



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:23 AM UTC

PDB ID : 6SLL / pdb_00006sll
Title : Diaminobutyrate acetyltransferase EctA from *Paenibacillus lautus* in complex with its substrate L-2,4-diaminobutyric acid (DAB) and coenzyme A
Authors : Richter, A.A.; Kobus, S.; Czech, L.; Hoepfner, A.; Bremer, E.; Smits, S.H.J.
Deposited on : 2019-08-20
Resolution : 1.20 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

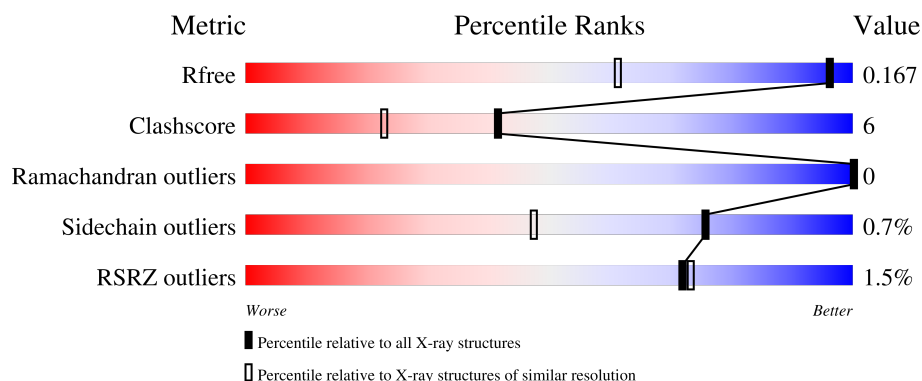
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.20 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	1216 (1.20-1.20)
Clashscore	190562	1265 (1.20-1.20)
Ramachandran outliers	187476	1226 (1.20-1.20)
Sidechain outliers	187428	1226 (1.20-1.20)
RSRZ outliers	180081	1214 (1.20-1.20)

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	186	% 75% 11% .. 12%
1	B	186	2% 76% 11% .. 11%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3398 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 L-2,4-diaminobutyric acid acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	164	Total	C	N	O	S	0	12	0
			1363	865	232	256	10			
1	B	166	Total	C	N	O	S	0	13	0
			1379	873	236	259	11			

There are 32 discrepancies between the modelled and reference sequences:

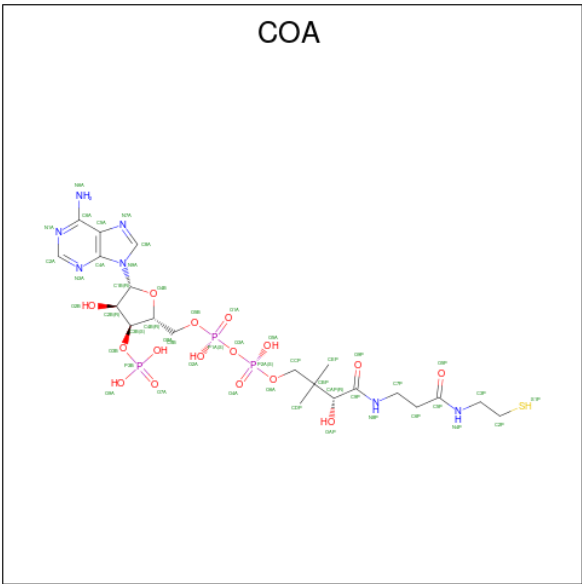
Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	TRP	-	expression tag	UNP D3EKC1
A	-8	SER	-	expression tag	UNP D3EKC1
A	-7	HIS	-	expression tag	UNP D3EKC1
A	-6	PRO	-	expression tag	UNP D3EKC1
A	-5	GLN	-	expression tag	UNP D3EKC1
A	-4	PHE	-	expression tag	UNP D3EKC1
A	-3	GLU	-	expression tag	UNP D3EKC1
A	-2	LYS	-	expression tag	UNP D3EKC1
A	-1	SER	-	expression tag	UNP D3EKC1
A	0	GLY	-	expression tag	UNP D3EKC1
A	171	GLY	-	expression tag	UNP D3EKC1
A	172	SER	-	expression tag	UNP D3EKC1
A	173	ALA	-	expression tag	UNP D3EKC1
A	174	TRP	-	expression tag	UNP D3EKC1
A	175	SER	-	expression tag	UNP D3EKC1
A	176	HIS	-	expression tag	UNP D3EKC1
B	-9	TRP	-	expression tag	UNP D3EKC1
B	-8	SER	-	expression tag	UNP D3EKC1
B	-7	HIS	-	expression tag	UNP D3EKC1
B	-6	PRO	-	expression tag	UNP D3EKC1
B	-5	GLN	-	expression tag	UNP D3EKC1
B	-4	PHE	-	expression tag	UNP D3EKC1
B	-3	GLU	-	expression tag	UNP D3EKC1
B	-2	LYS	-	expression tag	UNP D3EKC1
B	-1	SER	-	expression tag	UNP D3EKC1

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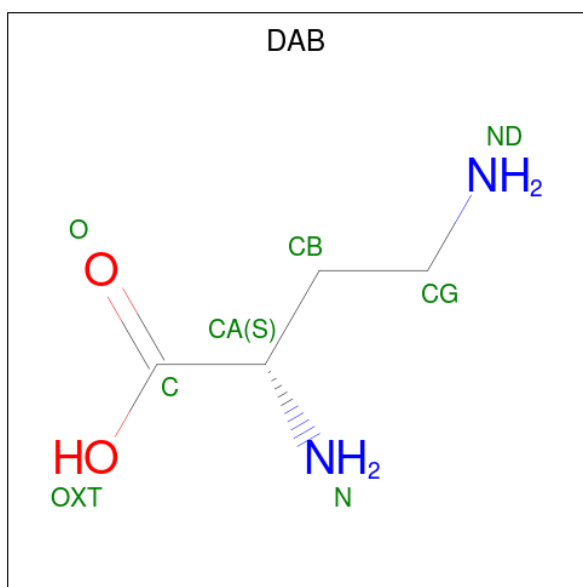
Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP D3EKC1
B	171	GLY	-	expression tag	UNP D3EKC1
B	172	SER	-	expression tag	UNP D3EKC1
B	173	ALA	-	expression tag	UNP D3EKC1
B	174	TRP	-	expression tag	UNP D3EKC1
B	175	SER	-	expression tag	UNP D3EKC1
B	176	HIS	-	expression tag	UNP D3EKC1

