Full wwPDB X-ray Structure Validation Report ⓘ

Oct 26, 2026 – 04:09 PM EDT

PDB ID	:	39RG / pdb_000039rg
Title	:	Crystal structure of a formate acetyltransferase (PFL) from Amygdalobacter indicium (CoA complex)
Authors	:	Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on	:	2026-10-22
Resolution	:	2.20 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:

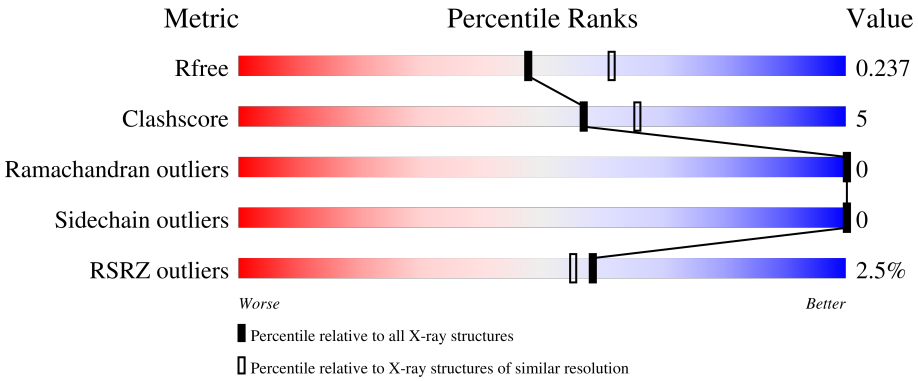
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

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 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 <=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	688	2% 81% 8% 11%
1	B	688	2% 77% 11% 12%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 10042 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 Formate acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	614	Total	C	N	O	S	0	0	0
			4844	3071	820	920	33			
1	B	607	Total	C	N	O	S	0	0	0
			4758	3015	816	895	32			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	initiating methionine	UNP A0ABY8C369
A	-20	ALA	-	expression tag	UNP A0ABY8C369
A	-19	HIS	-	expression tag	UNP A0ABY8C369
A	-18	HIS	-	expression tag	UNP A0ABY8C369
A	-17	HIS	-	expression tag	UNP A0ABY8C369
A	-16	HIS	-	expression tag	UNP A0ABY8C369
A	-15	HIS	-	expression tag	UNP A0ABY8C369
A	-14	HIS	-	expression tag	UNP A0ABY8C369
A	-13	MET	-	expression tag	UNP A0ABY8C369
A	-12	GLY	-	expression tag	UNP A0ABY8C369
A	-11	THR	-	expression tag	UNP A0ABY8C369
A	-10	LEU	-	expression tag	UNP A0ABY8C369
A	-9	GLU	-	expression tag	UNP A0ABY8C369
A	-8	ALA	-	expression tag	UNP A0ABY8C369
A	-7	GLN	-	expression tag	UNP A0ABY8C369
A	-6	THR	-	expression tag	UNP A0ABY8C369
A	-5	GLN	-	expression tag	UNP A0ABY8C369
A	-4	GLY	-	expression tag	UNP A0ABY8C369
A	-3	PRO	-	expression tag	UNP A0ABY8C369
A	-2	GLY	-	expression tag	UNP A0ABY8C369
A	-1	SER	-	expression tag	UNP A0ABY8C369
A	0	MET	-	expression tag	UNP A0ABY8C369
B	-21	MET	-	initiating methionine	UNP A0ABY8C369
B	-20	ALA	-	expression tag	UNP A0ABY8C369
B	-19	HIS	-	expression tag	UNP A0ABY8C369

