



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:08 AM UTC

PDB ID : 4UBU / pdb_00004ubu
Title : Structure of a modified C93S variant of the 3-ketoacyl-CoA thiolase FadA5 from *M. tuberculosis* in complex with CoA
Authors : Schaefer, C.M.; Kisker, C.
Deposited on : 2014-08-13
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) : 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

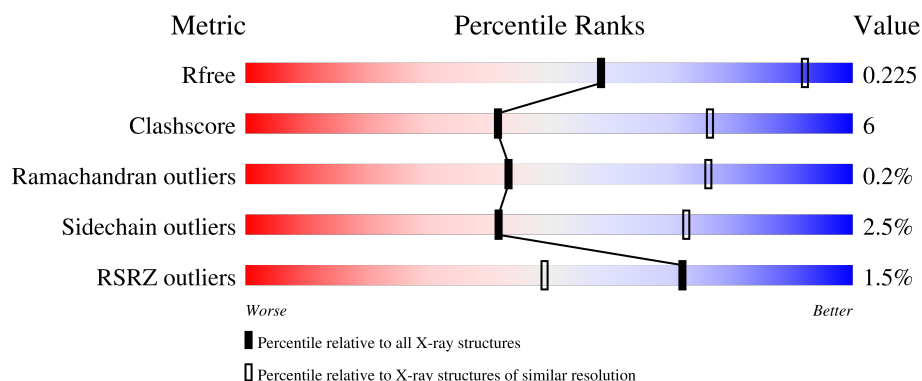
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (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 $\leq 5\%$. 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	399	0% 84% 13% ..
1	B	399	0% 83% 14% ..
1	C	399	2% 84% 13% ..
1	D	399	0% 86% 11% ..
1	E	399	2% 81% 13% . 5%

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Mol	Chain	Length	Quality of chain
1	F	399	2% 81% 15% ••
1	G	399	2% 85% 13% ••
1	H	399	% 85% 13% ••

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23649 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 Acetyl-CoA acetyltransferase FadA5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	0	1	0
			2911	1801	539	559	12			
1	B	391	Total	C	N	O	S	0	2	0
			2918	1806	541	559	12			
1	C	391	Total	C	N	O	S	0	0	0
			2903	1796	536	559	12			
1	D	391	Total	C	N	O	S	0	1	0
			2911	1801	539	559	12			
1	E	383	Total	C	N	O	S	0	1	2
			2827	1750	524	541	12			
1	F	386	Total	C	N	O	S	0	2	2
			2872	1777	530	554	11			
1	G	393	Total	C	N	O	S	0	1	0
			2921	1806	541	562	12			
1	H	394	Total	C	N	O	S	0	2	0
			2939	1820	543	564	12			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	HIS	-	expression tag	UNP I6XHI4
A	-6	HIS	-	expression tag	UNP I6XHI4
A	-5	HIS	-	expression tag	UNP I6XHI4
A	-4	HIS	-	expression tag	UNP I6XHI4
A	-3	HIS	-	expression tag	UNP I6XHI4
A	-2	HIS	-	expression tag	UNP I6XHI4
A	-1	GLY	-	expression tag	UNP I6XHI4
A	0	SER	-	expression tag	UNP I6XHI4
B	-7	HIS	-	expression tag	UNP I6XHI4
B	-6	HIS	-	expression tag	UNP I6XHI4
B	-5	HIS	-	expression tag	UNP I6XHI4
B	-4	HIS	-	expression tag	UNP I6XHI4
B	-3	HIS	-	expression tag	UNP I6XHI4

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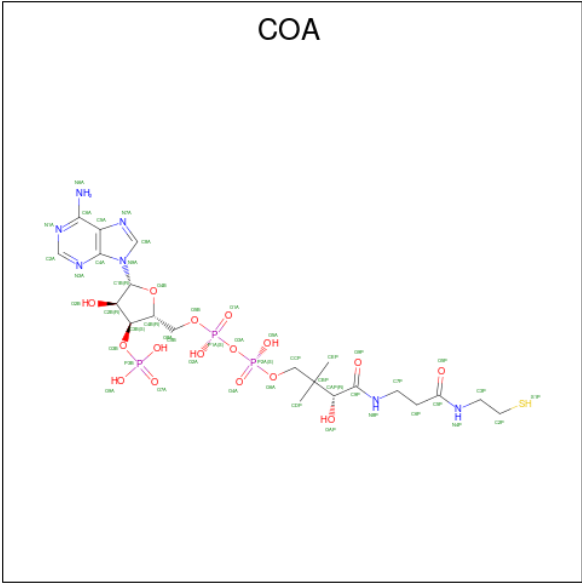
Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	HIS	-	expression tag	UNP I6XHI4
B	-1	GLY	-	expression tag	UNP I6XHI4
B	0	SER	-	expression tag	UNP I6XHI4
C	-7	HIS	-	expression tag	UNP I6XHI4
C	-6	HIS	-	expression tag	UNP I6XHI4
C	-5	HIS	-	expression tag	UNP I6XHI4
C	-4	HIS	-	expression tag	UNP I6XHI4
C	-3	HIS	-	expression tag	UNP I6XHI4
C	-2	HIS	-	expression tag	UNP I6XHI4
C	-1	GLY	-	expression tag	UNP I6XHI4
C	0	SER	-	expression tag	UNP I6XHI4
D	-7	HIS	-	expression tag	UNP I6XHI4
D	-6	HIS	-	expression tag	UNP I6XHI4
D	-5	HIS	-	expression tag	UNP I6XHI4
D	-4	HIS	-	expression tag	UNP I6XHI4
D	-3	HIS	-	expression tag	UNP I6XHI4
D	-2	HIS	-	expression tag	UNP I6XHI4
D	-1	GLY	-	expression tag	UNP I6XHI4
D	0	SER	-	expression tag	UNP I6XHI4
E	-7	HIS	-	expression tag	UNP I6XHI4
E	-6	HIS	-	expression tag	UNP I6XHI4
E	-5	HIS	-	expression tag	UNP I6XHI4
E	-4	HIS	-	expression tag	UNP I6XHI4
E	-3	HIS	-	expression tag	UNP I6XHI4
E	-2	HIS	-	expression tag	UNP I6XHI4
E	-1	GLY	-	expression tag	UNP I6XHI4
E	0	SER	-	expression tag	UNP I6XHI4
F	-7	HIS	-	expression tag	UNP I6XHI4
F	-6	HIS	-	expression tag	UNP I6XHI4
F	-5	HIS	-	expression tag	UNP I6XHI4
F	-4	HIS	-	expression tag	UNP I6XHI4
F	-3	HIS	-	expression tag	UNP I6XHI4
F	-2	HIS	-	expression tag	UNP I6XHI4
F	-1	GLY	-	expression tag	UNP I6XHI4
F	0	SER	-	expression tag	UNP I6XHI4
G	-7	HIS	-	expression tag	UNP I6XHI4
G	-6	HIS	-	expression tag	UNP I6XHI4
G	-5	HIS	-	expression tag	UNP I6XHI4
G	-4	HIS	-	expression tag	UNP I6XHI4
G	-3	HIS	-	expression tag	UNP I6XHI4
G	-2	HIS	-	expression tag	UNP I6XHI4
G	-1	GLY	-	expression tag	UNP I6XHI4

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Chain	Residue	Modelled	Actual	Comment	Reference
G	0	SER	-	expression tag	UNP I6XHI4
H	-7	HIS	-	expression tag	UNP I6XHI4
H	-6	HIS	-	expression tag	UNP I6XHI4
H	-5	HIS	-	expression tag	UNP I6XHI4
H	-4	HIS	-	expression tag	UNP I6XHI4
H	-3	HIS	-	expression tag	UNP I6XHI4
H	-2	HIS	-	expression tag	UNP I6XHI4
H	-1	GLY	-	expression tag	UNP I6XHI4
H	0	SER	-	expression tag	UNP I6XHI4