- Molecule 2 is COENZYME A (CCD ID: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is 2,4-DIAMINO BUTYRIC ACID (CCD ID: DAB) (formula: C₄H₁₀N₂O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	4	2	2		
3	B	1	Total	C	N	O	0	0
			8	4	2	2		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

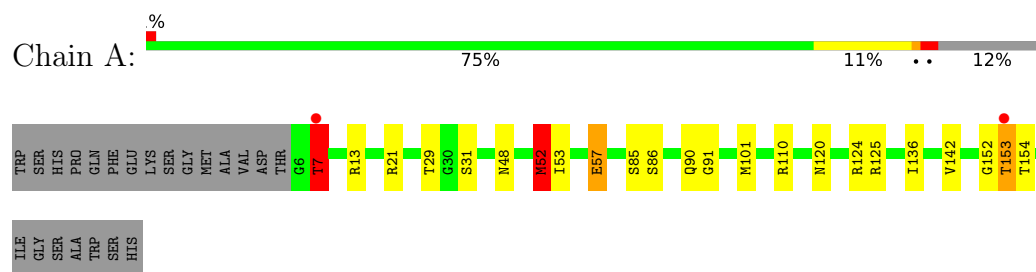
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	277	Total	O	0	0
			277	277		
5	B	266	Total	O	0	0
			266	266		

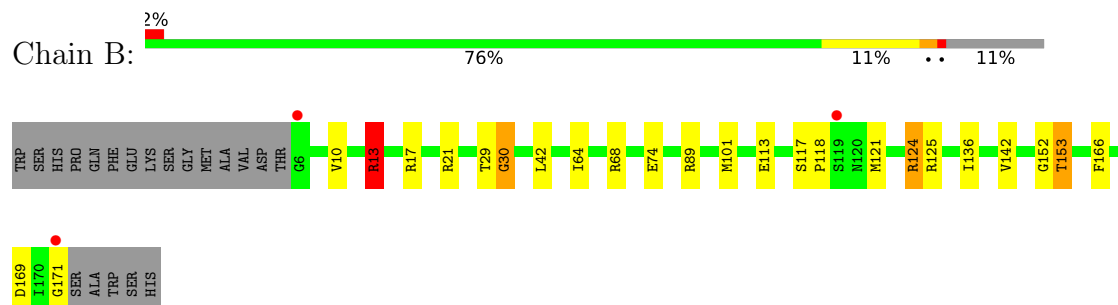
3 Residue-property plots [i](#)

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: L-2,4-diaminobutyric acid acetyltransferase



- Molecule 1: L-2,4-diaminobutyric acid acetyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	151.05Å 151.05Å 46.17Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	75.52 – 1.20 75.52 – 1.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (75.52-1.20) 99.5 (75.52-1.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 1.20Å)	Xtriage
Refinement program	REFMAC 5.8.0131	Depositor
R, R_{free}	0.122 , 0.150 (Not available) , 0.167	Depositor DCC
R_{free} test set	6191 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	12.7	Xtriage
Anisotropy	0.353	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.026 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	3398	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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, MG, DAB

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	1.73	16/1419 (1.1%)	1.33	10/1923 (0.5%)
1	B	2.04	28/1440 (1.9%)	1.59	19/1952 (1.0%)
All	All	1.89	44/2859 (1.5%)	1.47	29/3875 (0.7%)

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
1	B	0	5
All	All	0	6

All (44) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	13[A]	ARG	CZ-NH1	-22.03	1.01	1.32
1	B	13[B]	ARG	CZ-NH1	-22.03	1.01	1.32
1	B	30	GLY	C-O	12.82	1.36	1.23
1	B	101	MET	SD-CE	-9.97	1.54	1.79
1	A	124	ARG	CZ-NH2	9.09	1.45	1.33
1	A	101	MET	SD-CE	-9.07	1.56	1.79
1	B	10	VAL	CA-CB	-8.86	1.43	1.54
1	B	89	ARG	CD-NE	8.23	1.57	1.46
1	B	13[A]	ARG	CD-NE	7.52	1.56	1.46
1	B	13[B]	ARG	CD-NE	7.52	1.56	1.46
1	B	125	ARG	NE-CZ	7.52	1.41	1.33
1	B	89	ARG	NE-CZ	7.46	1.41	1.33
1	B	21	ARG	CD-NE	-7.17	1.36	1.46
1	A	154	THR	N-CA	7.09	1.54	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	68	ARG	CG-CD	-7.07	1.31	1.52
1	B	142	VAL	CA-C	6.94	1.61	1.52
1	A	13	ARG	CZ-NH2	6.81	1.42	1.33
1	A	90	GLN	CD-OE1	6.78	1.36	1.23
1	B	117[A]	SER	N-CA	-6.64	1.36	1.46
1	B	117[B]	SER	N-CA	-6.64	1.36	1.46
1	B	64	ILE	CB-CG1	-6.22	1.41	1.53
1	B	121[A]	MET	N-CA	-6.21	1.38	1.46
1	B	121[B]	MET	N-CA	-6.21	1.38	1.46
1	B	118	PRO	C-O	-6.10	1.16	1.24
1	B	74	GLU	CD-OE2	6.03	1.36	1.25
1	B	30	GLY	CA-C	6.02	1.64	1.51
1	B	124	ARG	NE-CZ	-5.93	1.26	1.33
1	B	29	THR	N-CA	5.88	1.53	1.46
1	A	31	SER	C-N	-5.78	1.25	1.33
1	A	13	ARG	CZ-NH1	5.75	1.40	1.32
1	A	52[A]	MET	CG-SD	-5.67	1.66	1.80
1	A	52[B]	MET	CG-SD	-5.67	1.66	1.80
1	A	124	ARG	NE-CZ	5.66	1.39	1.33
1	A	86	SER	CA-CB	5.55	1.63	1.53
1	B	153	THR	CB-CG2	5.40	1.70	1.52
1	A	91	GLY	CA-C	5.21	1.58	1.51
1	A	166	PHE	CG-CD2	-5.18	1.27	1.38
1	B	171	GLY	CA-C	5.16	1.61	1.52
1	A	57	GLU	CA-C	5.16	1.59	1.53
1	A	125	ARG	NE-CZ	5.10	1.38	1.33
1	B	42	LEU	CB-CG	-5.08	1.43	1.53
1	A	153	THR	C-O	5.05	1.30	1.23
1	B	142	VAL	C-N	-5.00	1.26	1.33
1	B	30	GLY	C-N	-5.00	1.26	1.33

