45	108.14
1	B	208	VAL	N-CA-CB	-5.40	104.22	112.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	239	ASP	CA-CB-CG	5.38	117.98	112.60
1	B	97	PHE	CA-C-N	-5.38	113.43	121.87
1	B	97	PHE	C-N-CA	-5.38	113.43	121.87
1	A	258	GLY	N-CA-C	5.26	119.25	112.83
1	B	48	TRP	CA-CB-CG	5.25	123.57	113.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	98	GLY	Peptide

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	3313	0	3252	53	0
1	B	3313	0	3252	52	0
2	A	48	0	32	3	0
2	B	48	0	32	2	0
3	B	5	0	0	2	0
4	A	62	0	0	2	0
4	B	68	0	0	5	0
All	All	6857	0	6568	96	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 7.

All (96) 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:126:GLY:HA3	1:A:138:VAL:HG11	1.52	0.92
1:B:87:ILE:HG12	1:B:93:MET:SD	2.11	0.89
1:A:167:LEU:HD21	1:B:167:LEU:HD21	1.58	0.86
1:A:126:GLY:CA	1:A:138:VAL:HG11	2.13	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:ILE:HD13	1:A:213:ARG:HH12	1.56	0.70
1:B:17:GLN:HG3	1:B:63:PHE:CE2	2.26	0.70
1:A:186:ILE:HG21	1:A:194:GLN:HE21	1.59	0.67
1:A:11:LEU:CD2	1:A:36:ILE:HD11	2.27	0.64
1:A:17:GLN:HG3	1:A:63:PHE:CE2	2.35	0.61
1:A:32:ASN:HB2	4:A:656:HOH:O	2.01	0.60
1:A:210:ILE:HD13	1:A:213:ARG:NH1	2.16	0.60
1:A:11:LEU:CD2	1:A:36:ILE:CD1	2.80	0.59
1:A:267:LEU:HD12	1:A:328:PHE:HE2	1.68	0.58
1:B:140:GLU:N	3:B:1002:OXD:O6	2.34	0.58
1:A:44:MET:O	1:A:46:ARG:N	2.37	0.58
1:A:32:ASN:ND2	4:A:603:HOH:O	2.36	0.57
1:B:105:MET:HE1	2:B:1001:COA:O2B	2.04	0.57
1:B:48:TRP:HD1	1:B:49:LEU:HG	1.69	0.56
1:A:167:LEU:HD21	1:B:167:LEU:CD2	2.34	0.55
1:A:298:LYS:HE3	1:A:323:PHE:CZ	2.42	0.55
1:B:11:LEU:HB3	1:B:93:MET:HG2	1.89	0.55
1:A:186:ILE:HG21	1:A:194:GLN:NE2	2.22	0.54
1:A:167:LEU:CD2	1:B:167:LEU:HD21	2.33	0.54
1:A:374:GLU:CD	1:A:374:GLU:H	2.15	0.53
1:A:4:PRO:HA	1:B:186:ILE:HG13	1.91	0.53
1:A:267:LEU:HD12	1:A:328:PHE:CE2	2.43	0.53
1:A:9:ASN:ND2	1:A:89:LYS:O	2.32	0.53
1:B:11:LEU:HD11	1:B:36:ILE:HD11	1.92	0.52
1:A:11:LEU:HD21	1:A:36:ILE:HD11	1.90	0.52
1:B:44:MET:HG3	1:B:48:TRP:CZ3	2.43	0.52
1:A:10:VAL:HG13	1:A:92:VAL:HG22	1.92	0.51
1:A:333:LYS:HE2	1:A:350:VAL:O	2.11	0.51
1:B:93:MET:O	1:B:121:LEU:HD12	2.10	0.51
1:B:231:GLN:HB2	4:B:1163:HOH:O	2.11	0.51
1:B:307:TYR:O	1:B:313:ARG:HD3	2.10	0.51
1:A:186:ILE:HD13	1:A:194:GLN:HE22	1.76	0.51
1:A:205:LEU:HA	1:A:208:VAL:HG22	1.92	0.51
1:B:120:ILE:HD13	1:B:183:ALA:HB3	1.93	0.51
1:A:188:HIS:HE1	1:B:185:GLU:OE1	1.95	0.50
1:A:186:ILE:HD13	1:A:194:GLN:NE2	2.27	0.50
1:A:210:ILE:CD1	1:A:213:ARG:NH1	2.75	0.49
1:B:122:ALA:HA	1:B:196:VAL:O	2.12	0.49