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



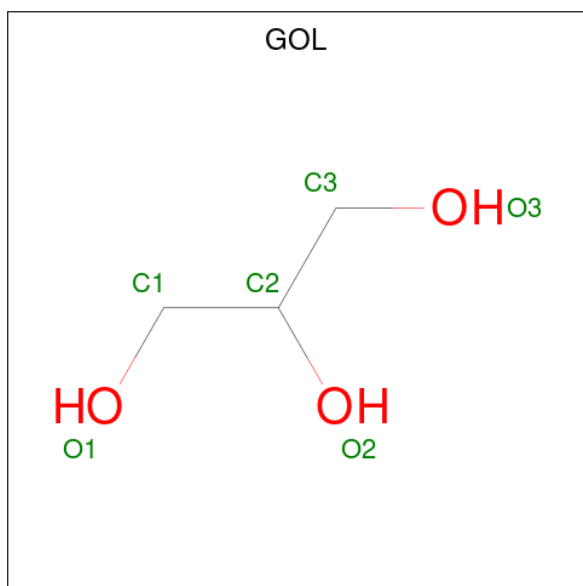
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	B	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	C	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	D	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	E	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	F	1	Total 48	C 21	N 7	O 16	P 3	S 1	0	0
2	G	1	Total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

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

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	O	0	0
			9	9		
4	B	6	Total	O	0	0
			6	6		
4	C	5	Total	O	0	0
			5	5		
4	E	3	Total	O	0	0
			3	3		

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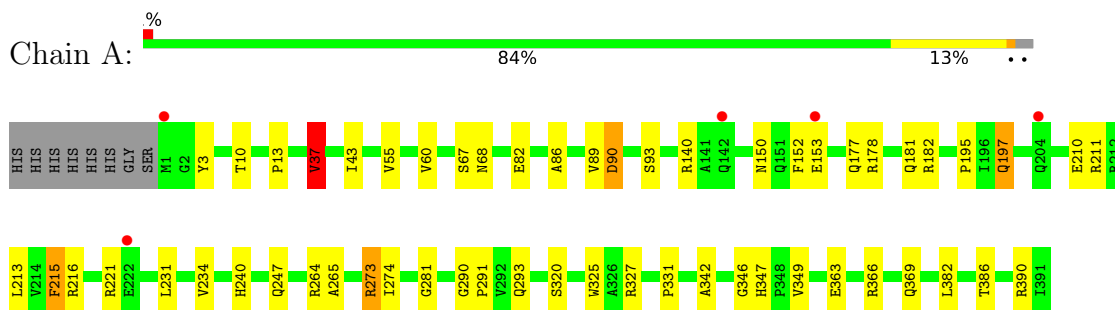
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	13	Total 13	O 13	0	0
4	G	1	Total 1	O 1	0	0
4	H	2	Total 2	O 2	0	0

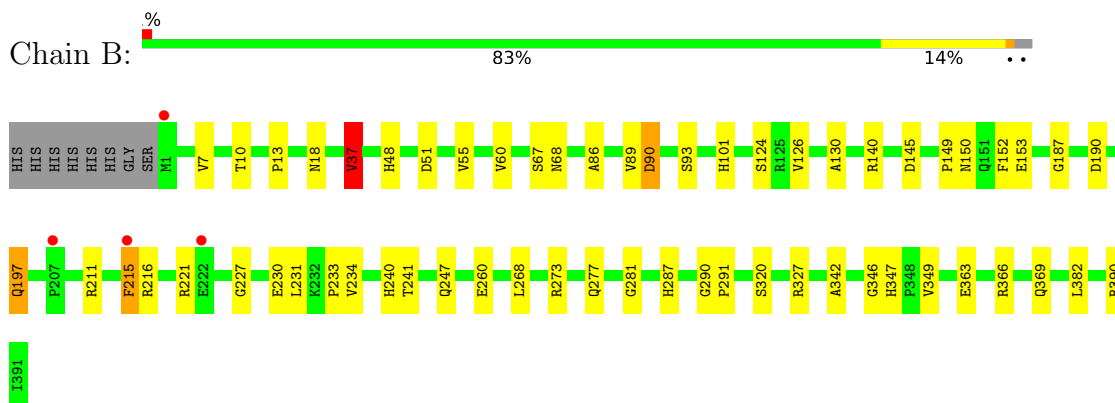
3 Residue-property plots [i](#)

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.

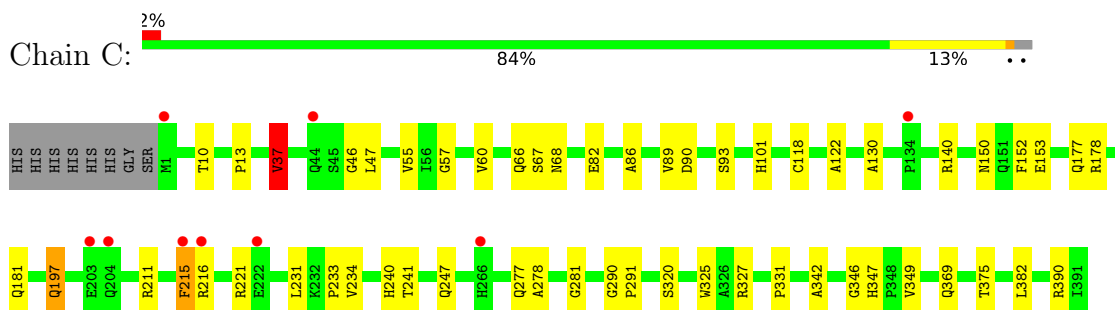
- Molecule 1: Acetyl-CoA acetyltransferase FadA5



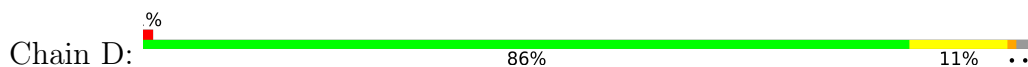
- Molecule 1: Acetyl-CoA acetyltransferase FadA5

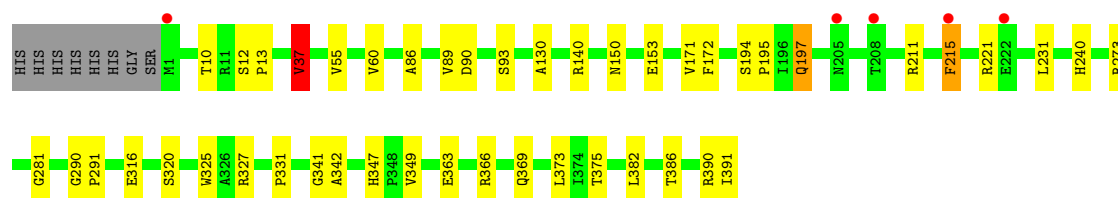


- Molecule 1: Acetyl-CoA acetyltransferase FadA5

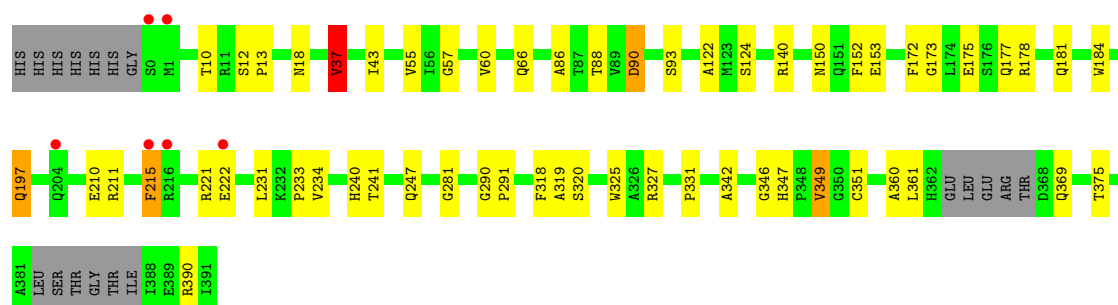
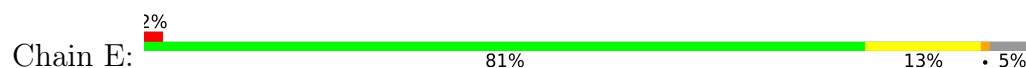