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	13[A]	ARG	NE-CZ-NH2	-14.51	106.14	119.20
1	B	13[B]	ARG	NE-CZ-NH2	-14.51	106.14	119.20
1	B	13[A]	ARG	NE-CZ-NH1	-12.30	109.20	121.50
1	B	13[B]	ARG	NE-CZ-NH1	-12.30	109.20	121.50
1	B	13[A]	ARG	NH1-CZ-NH2	-9.19	107.36	119.30
1	B	13[B]	ARG	NH1-CZ-NH2	-9.19	107.36	119.30
1	B	21	ARG	NE-CZ-NH1	8.71	130.21	121.50
1	B	121[A]	MET	N-CA-CB	-7.00	99.81	110.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	121[B]	MET	N-CA-CB	-7.00	99.81	110.16
1	A	7	THR	CB-CA-C	6.62	120.60	109.48
1	A	7	THR	CA-CB-CG2	6.60	121.72	110.50
1	A	142	VAL	N-CA-C	6.29	117.03	110.62
1	B	113	GLU	CA-C-O	-5.94	114.41	120.71
1	B	124	ARG	NE-CZ-NH2	-5.93	113.87	119.20
1	B	101	MET	CG-SD-CE	-5.83	88.09	100.90
1	B	68	ARG	CB-CG-CD	5.82	124.69	111.30
1	B	169	ASP	CA-CB-CG	-5.71	106.89	112.60
1	B	21	ARG	NE-CZ-NH2	-5.64	114.12	119.20
1	B	10	VAL	CB-CA-C	5.53	119.21	110.81
1	B	124	ARG	CD-NE-CZ	-5.49	116.71	124.40
1	B	142	VAL	N-CA-C	5.47	116.86	110.62
1	A	91	GLY	O-C-N	5.45	128.83	122.45
1	B	30	GLY	O-C-N	-5.28	116.18	122.31
1	A	152	GLY	N-CA-C	-5.26	108.36	115.36
1	A	7	THR	OG1-CB-CG2	5.22	119.74	109.30
1	A	168	ASN	CB-CA-C	-5.18	102.63	111.02
1	A	85	SER	CB-CA-C	-5.11	100.55	110.46
1	A	153	THR	CA-C-O	5.08	128.28	121.78
1	A	29	THR	N-CA-CB	5.02	117.95	110.22

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	166	PHE	Sidechain
1	B	13[A]	ARG	Sidechain
1	B	13[B]	ARG	Sidechain
1	B	152	GLY	Peptide
1	B	153	THR	Mainchain
1	B	30	GLY	Mainchain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1363	0	1335	24	0
1	B	1379	0	1354	12	0
2	A	48	0	32	2	0
2	B	48	0	31	0	0
3	A	8	0	9	0	0
3	B	8	0	9	0	0
4	A	1	0	0	0	0
5	A	277	0	0	8	0
5	B	266	0	0	5	0
All	All	3398	0	2770	36	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 6.

All (36) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:ARG:HB3	5:B:490:HOH:O	1.39	1.17
1:A:120[B]:ASN:HD21	2:A:201:COA:H32	1.21	1.01
1:A:120[B]:ASN:ND2	2:A:201:COA:H32	1.84	0.91
1:B:17:ARG:CB	5:B:490:HOH:O	2.05	0.89
1:A:136[B]:ILE:CD1	1:A:166:PHE:CE1	2.57	0.87
1:A:52[A]:MET:HG3	1:A:53:ILE:N	2.01	0.74
1:B:136[B]:ILE:CD1	1:B:166:PHE:CE1	2.74	0.70
1:A:136[B]:ILE:CD1	1:A:166:PHE:HE1	2.04	0.70
1:A:7:THR:HG23	1:A:57:GLU:OE2	1.92	0.70
1:A:136[B]:ILE:HD11	1:A:166:PHE:HE1	1.58	0.67
1:A:48[B]:ASN:ND2	5:A:304:HOH:O	2.30	0.65
1:A:136[B]:ILE:HD13	1:A:166:PHE:CD1	2.32	0.65
1:A:136[B]:ILE:CD1	1:A:166:PHE:CD1	2.84	0.60
1:A:110[B]:ARG:NH1	5:A:303:HOH:O	2.30	0.59
1:A:110[B]:ARG:NH2	5:A:303:HOH:O	2.29	0.58
1:B:136[B]:ILE:HD11	1:B:166:PHE:CE1	2.41	0.55
1:A:52[A]:MET:CG	1:A:53:ILE:N	2.69	0.55
1:B:124:ARG:HD3	5:B:477:HOH:O	2.07	0.53
1:A:136[B]:ILE:HD13	1:A:166:PHE:HD1	1.73	0.52
1:A:110[B]:ARG:CZ	5:A:303:HOH:O	2.60	0.50
1:B:13[B]:ARG:HG2	5:B:439:HOH:O	2.11	0.50
1:B:13[B]:ARG:CG	5:B:439:HOH:O	2.59	0.50
1:A:153:THR:OG1	5:A:301:HOH:O	2.20	0.47
1:B:136[B]:ILE:CD1	1:B:166:PHE:CD1	2.98	0.46
1:A:136[B]:ILE:HD11	1:A:166:PHE:CE1	2.38	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136[B]:ILE:HD12	1:A:166:PHE:CE1	2.47	0.45
1:A:21:ARG:HD3	5:A:355:HOH:O	2.17	0.44
1:A:21:ARG:NH1	5:A:306:HOH:O	2.42	0.43
1:A:153:THR:HG21	5:A:497:HOH:O	2.19	0.42
1:A:7:THR:CG2	1:A:57:GLU:OE2	2.66	0.41