1:A:289:TRP:NE1	1:A:293:CYS:SG	2.85	0.49
1:A:333:LYS:NZ	1:A:355:GLU:OE2	2.46	0.49
1:A:4:PRO:HG3	1:B:186:ILE:HD12	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:339:TRP:O	1:B:342:GLN:HB2	2.12	0.48
1:B:139:TYR:HE2	2:B:1001:COA:H143	1.79	0.47
1:A:392:GLN:HA	1:A:392:GLN:HE21	1.80	0.47
1:A:389:SER:HB3	1:B:195:LYS:HB3	1.95	0.47
1:B:9:ASN:OD1	1:B:9:ASN:C	2.57	0.47
1:B:62:MET:HE3	1:B:381:THR:HA	1.96	0.47
1:A:410:LYS:NZ	1:A:416:ASP:OD1	2.48	0.47
1:B:167:LEU:O	1:B:171:ASN:HB3	2.16	0.46
1:B:84:GLU:OE2	1:B:111:TYR:OH	2.29	0.46
1:B:139:TYR:HB3	3:B:1002:OXD:O6	2.16	0.46
1:B:295:MET:HE1	1:B:340:ALA:HB2	1.98	0.46
1:B:120:ILE:HD13	1:B:183:ALA:CB	2.46	0.46
1:A:147:GLY:O	1:A:205:LEU:HD12	2.16	0.45
1:A:405:THR:HG21	1:A:428:VAL:HG13	1.98	0.45
1:B:378:ASN:ND2	4:B:1108:HOH:O	2.50	0.45
1:B:9:ASN:HB2	4:B:1165:HOH:O	2.17	0.45
1:B:373:ASP:H	1:B:378:ASN:ND2	2.15	0.45
1:B:44:MET:HG3	1:B:48:TRP:HZ3	1.81	0.45
1:A:44:MET:O	1:A:45:THR:C	2.60	0.44
1:A:139:TYR:CZ	2:A:501:COA:H143	2.52	0.44
1:B:208:VAL:O	1:B:208:VAL:HG12	2.16	0.44
1:B:186:ILE:HG21	1:B:194:GLN:HE21	1.82	0.44
1:B:210:ILE:O	1:B:213:ARG:HB3	2.17	0.44
1:B:206:ASN:ND2	1:B:209:ARG:HH21	2.15	0.44
1:B:378:ASN:HD22	1:B:378:ASN:N	2.17	0.43
1:A:49:LEU:HD21	1:B:261:GLY:HA2	2.00	0.43
1:A:12:ASP:CG	1:A:23:THR:HG21	2.43	0.43
1:A:133:ASN:CG	1:A:136:LEU:HD12	2.44	0.43
2:A:501:COA:H131	2:A:501:COA:N4P	2.34	0.43
1:A:210:ILE:HD12	1:A:260:GLY:CA	2.48	0.43
1:B:161:THR:HG22	1:B:162:VAL:O	2.19	0.42
1:A:63:PHE:CZ	1:B:209:ARG:HD2	2.54	0.42
1:B:50:GLN:HG2	1:B:55:VAL:O	2.19	0.42
1:B:108:THR:HB	4:B:1156:HOH:O	2.19	0.42
1:A:75:LYS:HG3	2:A:501:COA:H2A	2.02	0.42
1:B:372:VAL:O	1:B:372:VAL:HG12	2.19	0.42
1:B:231:GLN:CB	4:B:1163:HOH:O	2.68	0.42
1:A:11:LEU:HD23	1:A:36:ILE:CD1	2.50	0.41
1:A:343:TYR:HB2	1:A:345:ILE:HD12	2.02	0.41
1:B:11:LEU:HD12	1:B:11:LEU:HA	1.95	0.41
1:A:213:ARG:CZ	1:A:214:ASP:OD1	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:MET:HB3	1:A:356:LEU:HD13	2.03	0.41
1:B:3:LYS:HA	1:B:4:PRO:HD3	1.94	0.41
1:A:43:ASP:O	1:A:46:ARG:HG3	2.21	0.41
1:A:62:MET:HE1	1:B:212:LEU:HB3	2.03	0.41
1:A:316:LYS:O	1:A:317:LEU:C	2.64	0.41
1:B:250:SER:O	1:B:251:VAL:C	2.63	0.41
1:B:133:ASN:HB3	1:B:136:LEU:HD12	2.03	0.40
1:B:295:MET:HE3	1:B:340:ALA:HA	2.03	0.40
1:A:14:THR:CG2	1:A:94:VAL:HG12	2.51	0.40
1:B:76:THR:HG22	1:B:78:GLU:H	1.86	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	425/427 (100%)	403 (95%)	18 (4%)	4 (1%)	14	28
1	B	425/427 (100%)	400 (94%)	21 (5%)	4 (1%)	14	28
All	All	850/854 (100%)	803 (94%)	39 (5%)	8 (1%)	14	28