- Molecule 1: Acetyl-CoA acetyltransferase FadA5

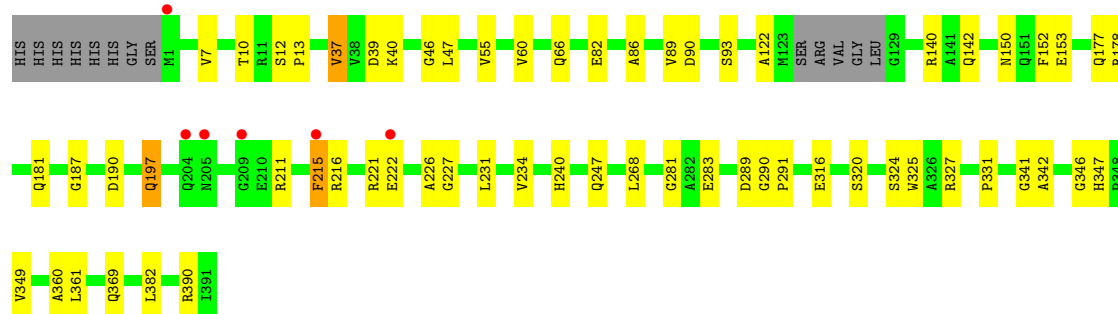
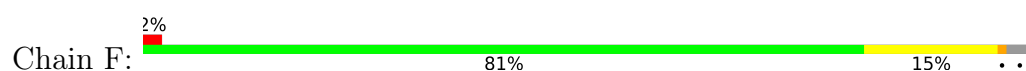




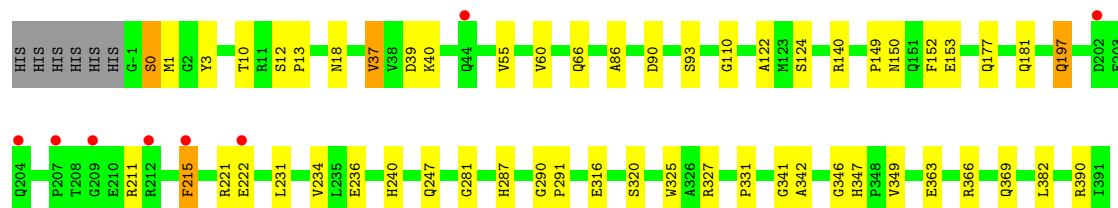
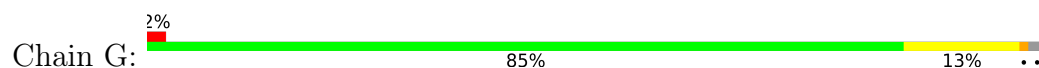
• Molecule 1: Acetyl-CoA acetyltransferase FadA5



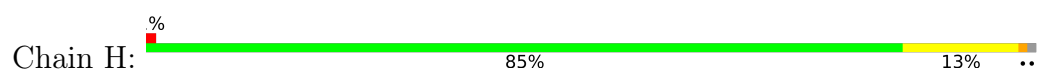
• Molecule 1: Acetyl-CoA acetyltransferase FadA5

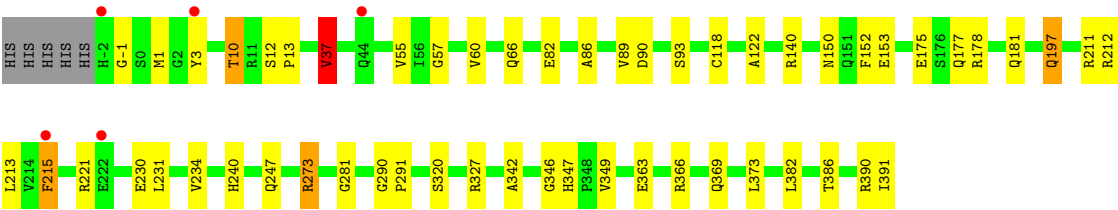


• Molecule 1: Acetyl-CoA acetyltransferase FadA5



• Molecule 1: Acetyl-CoA acetyltransferase FadA5





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	127.61Å 105.16Å 147.66Å 90.00° 107.33° 90.00°	Depositor
Resolution (Å)	58.82 – 3.00 58.82 – 3.00	Depositor EDS
% Data completeness (in resolution range)	94.7 (58.82-3.00) 94.7 (58.82-3.00)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 3.01Å)	Xtrriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.247 , 0.287 (Not available) , 0.225	Depositor DCC
R_{free} test set	3551 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	41.4	Xtrriage
Anisotropy	0.546	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 64.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	23649	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.91 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.9909e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: OAS, COA, GOL

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	0.92	0/2948	0.93	2/3996 (0.1%)
1	B	0.91	1/2959 (0.0%)	0.94	1/4011 (0.0%)
1	C	0.92	1/2937 (0.0%)	0.95	1/3982 (0.0%)
1	D	0.91	0/2948	0.94	4/3996 (0.1%)
1	E	0.91	0/2870	0.96	3/3890 (0.1%)
1	F	0.87	0/2921	0.93	2/3960 (0.1%)
1	G	0.89	0/2958	0.93	2/4009 (0.0%)
1	H	0.82	0/2982	0.92	3/4044 (0.1%)
All	All	0.89	2/23523 (0.0%)	0.94	18/31888 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	E	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	126	VAL	C-O	-5.67	1.17	1.24
1	C	278	ALA	CA-C	-5.35	1.45	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	12	SER	CA-C-N	-6.45	113.64	120.03
1	E	12	SER	C-N-CA	-6.45	113.64	120.03
1	D	12	SER	CA-C-N	-6.10	113.39	119.92
1	D	12	SER	C-N-CA	-6.10	113.39	119.92
1	E	37	VAL	CB-CA-C	-6.06	104.21	111.97
1	B	37	VAL	CB-CA-C	-5.82	104.52	111.97
1	C	37	VAL	CB-CA-C	-5.55	104.86	111.97
1	H	37	VAL	CB-CA-C	-5.45	105.00	111.97
1	A	37	VAL	N-CA-CB	5.40	116.87	110.55
1	G	12	SER	CA-C-N	-5.39	114.69	120.03
1	G	12	SER	C-N-CA	-5.39	114.69	120.03
1	H	12	SER	CA-C-N	-5.35	114.19	119.92
1	H	12	SER	C-N-CA	-5.35	114.19	119.92
1	D	37	VAL	CB-CA-C	-5.34	105.14	111.97
1	A	37	VAL	CB-CA-C	-5.31	105.17	111.97
1	D	37	VAL	N-CA-CB	5.29	116.74	110.55
1	F	12	SER	CA-C-N	-5.15	114.93	120.03
1	F	12	SER	C-N-CA	-5.15	114.93	120.03

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	90	ASP	Peptide
1	B	90	ASP	Peptide
1	E	90	ASP	Peptide