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	174/186 (94%)	172 (99%)	2 (1%)	0	100	100
1	B	177/186 (95%)	173 (98%)	4 (2%)	0	100	100
All	All	351/372 (94%)	345 (98%)	6 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	151/157 (96%)	148 (98%)	3 (2%)	48	12
1	B	153/157 (98%)	153 (100%)	0	100	100
All	All	304/314 (97%)	301 (99%)	3 (1%)	76	36

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	THR
1	A	52[A]	MET
1	A	52[B]	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	103	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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	COA	A	201	-	47,50,50	1.72	12 (25%)	69,75,75	1.57	11 (15%)
3	DAB	B	202	-	6,7,7	1.05	1 (16%)	4,8,8	0.45	0
3	DAB	A	202	-	6,7,7	1.33	1 (16%)	4,8,8	1.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	COA	B	201	-	47,50,50	1.60	8 (17%)	69,75,75	1.51	13 (18%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	COA	A	201	-	-	2/48/64/64	0/3/3/3
3	DAB	B	202	-	-	0/7/7/7	-
3	DAB	A	202	-	-	0/7/7/7	-
2	COA	B	201	-	-	1/48/64/64	0/3/3/3

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	201	COA	C5P-N4P	-4.97	1.21	1.33
2	A	201	COA	C4A-N9A	-4.51	1.28	1.37
2	A	201	COA	C8A-N9A	-4.37	1.30	1.37
2	A	201	COA	C3P-N4P	-4.08	1.37	1.46
2	A	201	COA	C5B-C4B	3.49	1.62	1.51
2	B	201	COA	C7P-N8P	-3.45	1.38	1.46
2	B	201	COA	C7P-C6P	3.19	1.62	1.51
2	A	201	COA	C4A-N3A	3.07	1.40	1.34
2	B	201	COA	C3P-N4P	2.93	1.52	1.46
2	B	201	COA	O5P-C5P	2.68	1.28	1.23
2	A	201	COA	C2B-C1B	2.49	1.61	1.53
3	A	202	DAB	OXT-C	-2.47	1.22	1.30
2	A	201	COA	OAP-CAP	2.42	1.46	1.42
3	B	202	DAB	OXT-C	-2.40	1.23	1.30
2	A	201	COA	C2A-N3A	2.34	1.38	1.33
2	A	201	COA	C3B-C4B	2.30	1.58	1.52
2	A	201	COA	P3B-O3B	2.28	1.63	1.59
2	B	201	COA	CEP-CBP	-2.26	1.49	1.53
2	B	201	COA	C1B-N9A	-2.21	1.40	1.46
2	B	201	COA	P2A-O3A	2.19	1.61	1.59
2	A	201	COA	O4B-C1B	-2.16	1.37	1.42
2	A	201	COA	C5A-C4A	2.00	1.42	1.39

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	COA	O9A-P3B-O8A	4.67	125.30	107.80
2	A	201	COA	O3B-C3B-C4B	-4.11	95.54	110.03
2	B	201	COA	O9A-P3B-O7A	-3.97	95.38	110.83
2	A	201	COA	P3B-O3B-C3B	-3.81	113.25	123.43
2	A	201	COA	N9A-C8A-N7A	3.79	119.31	113.94
2	A	201	COA	O9P-C9P-N8P	-3.72	115.10	122.98
2	A	201	COA	C5A-C4A-N3A	-3.53	121.86	126.72
2	A	201	COA	C5A-N7A-C8A	-3.28	98.29	103.45
2	B	201	COA	C7P-N8P-C9P	3.24	128.37	122.55
2	B	201	COA	O3B-P3B-O7A	-2.96	98.78	109.33
2	B	201	COA	O2B-C2B-C3B	2.68	118.69	111.19
2	B	201	COA	O4B-C4B-C5B	2.54	117.47	109.33
2	B	201	COA	C2P-C3P-N4P	-2.54	106.54	112.31
2	A	201	COA	N3A-C2A-N1A	-2.51	124.78	128.58
2	B	201	COA	C4A-N9A-C8A	2.49	108.35	105.74
2	B	201	COA	O8A-P3B-O7A	2.47	120.47	110.83
2	B	201	COA	N9A-C8A-N7A	-2.38	110.56	113.94
2	A	201	COA	C5B-C4B-C3B	-2.36	106.51	114.38
2	B	201	COA	C4A-C5A-N7A	-2.32	107.92	110.58
2	A	201	COA	CEP-CBP-CCP	2.32	112.05	108.22
2	A	201	COA	C4B-O4B-C1B	2.28	114.49	109.47
2	A	201	COA	C2P-C3P-N4P	2.21	117.32	112.31
2	B	201	COA	C5A-C4A-N3A	-2.06	123.89	126.72
2	B	201	COA	C6P-C5P-N4P	2.01	120.01	116.34

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	201	COA	CCP-O6A-P2A-O4A
2	B	201	COA	C3B-O3B-P3B-O9A
2	A	201	COA	S1P-C2P-C3P-N4P