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	333	LYS
1	A	16	VAL
1	B	16	VAL
1	A	45	THR
1	B	134	GLU
1	A	134	GLU
1	A	262	GLN
1	B	314	VAL

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	350/350 (100%)	313 (89%)	37 (11%)	6	13
1	B	350/350 (100%)	316 (90%)	34 (10%)	8	16
All	All	700/700 (100%)	629 (90%)	71 (10%)	7	14

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	32	ASN
1	A	39	ARG
1	A	49	LEU
1	A	52	LYS
1	A	57	SER
1	A	75	LYS
1	A	81	GLU
1	A	82	LEU
1	A	83	LEU
1	A	88	LYS
1	A	92	VAL
1	A	105	MET
1	A	125	LYS
1	A	137	LYS
1	A	162	VAL
1	A	195	LYS
1	A	210	ILE
1	A	213	ARG
1	A	262	GLN
1	A	274	THR
1	A	288	MET
1	A	297	ASP
1	A	313	ARG
1	A	316	LYS
1	A	326	THR
1	A	330	ASP

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Mol	Chain	Res	Type
1	A	333	LYS
1	A	335	GLU
1	A	356	LEU
1	A	374	GLU
1	A	386	PHE
1	A	392	GLN
1	A	394	GLU
1	A	418	LYS
1	A	421	GLU
1	A	428	VAL
1	B	2	THR
1	B	16	VAL
1	B	36	ILE
1	B	74	MET
1	B	75	LYS
1	B	80	LYS
1	B	83	LEU
1	B	92	VAL
1	B	105	MET
1	B	186	ILE
1	B	205	LEU
1	B	223	ILE
1	B	231	GLN
1	B	238	ARG
1	B	239	ASP
1	B	246	ASP
1	B	247	ASN
1	B	270	LYS
1	B	280	VAL
1	B	296	ILE
1	B	303	ASP
1	B	309	THR
1	B	313	ARG
1	B	333	LYS
1	B	386	PHE
1	B	392	GLN
1	B	394	GLU
1	B	396	THR
1	B	410	LYS
1	B	416	ASP
1	B	421	GLU
1	B	425	LYS

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Mol	Chain	Res	Type
1	B	427	VAL
1	B	428	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	32	ASN
1	A	188	HIS
1	A	194	GLN
1	A	342	GLN
1	A	392	GLN
1	B	188	HIS
1	B	206	ASN
1	B	231	GLN
1	B	241	ASN
1	B	262	GLN
1	B	358	HIS
1	B	378	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](#)

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
3	OXD	B	1002	1,2	4,4,5	1.89	1 (25%)	3,4,6	0.65	0
2	COA	A	501	-	47,50,50	2.02	6 (12%)	69,75,75	1.57	9 (13%)
2	COA	B	1001	3	47,50,50	1.69	5 (10%)	69,75,75	1.71	14 (20%)