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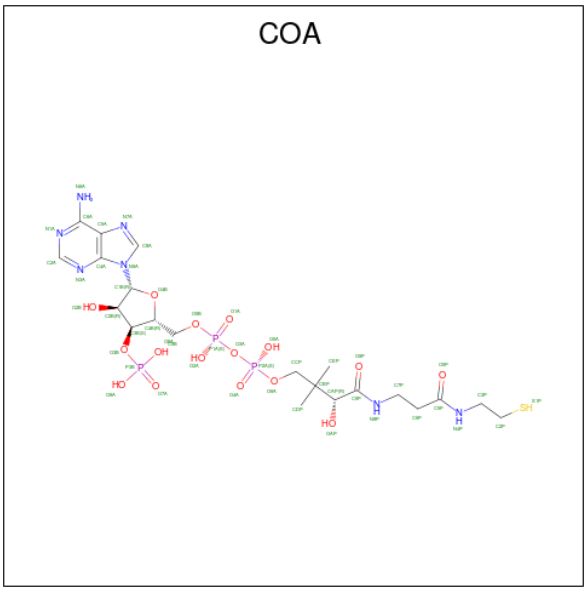
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	HIS	-	expression tag	UNP A0ABY8C369
B	-17	HIS	-	expression tag	UNP A0ABY8C369
B	-16	HIS	-	expression tag	UNP A0ABY8C369
B	-15	HIS	-	expression tag	UNP A0ABY8C369
B	-14	HIS	-	expression tag	UNP A0ABY8C369
B	-13	MET	-	expression tag	UNP A0ABY8C369
B	-12	GLY	-	expression tag	UNP A0ABY8C369
B	-11	THR	-	expression tag	UNP A0ABY8C369
B	-10	LEU	-	expression tag	UNP A0ABY8C369
B	-9	GLU	-	expression tag	UNP A0ABY8C369
B	-8	ALA	-	expression tag	UNP A0ABY8C369
B	-7	GLN	-	expression tag	UNP A0ABY8C369
B	-6	THR	-	expression tag	UNP A0ABY8C369
B	-5	GLN	-	expression tag	UNP A0ABY8C369
B	-4	GLY	-	expression tag	UNP A0ABY8C369
B	-3	PRO	-	expression tag	UNP A0ABY8C369
B	-2	GLY	-	expression tag	UNP A0ABY8C369
B	-1	SER	-	expression tag	UNP A0ABY8C369
B	0	MET	-	expression tag	UNP A0ABY8C369

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

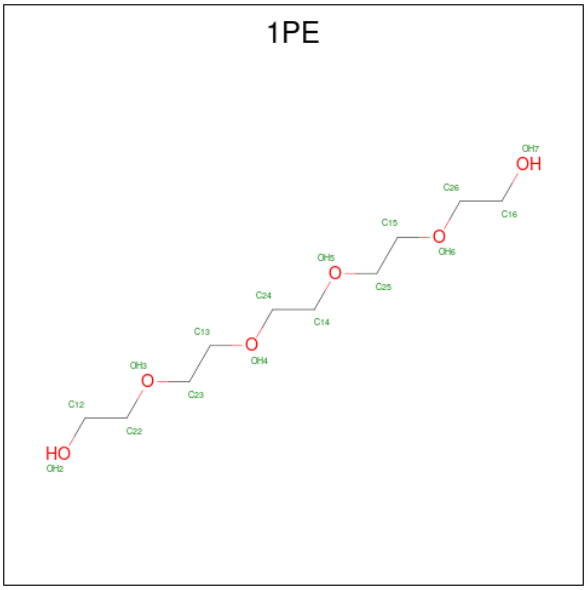
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Cl 2 2	0	0
2	B	3	Total Cl 3 3	0	0

- Molecule 3 is COENZYME A (CCD ID: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			16	10	6		
4	B	1	Total	C	O	0	0
			16	10	6		

- Molecule 5 is water.

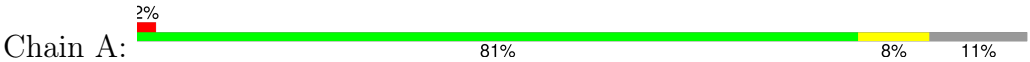
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	192	Total O 192 192	0	0
5	B	115	Total O 115 115	0	0