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	2911	0	2910	56	0
1	B	2918	0	2917	42	0
1	C	2903	0	2897	32	0
1	D	2911	0	2910	28	0
1	E	2827	0	2828	55	0
1	F	2872	0	2875	49	0
1	G	2921	0	2918	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	2939	0	2928	32	0
2	A	48	0	32	3	0
2	B	48	0	32	2	0
2	C	48	0	32	3	0
2	D	48	0	32	2	0
2	E	48	0	32	4	0
2	F	48	0	32	1	0
2	G	48	0	32	3	0
2	H	48	0	32	3	0
3	E	6	0	8	1	0
3	G	12	0	16	2	0
3	H	6	0	8	1	0
4	A	9	0	0	0	0
4	B	6	0	0	0	0
4	C	5	0	0	1	0
4	E	3	0	0	0	0
4	F	13	0	0	3	0
4	G	1	0	0	0	0
4	H	2	0	0	0	0
All	All	23649	0	23471	302	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 (302) 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:190:ASP:HB3	1:F:215:PHE:CE2	1.86	1.11
1:A:195:PRO:HG2	1:E:215:PHE:HZ	1.05	1.10
1:B:190:ASP:HB3	1:F:215:PHE:CZ	1.87	1.09
1:E:178[B]:ARG:HG2	1:E:178[B]:ARG:HH11	0.95	1.08
1:A:195:PRO:CG	1:E:215:PHE:HZ	1.69	1.06
1:A:265:ALA:N	1:G:1:MET:HE1	1.73	1.04
1:B:215:PHE:CE2	1:F:190:ASP:HB3	1.94	1.03
1:A:195:PRO:HG2	1:E:215:PHE:CZ	1.94	1.01
1:B:215:PHE:CZ	1:F:190:ASP:HB3	1.96	1.00
1:A:265:ALA:CA	1:G:1:MET:HE1	1.93	0.99
1:F:178[B]:ARG:HH11	1:F:178[B]:ARG:CG	1.79	0.95
1:E:178[B]:ARG:HG2	1:E:178[B]:ARG:NH1	1.72	0.88
1:F:178[B]:ARG:HH11	1:F:178[B]:ARG:HG2	1.43	0.83
1:A:195:PRO:CG	1:E:215:PHE:CZ	2.59	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:178[B]:ARG:HH11	1:E:178[B]:ARG:CG	1.86	0.82
1:B:190:ASP:CB	1:F:215:PHE:CZ	2.66	0.78
1:F:142:GLN:HG3	4:F:509:HOH:O	1.82	0.77
1:B:187:GLY:HA2	1:F:216:ARG:NH2	1.99	0.77
1:B:216:ARG:NH2	1:F:187:GLY:HA2	2.01	0.76
1:A:215:PHE:CE2	1:E:215:PHE:CD2	2.75	0.74
1:A:265:ALA:HA	1:G:1:MET:HE1	1.72	0.71
1:B:215:PHE:HD1	1:B:215:PHE:N	1.87	0.71
1:B:215:PHE:N	1:B:215:PHE:CD1	2.59	0.71
1:F:215:PHE:N	1:F:215:PHE:HD1	1.90	0.70
1:G:215:PHE:HD1	1:G:215:PHE:N	1.89	0.69
1:A:215:PHE:N	1:A:215:PHE:HD1	1.90	0.69
1:D:215:PHE:CD1	1:D:215:PHE:N	2.60	0.69
1:D:215:PHE:N	1:D:215:PHE:HD1	1.89	0.69
1:B:215:PHE:CZ	1:F:190:ASP:CB	2.75	0.68
1:A:215:PHE:N	1:A:215:PHE:CD1	2.61	0.68
1:H:215:PHE:N	1:H:215:PHE:HD1	1.91	0.68
1:F:215:PHE:N	1:F:215:PHE:CD1	2.61	0.68
1:C:240:HIS:HE1	1:C:320:SER:H	1.43	0.67
1:A:216:ARG:NH2	1:E:184:TRP:O	2.29	0.66
1:E:215:PHE:N	1:E:215:PHE:HD1	1.94	0.66
1:B:240:HIS:HE1	1:B:320:SER:H	1.44	0.66
1:H:215:PHE:N	1:H:215:PHE:CD1	2.63	0.66
1:C:215:PHE:CD1	1:C:215:PHE:N	2.64	0.66
1:C:215:PHE:N	1:C:215:PHE:HD1	1.93	0.65
1:A:240:HIS:HE1	1:A:320:SER:H	1.43	0.65
1:G:215:PHE:N	1:G:215:PHE:CD1	2.60	0.65
1:E:215:PHE:N	1:E:215:PHE:CD1	2.64	0.65
1:A:265:ALA:N	1:G:1:MET:CE	2.57	0.64
1:H:240:HIS:HE1	1:H:320:SER:H	1.45	0.64
1:A:215:PHE:CD2	1:E:215:PHE:CE2	2.86	0.64
1:D:90:ASP:OD1	1:D:90:ASP:C	2.41	0.63
1:F:90:ASP:OD1	1:F:90:ASP:C	2.40	0.63
1:E:240:HIS:HE1	1:E:320:SER:H	1.44	0.63
1:H:273:ARG:HD3	3:H:402:GOL:H2	1.80	0.63
1:F:178[B]:ARG:HG2	1:F:178[B]:ARG:NH1	2.11	0.62
1:F:215:PHE:HD1	1:F:215:PHE:H	1.47	0.62
1:F:240:HIS:HE1	1:F:320:SER:H	1.44	0.62
1:D:240:HIS:HE1	1:D:320:SER:H	1.47	0.62
1:G:240:HIS:HE1	1:G:320:SER:H	1.46	0.62
1:A:215:PHE:HD1	1:A:215:PHE:H	1.46	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:283:GLU:HG3	4:F:505:HOH:O	1.98	0.61
1:A:150:ASN:OD1	1:A:150:ASN:C	2.43	0.61
1:B:215:PHE:HD1	1:B:215:PHE:H	1.47	0.60
1:G:215:PHE:HD1	1:G:215:PHE:H	1.47	0.60
1:E:215:PHE:HD1	1:E:215:PHE:H	1.48	0.60
1:D:215:PHE:HD1	1:D:215:PHE:H	1.48	0.60
1:F:281:GLY:HA2	1:F:382:LEU:HD23	1.82	0.60
1:A:213:LEU:HD23	1:E:215:PHE:CD2	2.36	0.60
1:H:215:PHE:HD1	1:H:215:PHE:H	1.48	0.59
1:C:215:PHE:HD1	1:C:215:PHE:H	1.49	0.59
1:G:90:ASP:OD1	1:G:90:ASP:C	2.45	0.59
1:C:281:GLY:HA2	1:C:382:LEU:HD23	1.85	0.59
1:H:150:ASN:OD1	1:H:150:ASN:C	2.46	0.59
1:H:175:GLU:OE1	1:H:178:ARG:HD3	2.03	0.59
1:D:281:GLY:HA2	1:D:382:LEU:HD23	1.86	0.58
1:C:90:ASP:C	1:C:90:ASP:OD1	2.44	0.58
1:D:55:VAL:O	1:D:86:ALA:HA	2.04	0.58
1:B:281:GLY:HA2	1:B:382:LEU:HD23	1.85	0.58
1:A:43:ILE:HG13	1:E:210:GLU:OE2	2.04	0.57
1:B:89:VAL:HG12	1:C:89:VAL:HG12	1.86	0.57
1:H:363:GLU:OE2	1:H:366:ARG:NH2	2.38	0.57
1:F:178[B]:ARG:HH11	1:F:178[B]:ARG:HG3	1.66	0.56
1:G:18:ASN:N	1:G:124:SER:OG	2.36	0.56
1:F:55:VAL:O	1:F:86:ALA:HA	2.06	0.56
1:G:55:VAL:O	1:G:86:ALA:HA	2.05	0.55
1:A:281:GLY:HA2	1:A:382:LEU:HD23	1.88	0.55
1:G:342:ALA:HB1	1:G:347:HIS:HB2	1.88	0.55
1:G:281:GLY:HA2	1:G:382:LEU:HD23	1.87	0.55
1:A:89:VAL:HG12	1:D:89:VAL:HG12	1.88	0.54
1:F:82:GLU:HG3	4:F:506:HOH:O	2.07	0.54
1:C:342:ALA:HB1	1:C:347:HIS:HB2	1.88	0.54
1:A:342:ALA:HB1	1:A:347:HIS:HB2	1.90	0.54
1:D:342:ALA:HB1	1:D:347:HIS:HB2	1.90	0.54
1:A:197:GLN:HG2	1:A:211:ARG:HG2	1.90	0.53
1:A:3:TYR:HD2	1:G:3:TYR:CE2	2.27	0.53
1:E:342:ALA:HB1	1:E:347:HIS:HB2	1.90	0.53
1:F:342:ALA:HB1	1:F:347:HIS:HB2	1.90	0.53
1:F:89:VAL:HG12	1:H:89:VAL:HG12	1.90	0.53