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	OXD	B	1002	1,2	-	0/0/2/4	-
2	COA	A	501	-	-	11/48/64/64	0/3/3/3
2	COA	B	1001	3	-	6/48/64/64	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	COA	O9P-C9P	10.32	1.43	1.23
2	B	1001	COA	O9P-C9P	9.05	1.40	1.23
2	A	501	COA	P1A-O3A	4.26	1.64	1.59
2	A	501	COA	P2A-O3A	4.19	1.64	1.59
3	B	1002	OXD	O6-C2	-3.58	1.21	1.30
2	A	501	COA	C2A-N3A	3.04	1.39	1.33
2	B	1001	COA	P2A-O3A	2.93	1.62	1.59
2	B	1001	COA	C2A-N1A	2.87	1.39	1.33
2	A	501	COA	C2A-N1A	2.86	1.39	1.33
2	B	1001	COA	C2A-N3A	2.83	1.39	1.33
2	A	501	COA	C8A-N7A	2.46	1.36	1.31
2	B	1001	COA	P1A-O3A	2.14	1.61	1.59

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	COA	N3A-C2A-N1A	-5.72	119.92	128.58
2	B	1001	COA	C5A-C4A-N3A	-5.70	118.87	126.72
2	A	501	COA	C5A-C4A-N3A	-5.06	119.74	126.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	COA	C5A-N7A-C8A	4.09	109.88	103.45
2	B	1001	COA	N3A-C2A-N1A	-4.06	122.44	128.58
2	B	1001	COA	C4A-C5A-N7A	-3.84	106.19	110.58
2	B	1001	COA	N9A-C8A-N7A	-3.55	108.90	113.94
2	A	501	COA	C2A-N3A-C4A	3.48	120.33	111.83
2	A	501	COA	C5A-N7A-C8A	3.28	108.60	103.45
2	B	1001	COA	C2A-N3A-C4A	3.24	119.74	111.83
2	B	1001	COA	N3A-C4A-N9A	3.23	132.65	127.17
2	A	501	COA	N3A-C4A-N9A	3.18	132.57	127.17
2	B	1001	COA	C6A-C5A-C4A	3.15	121.48	117.18
2	A	501	COA	N9A-C8A-N7A	-3.02	109.65	113.94
2	B	1001	COA	O9A-P3B-O8A	2.97	118.93	107.80
2	A	501	COA	CDP-CBP-CCP	-2.58	103.97	108.22
2	A	501	COA	C4A-C5A-N7A	-2.58	107.64	110.58
2	B	1001	COA	C5A-C4A-N9A	2.44	108.47	105.81
2	B	1001	COA	CDP-CBP-CCP	-2.43	104.22	108.22
2	B	1001	COA	O4B-C1B-N9A	2.26	112.43	108.09
2	B	1001	COA	O2A-P1A-O3A	2.23	113.31	107.27
2	B	1001	COA	CDP-CBP-CAP	2.13	112.40	108.77
2	A	501	COA	C6A-C5A-C4A	2.07	120.00	117.18

There are no chirality outliers.

All (17) torsion outliers are listed below:

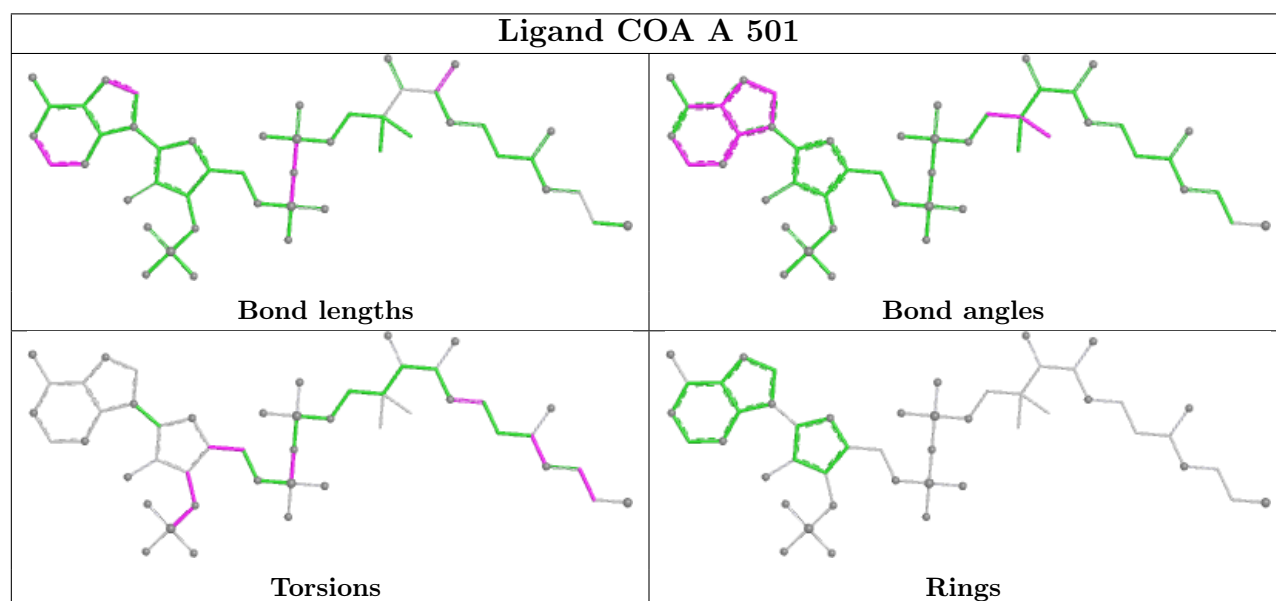
Mol	Chain	Res	Type	Atoms
2	A	501	COA	S1P-C2P-C3P-N4P
2	B	1001	COA	C5P-C6P-C7P-N8P
2	B	1001	COA	S1P-C2P-C3P-N4P
2	A	501	COA	O5P-C5P-N4P-C3P
2	A	501	COA	C6P-C5P-N4P-C3P
2	B	1001	COA	C6P-C5P-N4P-C3P
2	B	1001	COA	O5P-C5P-N4P-C3P
2	A	501	COA	C3B-C4B-C5B-O5B
2	A	501	COA	C6P-C7P-N8P-C9P
2	A	501	COA	O4B-C4B-C5B-O5B
2	B	1001	COA	P2A-O3A-P1A-O2A
2	A	501	COA	C4B-C3B-O3B-P3B
2	A	501	COA	C3B-O3B-P3B-O7A
2	A	501	COA	P2A-O3A-P1A-O2A
2	A	501	COA	C2B-C3B-O3B-P3B
2	B	1001	COA	P2A-O3A-P1A-O1A
2	A	501	COA	P2A-O3A-P1A-O1A

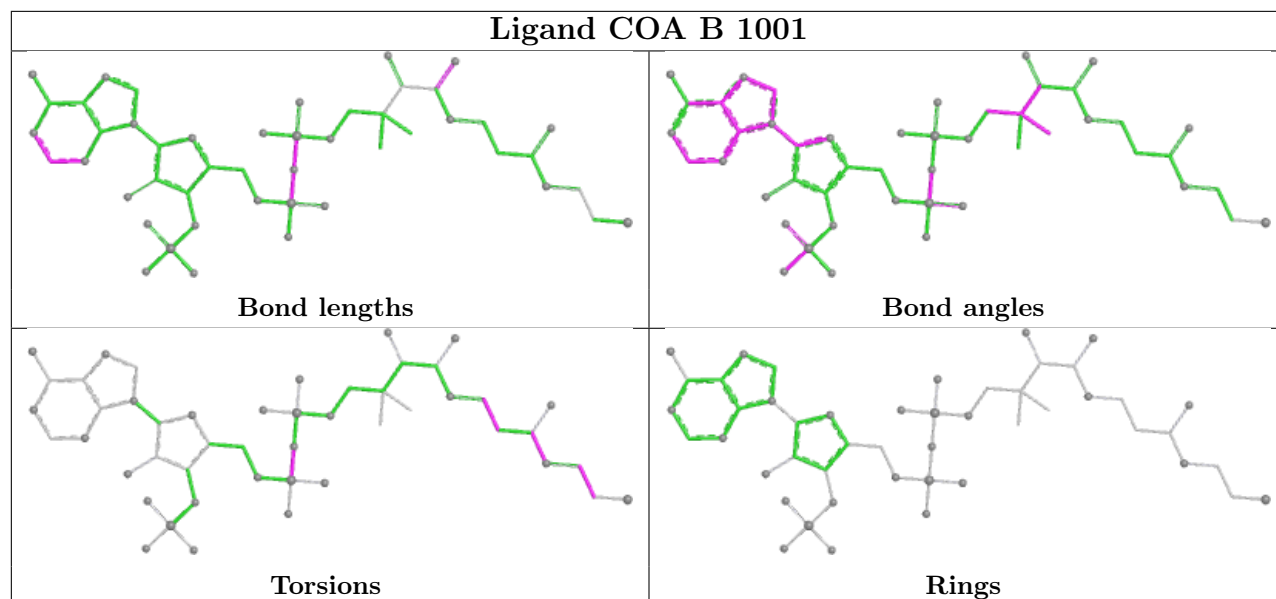
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1002	OXD	2	0
2	A	501	COA	3	0
2	B	1001	COA	2	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	427/427 (100%)	-0.47	3 (0%) 84 82	16, 28, 56, 64	0
1	B	426/427 (99%)	-0.47	1 (0%) 91 90	12, 28, 56, 64	0
All	All	853/854 (99%)	-0.47	4 (0%) 87 85	12, 28, 56, 64	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	135	HIS	2.6
1	B	93	MET	2.5
1	A	260	GLY	2.2
1	A	261	GLY	2.1