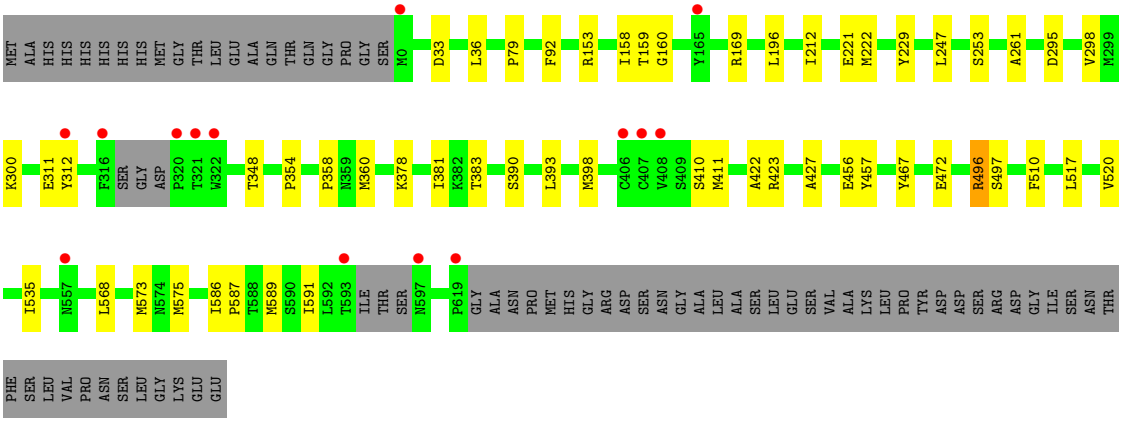
3 Residue-property plots (i)

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: Formate acetyltransferase

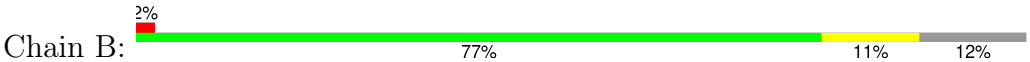
Chain A:

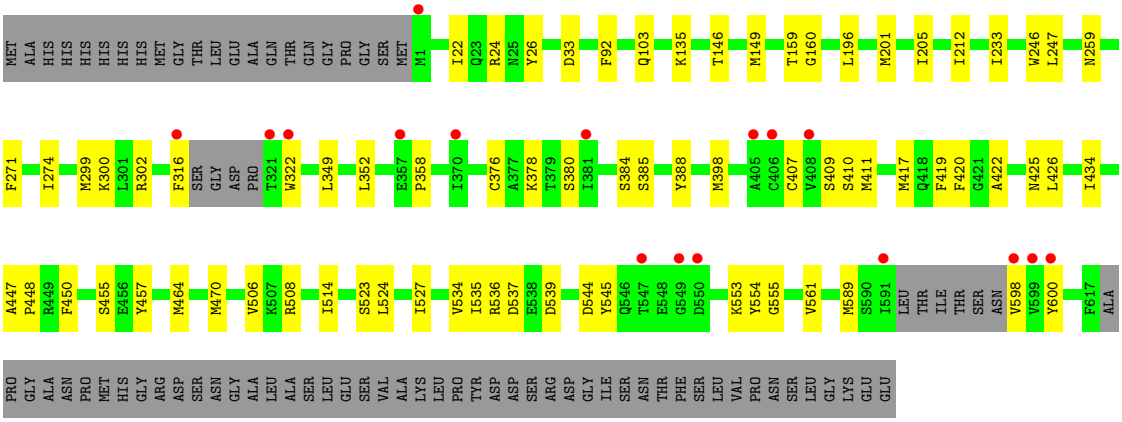




• Molecule 1: Formate acetyltransferase

Chain B:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	102.53Å 102.53Å 228.97Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.11 – 2.20 48.11 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.11-2.20) 100.0 (48.11-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.66 (at 2.20Å)	Xtriage
Refinement program	PHENIX 2.2.1_6174	Depositor
R, R_{free}	0.184 , 0.238 0.189 , 0.237	Depositor DCC
R_{free} test set	3644 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å^2)	51.2	Xtriage
Anisotropy	0.145	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 40.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	10042	wwPDB-VP
Average B, all atoms (Å^2)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 1PE, CL, COA

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	0.37	0/4947	0.50	0/6690
1	B	0.33	0/4859	0.47	0/6570
All	All	0.35	0/9806	0.49	0/13260