1:D:231:LEU:HD21	2:D:400:COA:H1B	1.91	0.52
1:H:281:GLY:HA2	1:H:382:LEU:HD23	1.91	0.52
1:F:152:PHE:CD1	1:F:234:VAL:HG11	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:TYR:HD2	1:G:3:TYR:CZ	2.27	0.52
1:C:60:VAL:HG21	1:C:349:VAL:CG1	2.39	0.52
1:E:150:ASN:OD1	1:E:150:ASN:C	2.52	0.52
1:D:60:VAL:HG21	1:D:349:VAL:CG1	2.39	0.52
1:F:281:GLY:HA3	1:H:82:GLU:O	2.09	0.52
1:H:231:LEU:HD21	2:H:401:COA:H1B	1.91	0.52
1:B:231:LEU:HD21	2:B:400:COA:H1B	1.90	0.52
1:A:67:SER:OG	1:A:68:ASN:ND2	2.42	0.52
1:H:55:VAL:O	1:H:86:ALA:HA	2.10	0.52
1:F:60:VAL:HG21	1:F:349:VAL:CG1	2.40	0.51
1:D:13:PRO:HG3	1:D:215:PHE:HA	1.93	0.51
1:D:363:GLU:OE2	1:D:366:ARG:NH2	2.43	0.51
1:E:60:VAL:HG21	1:E:349:VAL:CG1	2.40	0.51
1:H:197:GLN:HG2	1:H:211:ARG:HG2	1.91	0.51
1:H:373:LEU:HD12	1:H:386:THR:O	2.10	0.51
1:E:231:LEU:HD21	2:E:401:COA:H1B	1.92	0.51
1:H:342:ALA:HB1	1:H:347:HIS:HB2	1.91	0.51
1:B:197:GLN:HG2	1:B:211:ARG:HG2	1.92	0.51
1:A:210:GLU:HG3	1:E:43:ILE:CG1	2.41	0.51
1:F:197:GLN:HG2	1:F:211:ARG:HG2	1.93	0.51
1:H:212:ARG:HG2	1:H:213:LEU:N	2.25	0.51
1:D:150:ASN:C	1:D:150:ASN:OD1	2.53	0.51
1:E:197:GLN:HG2	1:E:211:ARG:HG2	1.93	0.51
1:D:197:GLN:HG2	1:D:211:ARG:HG2	1.93	0.50
1:A:210:GLU:CG	1:E:43:ILE:HG12	2.41	0.50
1:G:197:GLN:HG2	1:G:211:ARG:HG2	1.94	0.50
1:A:60:VAL:HG21	1:A:349:VAL:CG1	2.41	0.49
1:C:67:SER:OG	1:C:68:ASN:ND2	2.44	0.49
1:G:60:VAL:HG21	1:G:349:VAL:CG1	2.42	0.49
1:B:90:ASP:OD1	1:B:90:ASP:C	2.55	0.49
1:C:197:GLN:HG2	1:C:211:ARG:HG2	1.94	0.49
1:C:231:LEU:HD21	2:C:400:COA:H1B	1.93	0.49
1:G:363:GLU:OE2	1:G:366:ARG:NH2	2.44	0.49
1:C:150:ASN:OD1	1:C:150:ASN:C	2.54	0.49
1:G:150:ASN:C	1:G:150:ASN:OD1	2.55	0.49
1:A:215:PHE:CE2	1:E:215:PHE:CE2	3.01	0.49
1:H:60:VAL:HG21	1:H:349:VAL:CG1	2.43	0.49
1:A:247:GLN:HB2	1:A:346:GLY:HA2	1.95	0.49
1:G:231:LEU:HD21	2:G:401:COA:H1B	1.95	0.49
1:H:273:ARG:HG2	1:H:391:ILE:HD11	1.93	0.48
1:B:55:VAL:O	1:B:86:ALA:HA	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:ASN:OD1	1:B:150:ASN:C	2.56	0.48
1:E:369:GLN:O	1:E:390:ARG:HD3	2.13	0.48
1:A:90:ASP:C	1:A:90:ASP:OD1	2.56	0.48
1:B:13:PRO:HG3	1:B:215:PHE:HA	1.96	0.48
1:B:281:GLY:HA3	1:C:82:GLU:O	2.13	0.48
1:G:13:PRO:HG3	1:G:215:PHE:HA	1.95	0.48
1:B:342:ALA:HB1	1:B:347:HIS:HB2	1.96	0.48
1:E:90:ASP:C	1:E:90:ASP:OD1	2.57	0.47
1:C:13:PRO:HG3	1:C:215:PHE:HA	1.97	0.47
1:F:150:ASN:OD1	1:F:150:ASN:C	2.56	0.47
1:E:178[B]:ARG:NH1	1:E:178[B]:ARG:CG	2.54	0.47
1:B:60:VAL:HG21	1:B:349:VAL:CG1	2.44	0.47
1:C:247:GLN:HB2	1:C:346:GLY:HA2	1.96	0.47
1:F:152:PHE:CE1	1:F:234:VAL:HG11	2.50	0.47
1:F:177:GLN:OE1	1:F:181:GLN:NE2	2.44	0.47
1:F:231:LEU:HD21	2:F:400:COA:H1B	1.96	0.47
1:H:10:THR:HG22	1:H:37:VAL:HG13	1.97	0.47
1:E:13:PRO:HG3	1:E:215:PHE:HA	1.96	0.47
1:H:13:PRO:HG3	1:H:215:PHE:HA	1.97	0.47
1:H:247:GLN:HB2	1:H:346:GLY:HA2	1.97	0.47
1:C:101:HIS:HB3	1:C:277:GLN:OE1	2.15	0.46
1:E:152:PHE:CD1	1:E:234:VAL:HG11	2.51	0.46
1:F:10:THR:HG22	1:F:37:VAL:HG13	1.97	0.46
1:D:194:SER:O	1:D:195:PRO:C	2.58	0.46
1:G:10:THR:HG22	1:G:37:VAL:HG13	1.97	0.46
1:E:175:GLU:OE1	1:E:178[A]:ARG:HD3	2.15	0.46
1:E:66:GLN:HE22	1:E:122:ALA:H	1.64	0.46
1:A:13:PRO:HG3	1:A:215:PHE:HA	1.97	0.46
1:E:10:THR:HG22	1:E:37:VAL:HG13	1.98	0.46
1:F:13:PRO:HG3	1:F:215:PHE:HA	1.97	0.46
1:A:325:TRP:CZ3	1:A:331:PRO:HG3	2.51	0.46
1:B:10:THR:HG22	1:B:37:VAL:HG13	1.97	0.46
1:B:18:ASN:N	1:B:124:SER:OG	2.39	0.46
1:G:152:PHE:CD1	1:G:234:VAL:HG11	2.51	0.46
1:G:325:TRP:CZ3	1:G:331:PRO:HG3	2.51	0.46
1:D:369:GLN:O	1:D:390:ARG:HD3	2.16	0.46
1:B:67:SER:OG	1:B:68:ASN:ND2	2.48	0.45
1:H:230:GLU:OE1	1:H:230:GLU:HA	2.17	0.45
1:A:369:GLN:O	1:A:390:ARG:HD3	2.16	0.45
1:G:66:GLN:HE22	1:G:122:ALA:H	1.64	0.45
1:A:210:GLU:CG	1:E:43:ILE:CG1	2.94	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101:HIS:HB3	1:B:277:GLN:OE1	2.17	0.45
1:C:46:GLY:O	1:C:47:LEU:C	2.60	0.45
1:G:236:GLU:O	3:G:403:GOL:H32	2.16	0.45
1:A:231:LEU:HD21	2:A:400:COA:H1B	1.98	0.45
1:A:264:ARG:C	1:G:1:MET:HE1	2.38	0.45
1:E:318:PHE:O	1:E:319:ALA:C	2.59	0.45
1:A:177:GLN:OE1	1:A:181:GLN:NE2	2.46	0.45
1:A:265:ALA:CB	1:G:1:MET:HE1	2.46	0.45
1:D:197:GLN:HG2	1:D:211:ARG:CZ	2.47	0.45
1:H:57:GLY:HA2	1:H:118:CYS:O	2.16	0.45
1:E:18:ASN:N	1:E:124:SER:OG	2.44	0.45
1:H:290:GLY:N	1:H:291:PRO:CD	2.80	0.45
1:E:172:PHE:O	1:E:173:GLY:C	2.60	0.45
2:E:401:COA:O9P	2:E:401:COA:H62	2.17	0.45
1:A:210:GLU:HG3	1:E:43:ILE:HG13	1.97	0.45
1:D:373:LEU:HD12	1:D:386:THR:O	2.16	0.45
1:H:66:GLN:HE22	1:H:122:ALA:H	1.65	0.44
1:C:177:GLN:OE1	1:C:181:GLN:NE2	2.46	0.44
1:E:325:TRP:CZ3	1:E:331:PRO:HG3	2.52	0.44
1:A:82:GLU:O	1:D:281:GLY:HA3	2.17	0.44
1:B:48[A]:HIS:CE1	1:B:51:ASP:OD1	2.70	0.44
1:B:369:GLN:O	1:B:390:ARG:HD3	2.16	0.44
1:A:213:LEU:HD23	1:E:215:PHE:CE2	2.52	0.44
1:B:260:GLU:OE1	1:B:273:ARG:NH2	2.51	0.44