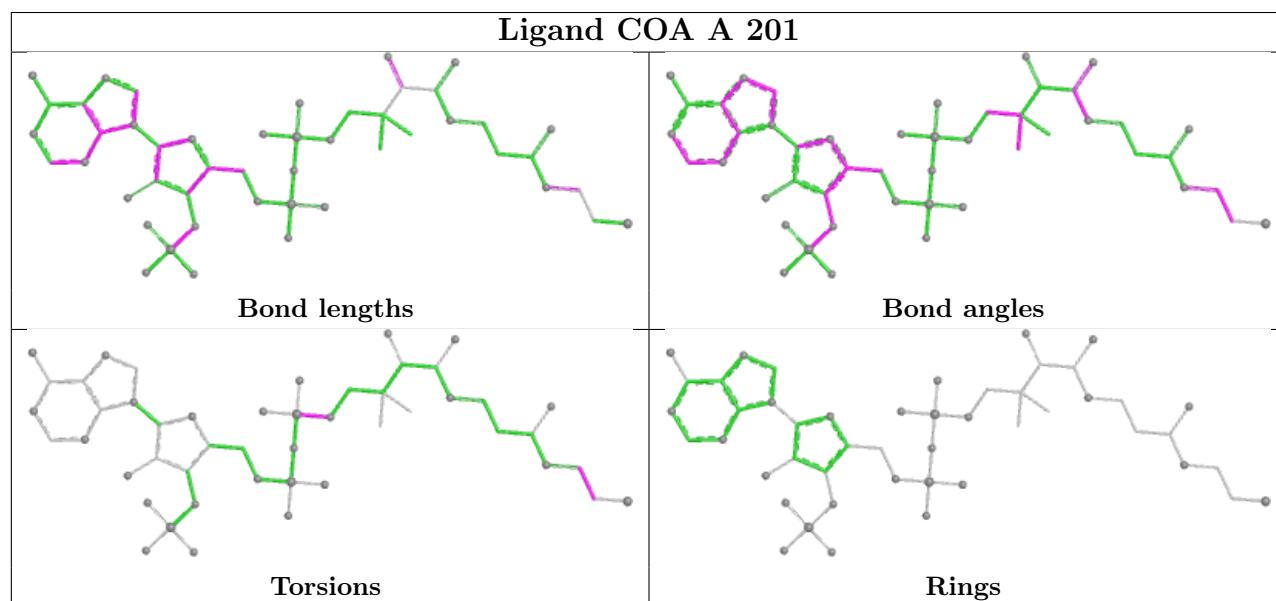
There are no ring outliers.

1 monomer is involved in 2 short contacts:

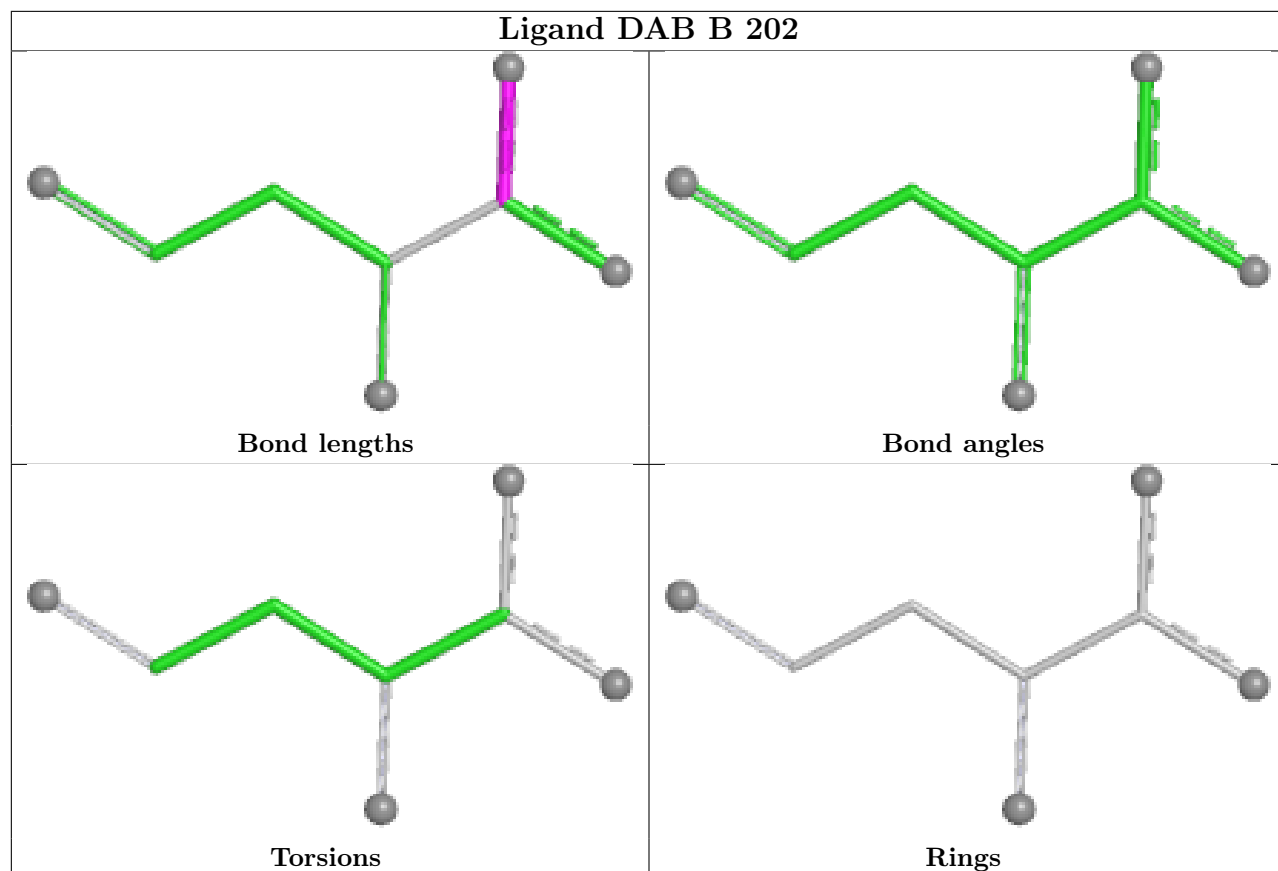
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201	COA	2	0