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.

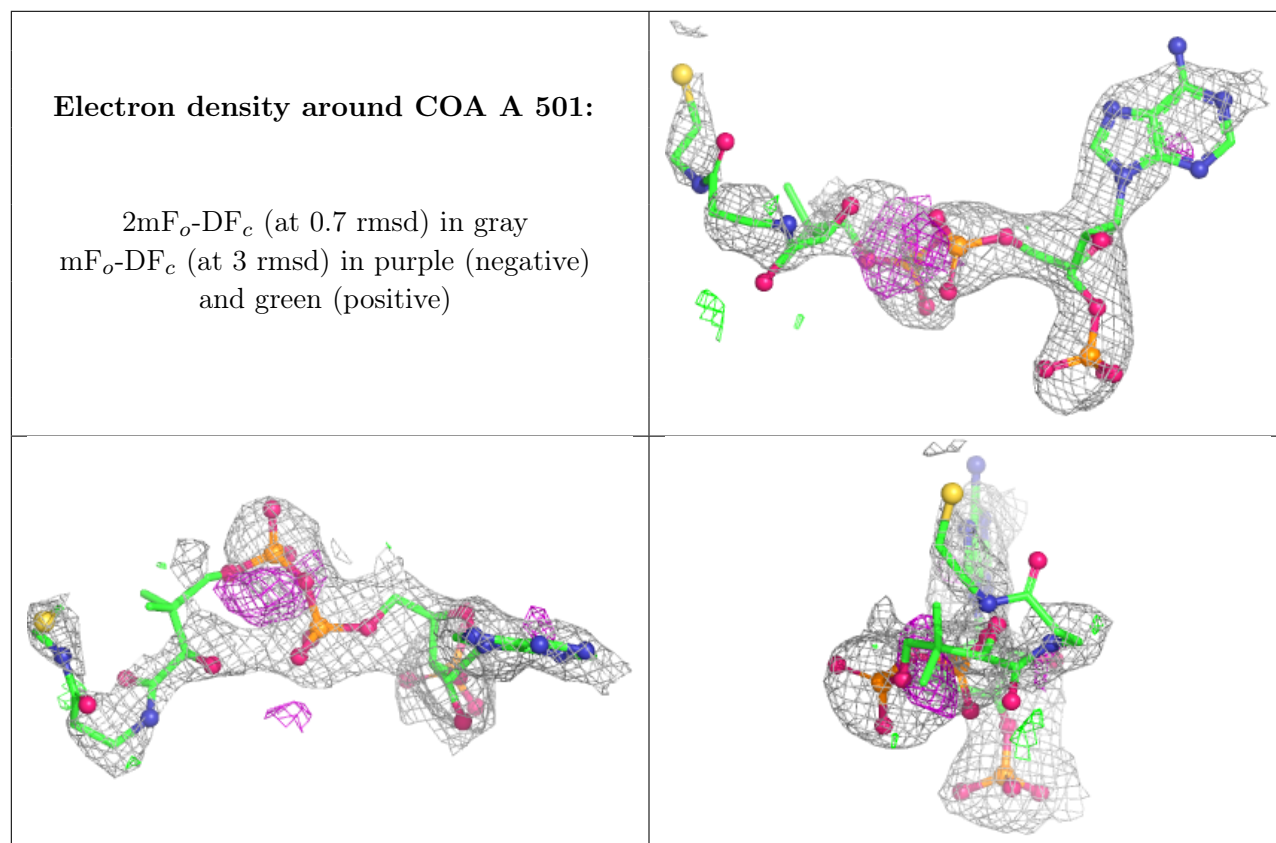
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	COA	A	501	48/48	0.77	0.14	98,102,106,107	0