1:F:316:GLU:CD	1:F:341:GLY:HA3	2.43	0.44
1:C:197:GLN:HG2	1:C:211:ARG:CZ	2.48	0.44
1:C:150:ASN:OD1	1:C:152:PHE:N	2.51	0.44
1:H:369:GLN:O	1:H:390:ARG:HD3	2.17	0.44
1:A:213:LEU:CD2	1:E:215:PHE:CE2	3.00	0.44
1:C:325:TRP:CZ3	1:C:331:PRO:HG3	2.53	0.44
1:F:360:ALA:O	1:F:361:LEU:C	2.60	0.44
1:B:347:HIS:O	1:B:347:HIS:CG	2.68	0.43
1:A:10:THR:HG22	1:A:37:VAL:HG13	2.00	0.43
1:C:66:GLN:HE22	1:C:122:ALA:H	1.66	0.43
1:D:10:THR:HG22	1:D:37:VAL:HG13	1.99	0.43
1:D:290:GLY:N	1:D:291:PRO:CD	2.81	0.43
1:F:369:GLN:O	1:F:390:ARG:HD3	2.18	0.43
1:F:289:ASP:OD1	1:F:324:SER:OG	2.20	0.43
1:A:178:ARG:HH22	1:A:182:ARG:NH1	2.17	0.43
1:E:247:GLN:HB2	1:E:346:GLY:HA2	2.00	0.43
1:E:290:GLY:N	1:E:291:PRO:CD	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:0:SER:O	1:G:110:GLY:HA3	2.19	0.43
1:A:210:GLU:HG3	1:E:43:ILE:HG12	2.00	0.43
1:E:177:GLN:OE1	1:E:181:GLN:NE2	2.48	0.43
1:F:197:GLN:HG2	1:F:211:ARG:CZ	2.48	0.43
1:A:363:GLU:OE2	1:A:366:ARG:NH2	2.48	0.43
1:C:152:PHE:CD1	1:C:234:VAL:HG11	2.54	0.43
1:G:247:GLN:HB2	1:G:346:GLY:HA2	2.01	0.43
1:H:152:PHE:CD1	1:H:234:VAL:HG11	2.53	0.43
1:A:55:VAL:O	1:A:86:ALA:HA	2.18	0.43
1:E:184:TRP:HB3	3:E:402:GOL:H2	2.00	0.43
1:G:152:PHE:CE1	1:G:234:VAL:HG11	2.53	0.43
1:H:90:ASP:OD1	1:H:90:ASP:C	2.62	0.43
1:F:290:GLY:N	1:F:291:PRO:CD	2.82	0.43
1:D:325:TRP:CZ3	1:D:331:PRO:HG3	2.53	0.42
1:E:221:ARG:O	1:E:222:GLU:C	2.62	0.42
1:F:46:GLY:O	1:F:47:LEU:C	2.62	0.42
1:B:152:PHE:CD1	1:B:234:VAL:HG11	2.55	0.42
1:B:227:GLY:O	1:B:230:GLU:HB2	2.20	0.42
1:C:57:GLY:HA2	1:C:118:CYS:O	2.20	0.42
1:E:55:VAL:O	1:E:86:ALA:HA	2.20	0.42
1:E:197:GLN:HG2	1:E:211:ARG:CZ	2.49	0.42
1:B:363:GLU:OE2	1:B:366:ARG:NH2	2.52	0.42
1:D:273[A]:ARG:HB3	1:D:391:ILE:HD11	2.01	0.42
1:A:195:PRO:HG3	1:E:215:PHE:CZ	2.52	0.42
1:G:369:GLN:O	1:G:390:ARG:HD3	2.19	0.42
1:C:55:VAL:O	1:C:86:ALA:HA	2.19	0.42
1:G:316:GLU:CD	1:G:341:GLY:HA3	2.45	0.42
1:E:57:GLY:O	1:E:88:THR:HA	2.19	0.42
1:F:66:GLN:HE22	1:F:122:ALA:H	1.66	0.42
1:G:221:ARG:O	1:G:222:GLU:C	2.61	0.42
1:A:197:GLN:HG2	1:A:211:ARG:CZ	2.50	0.42
2:A:400:COA:H141	2:A:400:COA:O9P	2.19	0.42
1:D:171:VAL:O	1:D:172:PHE:C	2.62	0.42
2:H:401:COA:H62	2:H:401:COA:O9P	2.20	0.42
1:B:221:ARG:CZ	2:B:400:COA:H2B	2.50	0.42
1:F:178[B]:ARG:CG	1:F:178[B]:ARG:NH1	2.52	0.42
1:A:273[B]:ARG:NH1	1:A:274:ILE:O	2.54	0.41
1:D:60:VAL:HG21	1:D:349:VAL:HG12	2.02	0.41
1:C:290:GLY:N	1:C:291:PRO:CD	2.84	0.41
1:F:325:TRP:CZ3	1:F:331:PRO:HG3	2.55	0.41
2:G:401:COA:O9P	2:G:401:COA:H62	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:369:GLN:O	1:C:390:ARG:HD3	2.19	0.41
1:D:316:GLU:CD	1:D:341:GLY:HA3	2.46	0.41
1:F:221:ARG:O	1:F:222:GLU:C	2.64	0.41
1:C:216:ARG:NH2	4:C:501:HOH:O	2.44	0.41
1:F:7:VAL:HG11	1:F:268:LEU:HD13	2.02	0.41
1:B:7:VAL:HG11	1:B:268:LEU:HD13	2.01	0.41
1:C:10:THR:HG22	1:C:37:VAL:HG13	2.01	0.41
1:C:233:PRO:HA	1:C:241:THR:HG22	2.02	0.41
1:E:240:HIS:CE1	1:E:320:SER:H	2.31	0.41
1:G:39:ASP:O	1:G:40:LYS:C	2.64	0.41
1:G:150:ASN:OD1	1:G:152:PHE:N	2.54	0.41
1:A:221:ARG:CZ	2:A:400:COA:H2B	2.50	0.41
1:A:290:GLY:O	1:A:291:PRO:C	2.64	0.41
1:B:149:PRO:O	1:B:287:HIS:ND1	2.53	0.41
1:B:290:GLY:N	1:B:291:PRO:CD	2.84	0.41
1:F:39:ASP:O	1:F:40:LYS:C	2.62	0.41
1:E:360:ALA:O	1:E:361:LEU:C	2.64	0.41
1:G:177:GLN:OE1	1:G:181:GLN:NE2	2.48	0.41
1:B:145:ASP:O	1:B:145:ASP:CG	2.63	0.41
1:B:197:GLN:HG2	1:B:211:ARG:CZ	2.51	0.41
1:E:233:PRO:HA	1:E:241:THR:HG22	2.03	0.41
1:G:221:ARG:CZ	2:G:401:COA:H2B	2.51	0.41
1:H:177:GLN:OE1	1:H:181:GLN:NE2	2.50	0.41
1:A:150:ASN:OD1	1:A:152:PHE:N	2.53	0.41
1:A:290:GLY:O	1:A:293:GLN:N	2.54	0.41
1:B:233:PRO:HA	1:B:241:THR:HG22	2.03	0.41
1:C:221:ARG:CZ	2:C:400:COA:H2B	2.51	0.41
2:E:401:COA:O9P	2:E:401:COA:C6P	2.69	0.41
1:G:290:GLY:N	1:G:291:PRO:CD	2.84	0.41
1:A:152:PHE:CD1	1:A:234:VAL:HG11	2.56	0.40
1:F:226:ALA:O	1:F:227:GLY:C	2.64	0.40
1:F:247:GLN:HB2	1:F:346:GLY:HA2	2.03	0.40
1:G:197:GLN:HG2	1:G:211:ARG:CZ	2.51	0.40
1:H:221:ARG:CZ	2:H:401:COA:H2B	2.51	0.40
1:B:247:GLN:HB2	1:B:346:GLY:HA2	2.03	0.40
2:C:400:COA:O9P	2:C:400:COA:H62	2.20	0.40
1:D:221:ARG:CZ	2:D:400:COA:H2B	2.51	0.40
1:E:221:ARG:CZ	2:E:401:COA:H2B	2.51	0.40
1:G:149:PRO:O	1:G:287:HIS:ND1	2.52	0.40
1:G:236:GLU:O	3:G:403:GOL:C3	2.70	0.40
1:H:150:ASN:OD1	1:H:152:PHE:N	2.54	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	389/399 (98%)	375 (96%)	14 (4%)	0	100	100
1	B	390/399 (98%)	376 (96%)	13 (3%)	1 (0%)	36	70
1	C	388/399 (97%)	373 (96%)	14 (4%)	1 (0%)	36	70
1	D	389/399 (98%)	377 (97%)	11 (3%)	1 (0%)	36	70
1	E	379/399 (95%)	366 (97%)	13 (3%)	0	100	100
1	F	385/399 (96%)	371 (96%)	14 (4%)	0	100	100
1	G	391/399 (98%)	376 (96%)	14 (4%)	1 (0%)	36	70
1	H	393/399 (98%)	375 (95%)	16 (4%)	2 (0%)	24	60
All	All	3104/3192 (97%)	2989 (96%)	109 (4%)	6 (0%)	43	76