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

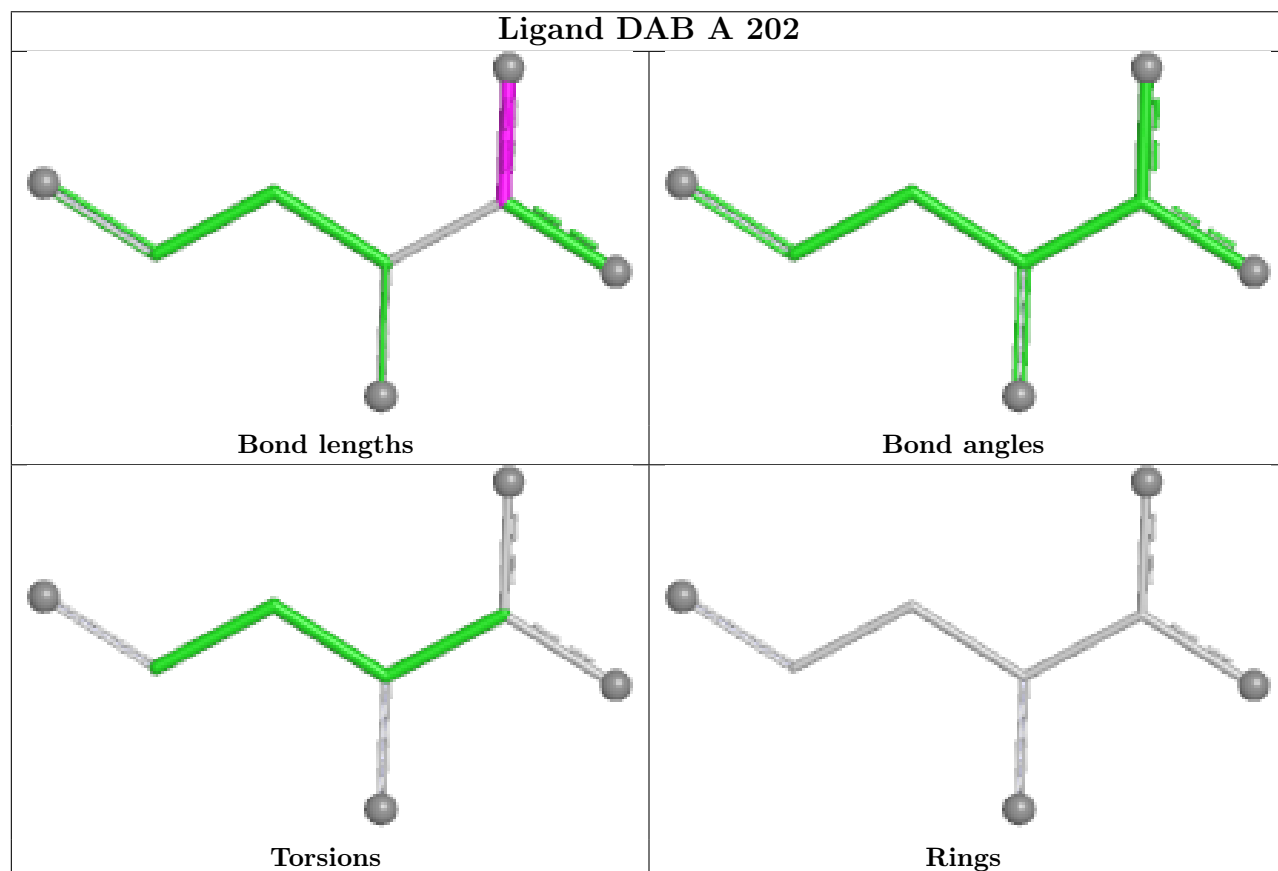
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted 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.

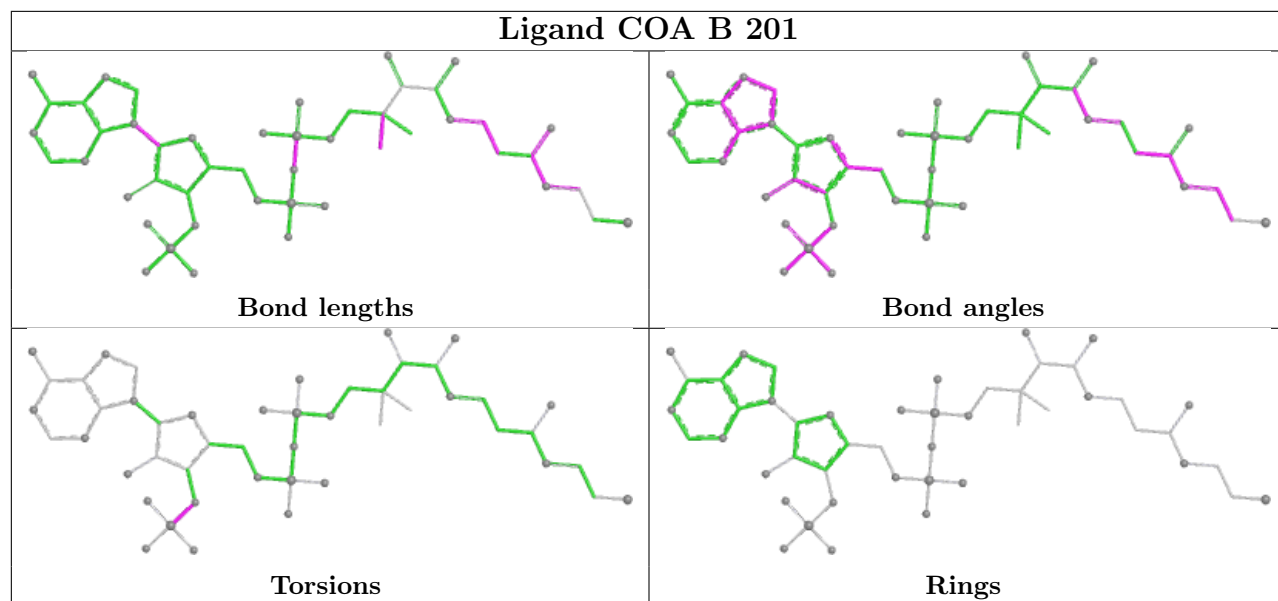


Ligand DAB B 202



Ligand DAB A 202





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	164/186 (88%)	-0.25	2 (1%) 76 78	5, 13, 31, 51	12 (7%)
1	B	166/186 (89%)	-0.18	3 (1%) 67 68	6, 14, 27, 51	13 (7%)
All	All	330/372 (88%)	-0.22	5 (1%) 72 73	5, 13, 31, 51	25 (7%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	153	THR	3.2
1	B	119	SER	2.8
1	B	6	GLY	2.5
1	B	171	GLY	2.1
1	A	7	THR	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](#)