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	496	ARG	Sidechain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4844	0	4715	39	0
1	B	4758	0	4600	48	0
2	A	2	0	0	1	0
2	B	3	0	0	0	0
3	A	48	0	32	0	0
3	B	48	0	32	0	0
4	A	16	0	22	1	0
4	B	16	0	22	1	0
5	A	192	0	0	1	0
5	B	115	0	0	2	0
All	All	10042	0	9423	88	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 (88) 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:573:MET:HE3	1:A:589:MET:HE2	1.56	0.87
1:A:390:SER:OG	1:A:393:LEU:HD23	1.80	0.81
1:A:358:PRO:HG2	1:A:360:MET:HE2	1.70	0.73
1:B:411:MET:CE	5:B:825:HOH:O	2.45	0.64
1:B:523:SER:O	1:B:527:ILE:HG13	1.96	0.64
1:A:591:ILE:HG22	1:A:591:ILE:O	2.01	0.60
1:B:299:MET:HE2	1:B:352:LEU:HD21	1.83	0.60
1:B:24:ARG:NE	1:B:24:ARG:HA	2.17	0.60
1:A:573:MET:CE	1:A:589:MET:HE2	2.30	0.59
1:B:316:PHE:CD2	1:B:322:TRP:CZ3	2.90	0.58
1:A:573:MET:HA	1:A:573:MET:HE2	1.86	0.58
1:B:196:LEU:HD12	1:B:212:ILE:HD12	1.86	0.57
1:B:146:THR:OG1	1:B:149:MET:HG3	2.05	0.56
1:B:426:LEU:HD23	1:B:514:ILE:HG23	1.86	0.56
1:A:92:PHE:CE2	1:A:300:LYS:HE3	2.40	0.56
1:B:450:PHE:HD2	1:B:470:MET:HE1	1.71	0.56
1:B:316:PHE:HD2	1:B:322:TRP:CH2	2.25	0.55
1:A:573:MET:HE3	1:A:589:MET:HB2	1.89	0.55
1:B:316:PHE:CD2	1:B:322:TRP:HZ3	2.26	0.54
1:B:247:LEU:HD23	1:B:247:LEU:C	2.33	0.52
1:B:553:LYS:C	1:B:561:VAL:HG21	2.33	0.52
1:A:378:LYS:O	1:A:381:ILE:HG22	2.09	0.52
1:B:135:LYS:HB2	4:B:705:1PE:H242	1.92	0.52
1:B:514:ILE:HD11	1:B:589:MET:HE3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:196:LEU:CD1	1:B:212:ILE:HD12	2.40	0.51
1:A:517:LEU:CD2	1:A:591:ILE:HG21	2.41	0.51
1:A:33:ASP:HB2	1:A:36:LEU:HD12	1.93	0.50
1:A:169:ARG:HH12	1:A:261:ALA:HB3	1.76	0.50
1:B:316:PHE:CD2	1:B:322:TRP:CH2	3.00	0.50
1:B:302:ARG:HG2	1:B:358:PRO:HD3	1.94	0.50
1:B:398:MET:HE3	1:B:410:SER:OG	2.12	0.50
1:A:586:ILE:N	1:A:586:ILE:HD12	2.26	0.50
1:B:411:MET:HE1	5:B:825:HOH:O	2.10	0.49
1:B:26:TYR:HA	1:B:274:ILE:HG12	1.94	0.49
1:B:103:GLN:OE1	1:B:259:ASN:HB3	2.13	0.48
1:A:354:PRO:HG3	1:A:383:THR:HG22	1.95	0.48
1:B:92:PHE:CE2	1:B:300:LYS:HE3	2.48	0.48
1:B:417:MET:HG2	1:B:506:VAL:HG11	1.95	0.48
1:B:33:ASP:OD2	1:B:378:LYS:NZ	2.47	0.47
1:A:222:MET:HG2	1:A:253:SER:OG	2.14	0.47
1:A:456:GLU:HG2	1:A:457:TYR:HD1	1.80	0.47
1:B:425:ASN:HB2	1:B:600:TYR:CE1	2.50	0.47
1:B:554:TYR:HD1	1:B:555:GLY:N	2.12	0.47
1:A:311:GLU:HG2	1:A:312:TYR:N	2.30	0.46
1:B:22:ILE:HD11	1:B:271:PHE:HB2	1.97	0.46
1:A:517:LEU:HD23	1:A:591:ILE:HG21	1.97	0.46
1:B:535:ILE:HB	1:B:544:ASP:HB3	1.97	0.46
1:A:456:GLU:HG2	1:A:457:TYR:CD1	2.51	0.46
1:B:464:MET:CE	1:B:524:LEU:HD13	2.46	0.46
1:B:316:PHE:HD2	1:B:322:TRP:HH2	1.61	0.45
1:B:349:LEU:HD13	1:B:385:SER:HB2	1.99	0.45
1:A:247:LEU:C	1:A:247:LEU:HD23	2.41	0.45
1:B:419:PHE:CZ	1:B:508:ARG:HD3	2.52	0.45