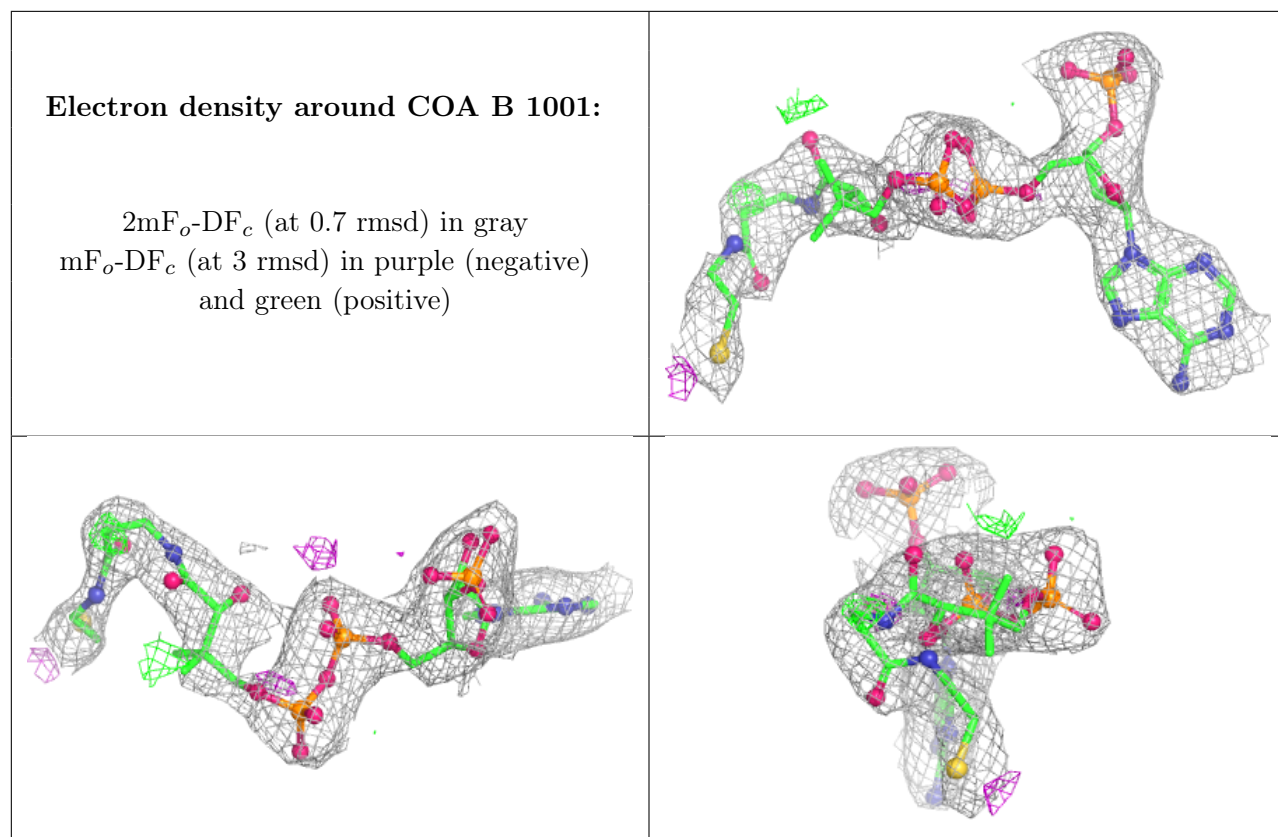
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	OXD	B	1002	5/6	0.80	0.27	40,44,47,50	0
2	COA	B	1001	48/48	0.86	0.11	62,75,88,89	0

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.