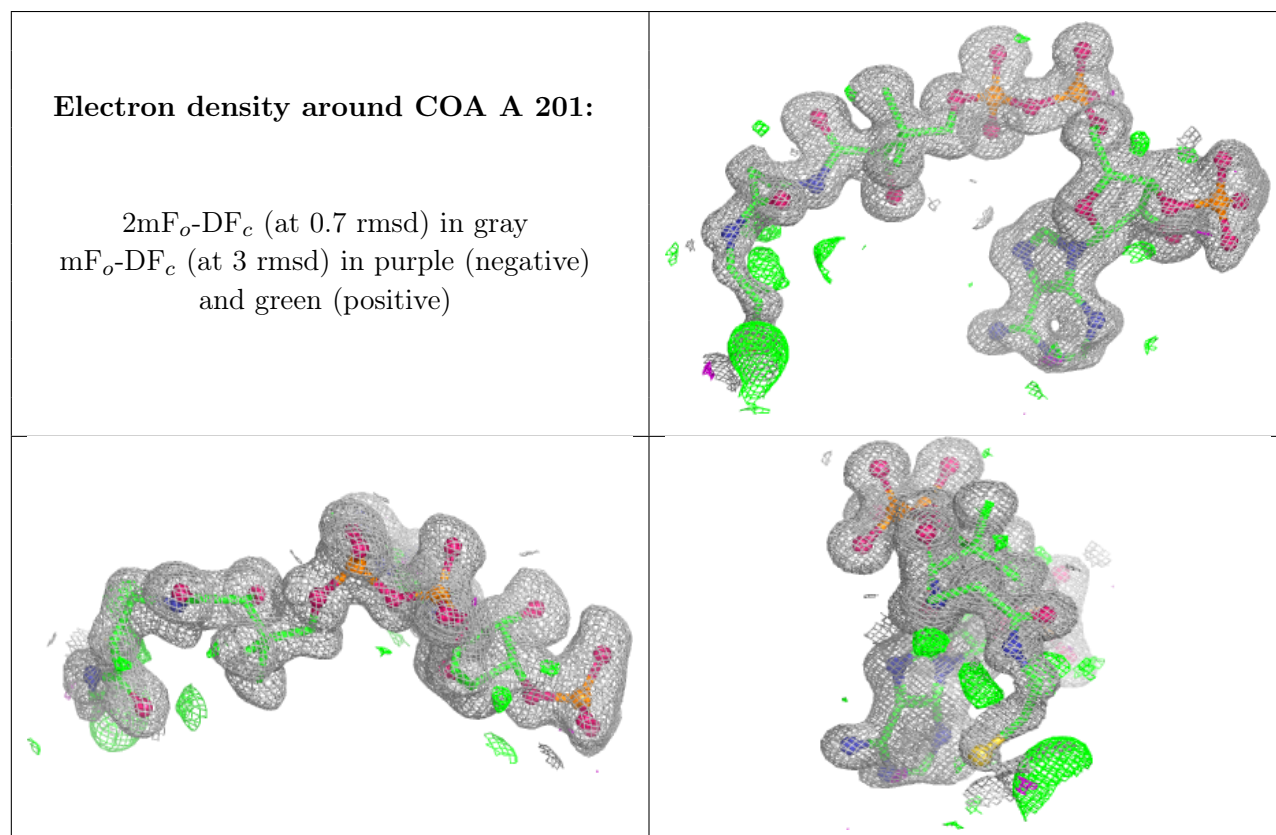
There are no oligosaccharides in this entry.

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.

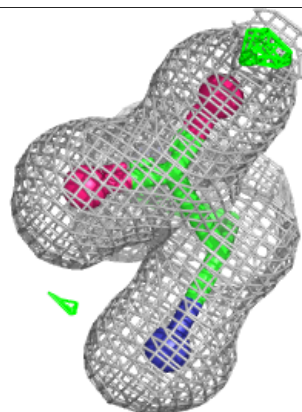
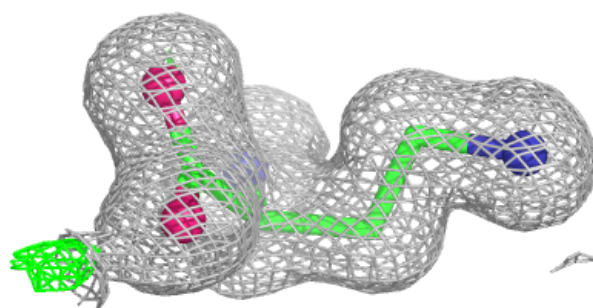
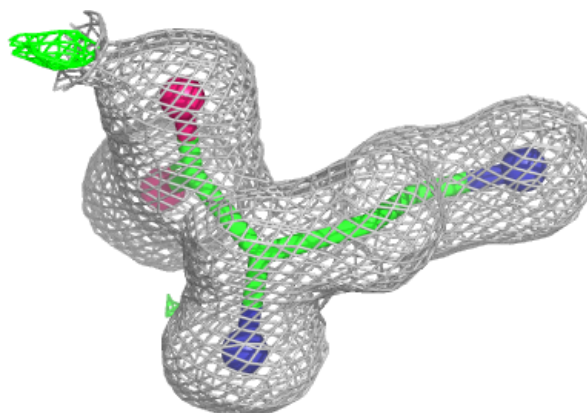
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	COA	A	201	48/48	0.98	0.06	12,18,37,59	5
4	MG	A	203	1/1	0.98	0.06	34,34,34,34	0
3	DAB	A	202	8/8	0.99	0.03	12,13,13,14	0
3	DAB	B	202	8/8	0.99	0.04	10,11,12,13	0
2	COA	B	201	48/48	0.99	0.05	10,14,29,54	9

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.