1:A:472:GLU:HG2	1:A:575:MET:HE2	1.99	0.45
1:A:222:MET:HE2	1:A:497:SER:HB2	1.98	0.44
1:B:201:MET:HE1	1:B:205:ILE:HG22	1.99	0.44
1:B:233:ILE:HG22	1:B:246:TRP:CD2	2.52	0.44
1:B:464:MET:HE1	1:B:524:LEU:HD13	1.98	0.44
1:B:247:LEU:HD23	1:B:247:LEU:O	2.17	0.44
1:B:537:ASP:OD2	1:B:539:ASP:HB2	2.18	0.44
1:B:534:VAL:HG11	1:B:536:ARG:NH2	2.33	0.44
1:B:598:VAL:HG23	1:B:600:TYR:H	1.84	0.43
1:A:427:ALA:HA	1:A:520:VAL:HG22	2.00	0.43
1:A:422:ALA:C	1:A:423:ARG:HG2	2.44	0.43
1:B:420:PHE:CZ	1:B:422:ALA:HA	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:VAL:HG13	1:A:358:PRO:HG3	2.00	0.42
1:A:467:TYR:CE2	1:A:568:LEU:HD11	2.54	0.42
4:A:704:1PE:H151	5:A:912:HOH:O	2.19	0.42
1:A:398:MET:HE3	1:A:410:SER:C	2.44	0.42
1:B:376:CYS:HB3	1:B:388:TYR:OH	2.20	0.42
1:B:447:ALA:HB1	1:B:448:PRO:HD2	2.01	0.42
1:A:196:LEU:HD12	1:A:212:ILE:HD12	2.02	0.41
1:A:222:MET:CE	1:A:497:SER:HB2	2.50	0.41
1:A:398:MET:CE	1:A:410:SER:C	2.93	0.41
1:A:159:THR:HG22	1:A:160:GLY:H	1.84	0.41
1:A:153:ARG:HA	1:A:158:ILE:O	2.21	0.41
1:A:411:MET:HE2	1:A:411:MET:HB3	1.92	0.41
1:A:160:GLY:O	2:A:701:CL:CL	2.76	0.41
1:A:535:ILE:HD12	1:A:535:ILE:N	2.35	0.41
1:B:407:CYS:C	1:B:409:SER:N	2.78	0.41
1:A:221:GLU:OE1	1:A:496:ARG:NH2	2.50	0.41
1:B:380:SER:O	1:B:384:SER:HA	2.20	0.41
1:B:434:ILE:HD13	1:B:545:TYR:OH	2.20	0.41
1:B:455:SER:OG	1:B:457:TYR:O	2.35	0.40
1:A:510:PHE:HB3	1:A:587:PRO:HA	2.04	0.40
1:A:79:PRO:HA	1:A:229:TYR:CZ	2.56	0.40
1:A:295:ASP:OD2	1:A:348:THR:HG22	2.21	0.40
1:B:159:THR:CG2	1:B:160:GLY:N	2.84	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	608/688 (88%)	587 (96%)	21 (4%)	0	100	100
1	B	601/688 (87%)	577 (96%)	24 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1209/1376 (88%)	1164 (96%)	45 (4%)	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	508/580 (88%)	508 (100%)	0	100 100
1	B	490/580 (84%)	490 (100%)	0	100 100
All	All	998/1160 (86%)	998 (100%)	0	100 100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	191	GLN
1	A	313	ASN
1	A	418	GLN
1	A	469	GLN
1	B	154	HIS
1	B	241	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 9 ligands modelled in this entry, 5 are 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	$\#\|Z\| > 2$
4	1PE	B	705	-	15,15,15	0.32	0	14,14,14	0.51	0
3	COA	B	704	-	47,50,50	0.59	1 (2%)	69,75,75	0.54	0
3	COA	A	703	-	47,50,50	0.54	1 (2%)	69,75,75	0.53	0
4	1PE	A	704	-	15,15,15	0.31	0	14,14,14	0.38	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
4	1PE	B	705	-	-	5/13/13/13	-
3	COA	B	704	-	-	15/48/64/64	0/3/3/3
3	COA	A	703	-	-	10/48/64/64	0/3/3/3
4	1PE	A	704	-	-	7/13/13/13	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	704	COA	P3B-O3B	2.14	1.63	1.59
3	A	703	COA	P3B-O3B	2.07	1.63	1.59