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	0	SER
1	H	1	MET
1	D	130	ALA
1	B	130	ALA
1	C	130	ALA
1	H	-1	GLY

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	296/302 (98%)	287 (97%)	9 (3%)	36	69
1	B	297/302 (98%)	291 (98%)	6 (2%)	48	76
1	C	295/302 (98%)	287 (97%)	8 (3%)	39	71
1	D	296/302 (98%)	289 (98%)	7 (2%)	43	73
1	E	287/302 (95%)	278 (97%)	9 (3%)	35	68
1	F	293/302 (97%)	287 (98%)	6 (2%)	48	76
1	G	297/302 (98%)	291 (98%)	6 (2%)	48	76
1	H	299/302 (99%)	289 (97%)	10 (3%)	33	67
All	All	2360/2416 (98%)	2299 (97%)	61 (3%)	42	72

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	VAL
1	A	140	ARG
1	A	153	GLU
1	A	197	GLN
1	A	215	PHE
1	A	273[A]	ARG
1	A	273[B]	ARG
1	A	327	ARG
1	A	386	THR
1	B	37	VAL
1	B	140	ARG
1	B	153	GLU
1	B	197	GLN
1	B	215	PHE
1	B	327	ARG
1	C	37	VAL
1	C	140	ARG
1	C	153	GLU
1	C	178	ARG
1	C	197	GLN
1	C	215	PHE
1	C	327	ARG
1	C	375	THR
1	D	37	VAL
1	D	140	ARG

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Mol	Chain	Res	Type
1	D	153	GLU
1	D	197	GLN
1	D	215	PHE
1	D	327	ARG
1	D	375	THR
1	E	37	VAL
1	E	140	ARG
1	E	153	GLU
1	E	197	GLN
1	E	215	PHE
1	E	327	ARG
1	E	349	VAL
1	E	351	CYS
1	E	375	THR
1	F	37	VAL
1	F	140	ARG
1	F	153	GLU
1	F	197	GLN
1	F	215	PHE
1	F	327	ARG
1	G	37	VAL
1	G	140	ARG
1	G	153	GLU
1	G	197	GLN
1	G	215	PHE
1	G	327	ARG
1	H	3[A]	TYR
1	H	3[B]	TYR
1	H	10	THR
1	H	37	VAL
1	H	140	ARG
1	H	153	GLU
1	H	197	GLN
1	H	215	PHE
1	H	273	ARG
1	H	327	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	68	ASN

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Mol	Chain	Res	Type
1	A	369	GLN
1	B	66	GLN
1	B	68	ASN
1	B	92	GLN
1	B	369	GLN
1	C	66	GLN
1	C	68	ASN
1	C	92	GLN
1	C	369	GLN
1	D	48	HIS
1	D	66	GLN
1	D	68	ASN
1	D	369	GLN
1	E	66	GLN
1	E	68	ASN
1	E	69	ASN
1	E	369	GLN
1	F	66	GLN
1	F	68	ASN
1	F	92	GLN
1	F	206	GLN
1	F	369	GLN
1	G	66	GLN
1	G	68	ASN
1	G	206	GLN
1	G	369	GLN
1	H	66	GLN
1	H	68	ASN
1	H	369	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 non-standard protein/DNA/RNA residues 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
1	OAS	E	93	1	7,8,9	1.37	1 (14%)	4,9,11	2.77	1 (25%)
1	OAS	G	93	1	7,8,9	1.57	1 (14%)	4,9,11	2.55	2 (50%)
1	OAS	H	93	1	7,8,9	1.26	1 (14%)	4,9,11	2.46	2 (50%)
1	OAS	B	93	1	7,8,9	1.39	1 (14%)	4,9,11	2.30	1 (25%)
1	OAS	D	93	1	7,8,9	1.43	1 (14%)	4,9,11	1.96	1 (25%)
1	OAS	F	93	1	7,8,9	1.51	1 (14%)	4,9,11	2.50	3 (75%)
1	OAS	A	93	1	7,8,9	1.46	1 (14%)	4,9,11	2.33	1 (25%)
1	OAS	C	93	1	7,8,9	1.50	1 (14%)	4,9,11	2.07	2 (50%)