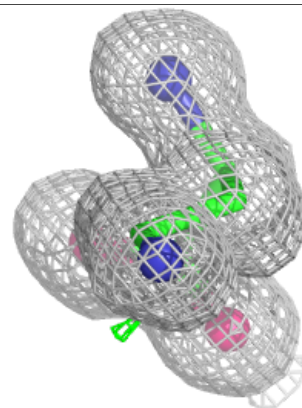
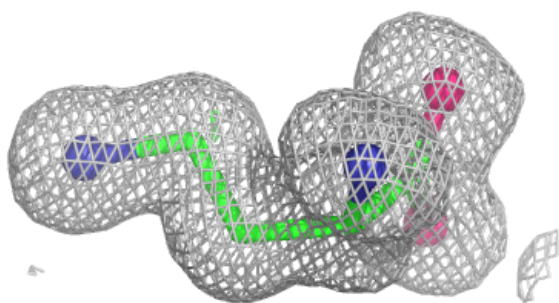
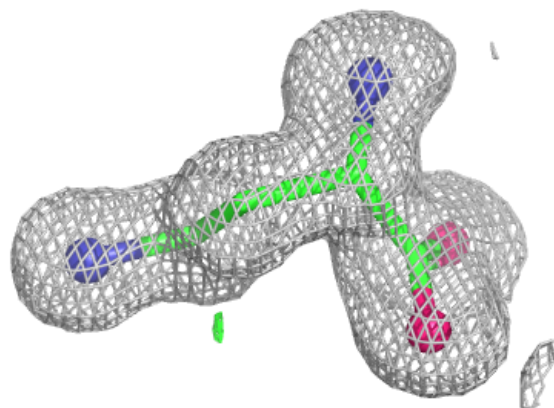


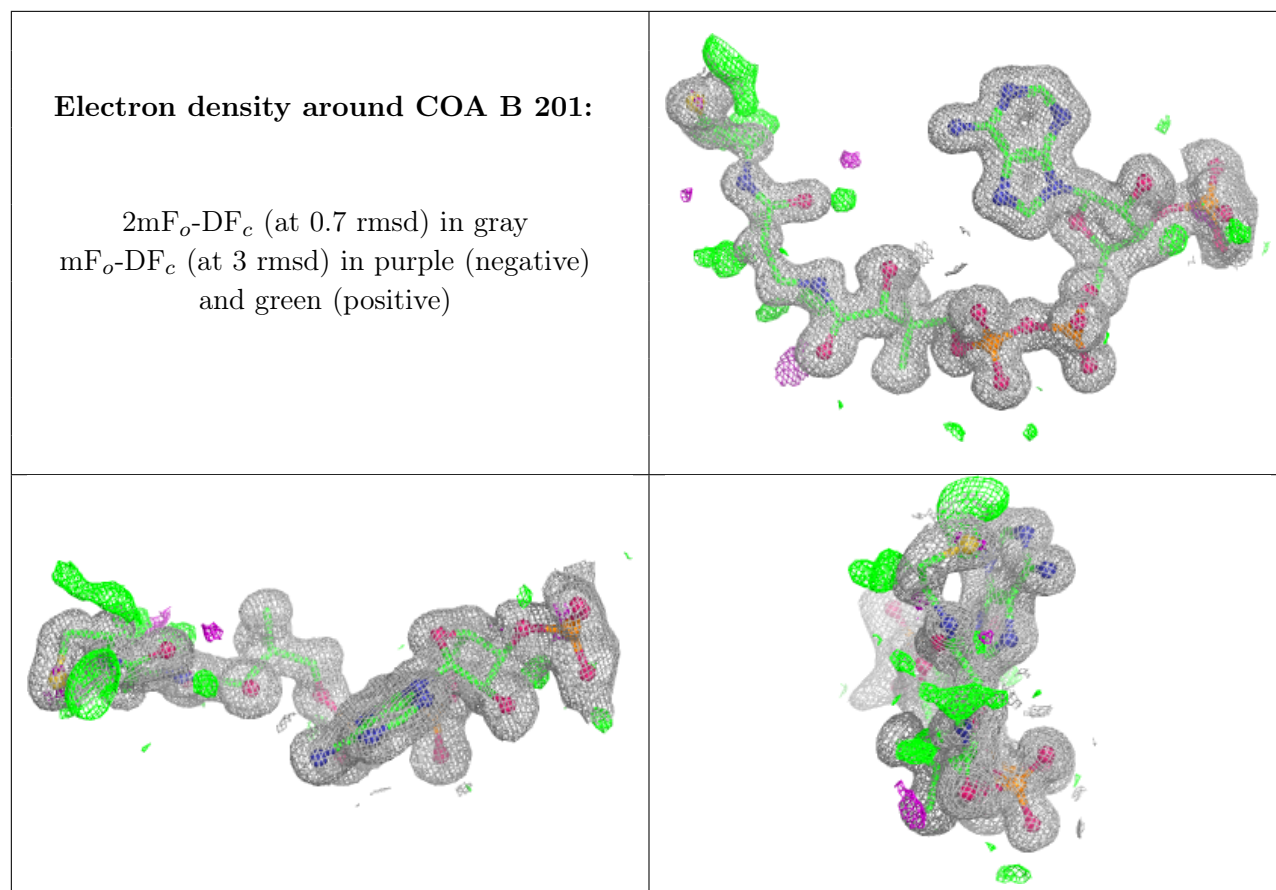
Electron density around DAB A 202:

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

**Electron density around DAB B 202:**

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





6.5 Other polymers ⓘ

There are no such residues in this entry.