There are no bond angle outliers.

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	703	COA	CCP-O6A-P2A-O3A
3	A	703	COA	CCP-O6A-P2A-O5A
3	A	703	COA	OAP-CAP-CBP-CCP
3	A	703	COA	C9P-CAP-CBP-CCP
3	A	703	COA	C9P-CAP-CBP-CDP
3	A	703	COA	OAP-CAP-CBP-CEP
3	A	703	COA	C9P-CAP-CBP-CEP
3	A	703	COA	C5P-C6P-C7P-N8P
3	B	704	COA	CCP-O6A-P2A-O3A
3	B	704	COA	CCP-O6A-P2A-O4A
3	B	704	COA	CCP-O6A-P2A-O5A
3	B	704	COA	OAP-CAP-CBP-CCP
3	B	704	COA	C9P-CAP-CBP-CCP
3	B	704	COA	C9P-CAP-CBP-CDP
3	B	704	COA	OAP-CAP-CBP-CEP
3	B	704	COA	C9P-CAP-CBP-CEP
3	B	704	COA	S1P-C2P-C3P-N4P
4	B	705	1PE	OH4-C13-C23-OH3
4	A	704	1PE	OH5-C14-C24-OH4
4	A	704	1PE	OH6-C15-C25-OH5
4	A	704	1PE	OH2-C12-C22-OH3
3	B	704	COA	O9P-C9P-CAP-OAP
4	A	704	1PE	C12-C22-OH3-C23
3	A	703	COA	OAP-CAP-CBP-CDP
3	B	704	COA	OAP-CAP-CBP-CDP
3	B	704	COA	C2P-C3P-N4P-C5P
3	B	704	COA	O5P-C5P-C6P-C7P
4	B	705	1PE	C24-C14-OH5-C25
4	A	704	1PE	C23-C13-OH4-C24
4	B	705	1PE	C15-C25-OH5-C14
4	A	704	1PE	C25-C15-OH6-C26
3	B	704	COA	N4P-C5P-C6P-C7P
3	A	703	COA	N8P-C9P-CAP-OAP
3	B	704	COA	N8P-C9P-CAP-OAP
4	B	705	1PE	OH6-C15-C25-OH5
4	B	705	1PE	OH2-C12-C22-OH3
4	A	704	1PE	OH4-C13-C23-OH3

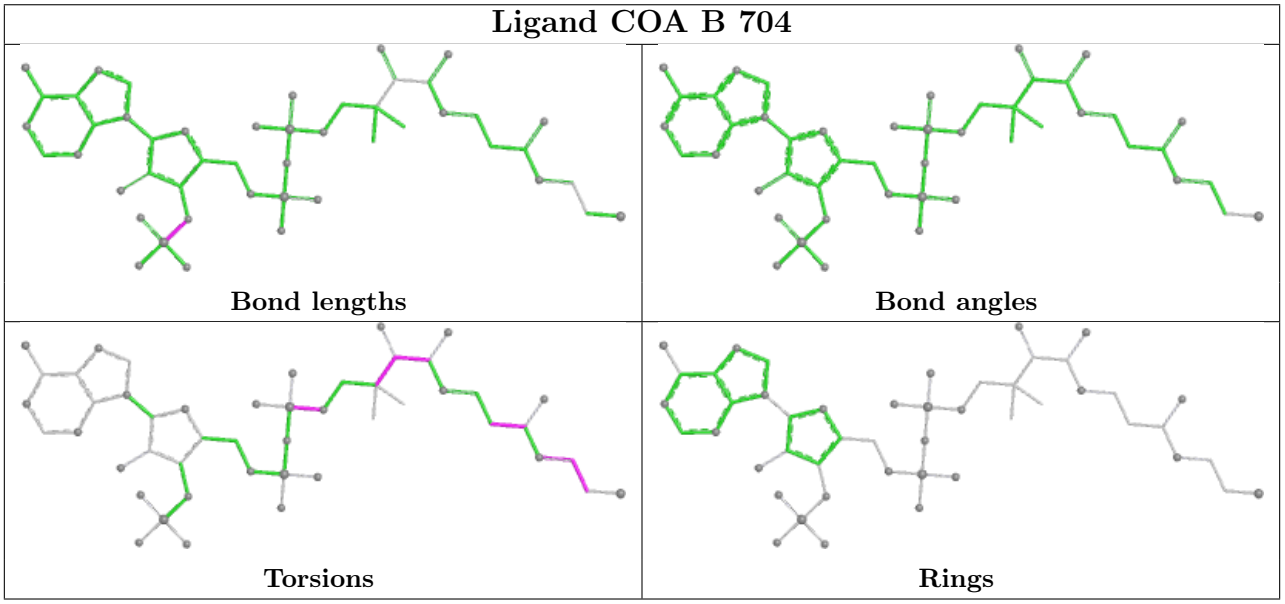
There are no ring outliers.

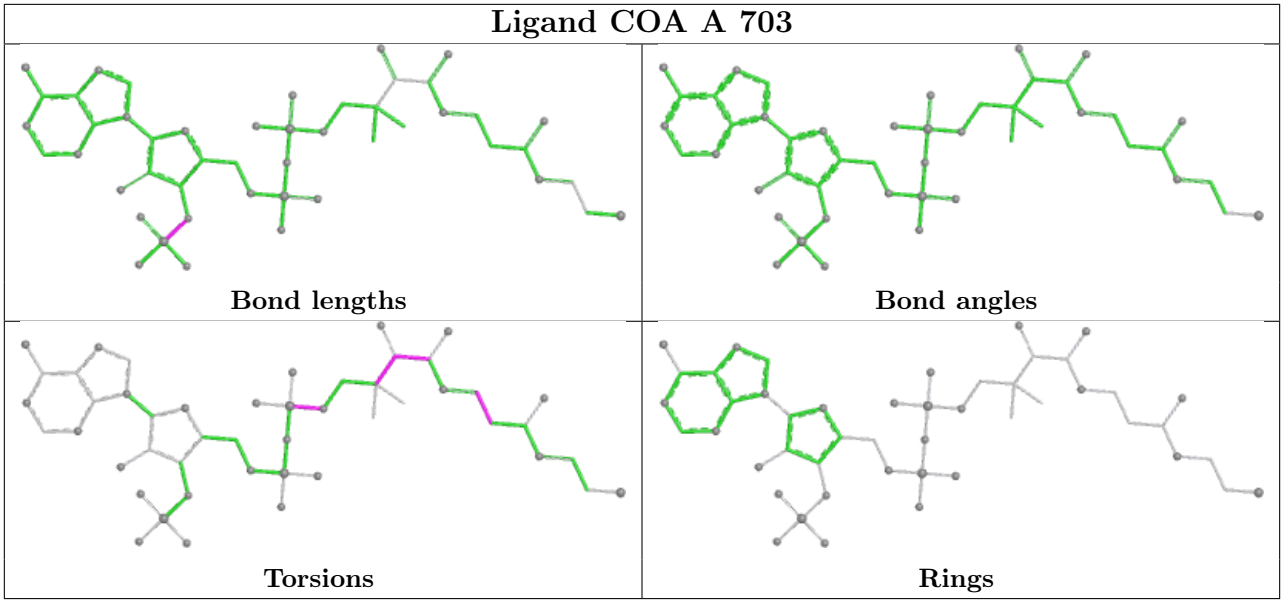
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	705	1PE	1	0
4	A	704	1PE	1	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 COA B 704





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

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, 95^{th} 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	614/688 (89%)	-0.06	14 (2%)	61	58	38, 56, 81, 124	0
1	B	607/688 (88%)	0.31	17 (2%)	55	52	37, 69, 99, 140	0
All	All	1221/1376 (88%)	0.13	31 (2%)	58	55	37, 61, 94, 140	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	321	THR	7.3
1	B	598	VAL	5.3
1	B	599	VAL	5.1
1	A	320	PRO	4.4
1	B	591	ILE	3.5
1	A	593	THR	3.5
1	A	316	PHE	3.4
1	B	406	CYS	3.2
1	A	321	THR	3.2
1	B	408	VAL	3.1
1	A	619	PRO	3.1
1	B	550	ASP	3.0
1	A	408	VAL	2.9
1	A	0	MET	2.7
1	B	405	ALA	2.7
1	A	406	CYS	2.5
1	B	322	TRP	2.5
1	B	357	GLU	2.5
1	A	407	CYS	2.5
1	A	597	ASN	2.3
1	A	312	TYR	2.3
1	B	600	TYR	2.3
1	A	322	TRP	2.3
1	B	1	MET	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	381	ILE	2.3
1	B	316	PHE	2.1
1	B	549	GLY	2.1
1	A	557	ASN	2.0
1	B	370	ILE	2.0
1	A	165	TYR	2.0
1	B	547	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)

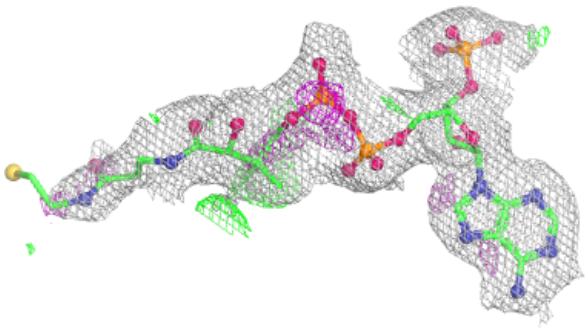
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

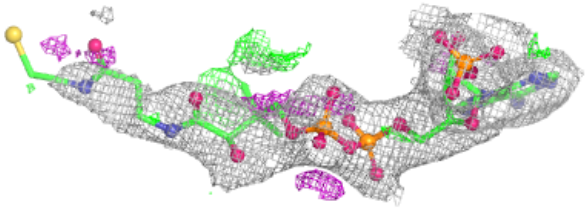
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	CL	A	701	1/1	0.77	0.25	74,74,74,74	0
2	CL	A	702	1/1	0.84	0.24	129,129,129,129	0
2	CL	B	702	1/1	0.85	0.16	77,77,77,77	0
2	CL	B	703	1/1	0.85	0.25	122,122,122,122	0
3	COA	B	704	48/48	0.85	0.12	54,80,98,114	0
3	COA	A	703	48/48	0.89	0.12	50,79,106,115	0
2	CL	B	701	1/1	0.92	0.10	76,76,76,76	0
4	1PE	A	704	16/16	0.92	0.10	48,59,69,73	0
4	1PE	B	705	16/16	0.92	0.11	46,58,67,80	0

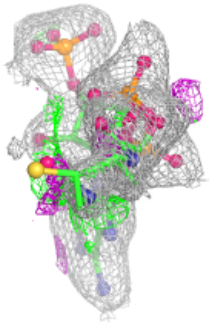
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around COA B 704:

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

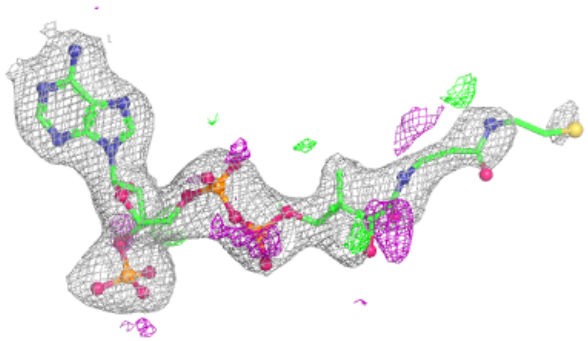


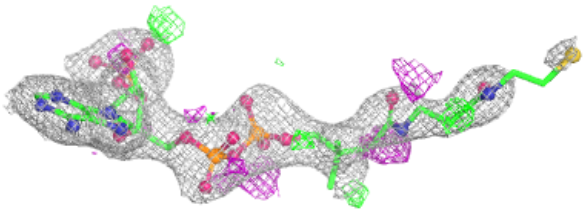


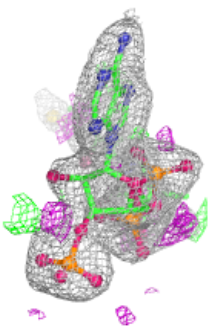


Electron density around COA A 703:

$2mF_o$-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.