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
1	OAS	E	93	1	-	2/5/7/9	-
1	OAS	G	93	1	-	2/5/7/9	-
1	OAS	H	93	1	-	2/5/7/9	-
1	OAS	B	93	1	-	2/5/7/9	-
1	OAS	D	93	1	-	2/5/7/9	-
1	OAS	F	93	1	-	2/5/7/9	-
1	OAS	A	93	1	-	2/5/7/9	-
1	OAS	C	93	1	-	2/5/7/9	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	93	OAS	OG-C1A	3.86	1.51	1.33
1	F	93	OAS	OG-C1A	3.68	1.51	1.33
1	C	93	OAS	OG-C1A	3.54	1.50	1.33
1	B	93	OAS	OG-C1A	3.48	1.50	1.33
1	A	93	OAS	OG-C1A	3.42	1.49	1.33
1	D	93	OAS	OG-C1A	3.28	1.49	1.33
1	E	93	OAS	OG-C1A	3.12	1.48	1.33
1	H	93	OAS	OG-C1A	3.05	1.48	1.33

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	93	OAS	CB-OG-C1A	5.33	130.23	117.08
1	G	93	OAS	CB-OG-C1A	4.35	127.80	117.08
1	A	93	OAS	CB-OG-C1A	4.16	127.36	117.08
1	H	93	OAS	CB-OG-C1A	4.15	127.32	117.08
1	B	93	OAS	CB-OG-C1A	4.15	127.31	117.08
1	F	93	OAS	CB-OG-C1A	3.91	126.73	117.08
1	D	93	OAS	CB-OG-C1A	3.43	125.55	117.08
1	C	93	OAS	CB-OG-C1A	3.12	124.78	117.08
1	F	93	OAS	OAC-C1A-C2A	-2.33	116.51	124.77
1	H	93	OAS	OG-C1A-C2A	2.20	121.73	112.38
1	G	93	OAS	OAC-C1A-C2A	-2.17	117.09	124.77
1	C	93	OAS	OG-C1A-C2A	2.05	121.06	112.38
1	F	93	OAS	OG-C1A-C2A	2.03	121.00	112.38

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	93	OAS	OAC-C1A-OG-CB
1	B	93	OAS	OAC-C1A-OG-CB
1	D	93	OAS	OAC-C1A-OG-CB
1	E	93	OAS	OAC-C1A-OG-CB
1	G	93	OAS	OAC-C1A-OG-CB
1	B	93	OAS	C2A-C1A-OG-CB
1	D	93	OAS	C2A-C1A-OG-CB
1	E	93	OAS	C2A-C1A-OG-CB
1	G	93	OAS	C2A-C1A-OG-CB
1	C	93	OAS	OAC-C1A-OG-CB
1	F	93	OAS	OAC-C1A-OG-CB
1	H	93	OAS	OAC-C1A-OG-CB
1	H	93	OAS	C2A-C1A-OG-CB
1	C	93	OAS	C2A-C1A-OG-CB
1	F	93	OAS	C2A-C1A-OG-CB
1	A	93	OAS	C2A-C1A-OG-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

12 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	COA	B	400	-	47,50,50	1.22	4 (8%)	69,75,75	1.78	14 (20%)
2	COA	G	401	-	47,50,50	1.21	4 (8%)	69,75,75	1.76	15 (21%)
3	GOL	E	402	-	5,5,5	0.64	0	5,5,5	1.90	2 (40%)
2	COA	C	400	-	47,50,50	1.23	5 (10%)	69,75,75	1.71	13 (18%)
2	COA	H	401	-	47,50,50	1.23	5 (10%)	69,75,75	1.71	13 (18%)
3	GOL	H	402	-	5,5,5	0.66	0	5,5,5	1.49	1 (20%)
2	COA	E	401	-	47,50,50	1.29	7 (14%)	69,75,75	1.77	15 (21%)
3	GOL	G	403	-	5,5,5	0.79	0	5,5,5	1.25	1 (20%)
2	COA	F	400	-	47,50,50	1.27	6 (12%)	69,75,75	1.74	13 (18%)
3	GOL	G	402	-	5,5,5	0.61	0	5,5,5	1.76	1 (20%)
2	COA	D	400	-	47,50,50	1.20	4 (8%)	69,75,75	1.79	14 (20%)
2	COA	A	400	-	47,50,50	1.21	5 (10%)	69,75,75	1.75	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
2	COA	B	400	-	-	14/48/64/64	0/3/3/3
2	COA	G	401	-	-	14/48/64/64	0/3/3/3
3	GOL	E	402	-	-	1/4/4/4	-
2	COA	C	400	-	-	14/48/64/64	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	COA	H	401	-	-	15/48/64/64	0/3/3/3
3	GOL	H	402	-	-	4/4/4/4	-
2	COA	E	401	-	-	15/48/64/64	0/3/3/3
3	GOL	G	403	-	-	2/4/4/4	-
2	COA	F	400	-	-	15/48/64/64	0/3/3/3
3	GOL	G	402	-	-	4/4/4/4	-
2	COA	D	400	-	-	14/48/64/64	0/3/3/3
2	COA	A	400	-	-	15/48/64/64	0/3/3/3

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	400	COA	C5A-C4A	4.79	1.47	1.39
2	F	400	COA	C5A-C4A	4.77	1.47	1.39
2	H	401	COA	C5A-C4A	4.74	1.47	1.39
2	D	400	COA	C5A-C4A	4.68	1.47	1.39
2	A	400	COA	C5A-C4A	4.62	1.47	1.39
2	G	401	COA	C5A-C4A	4.60	1.47	1.39
2	B	400	COA	C5A-C4A	4.59	1.47	1.39
2	E	401	COA	C5A-C4A	4.51	1.47	1.39
2	F	400	COA	C5A-C6A	3.14	1.49	1.41
2	D	400	COA	C5A-C6A	3.10	1.49	1.41
2	B	400	COA	C5A-C6A	3.05	1.49	1.41
2	E	401	COA	P1A-O3A	3.04	1.62	1.59
2	G	401	COA	C5A-C6A	3.03	1.49	1.41
2	E	401	COA	C5A-C6A	2.96	1.49	1.41
2	C	400	COA	C5A-C6A	2.91	1.49	1.41
2	A	400	COA	C5A-C6A	2.91	1.49	1.41
2	C	400	COA	C8A-N7A	2.91	1.37	1.31
2	H	401	COA	C5A-C6A	2.76	1.48	1.41
2	D	400	COA	C8A-N7A	2.72	1.36	1.31
2	F	400	COA	C8A-N7A	2.70	1.36	1.31
2	H	401	COA	C8A-N7A	2.69	1.36	1.31
2	G	401	COA	C8A-N7A	2.65	1.36	1.31
2	A	400	COA	C8A-N7A	2.61	1.36	1.31
2	E	401	COA	C8A-N7A	2.55	1.36	1.31
2	E	401	COA	P2A-O3A	2.53	1.62	1.59
2	B	400	COA	C8A-N7A	2.52	1.36	1.31
2	H	401	COA	C5A-N7A	-2.34	1.34	1.39
2	B	400	COA	C5A-N7A	-2.28	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	400	COA	C5A-N7A	-2.25	1.35	1.39
2	F	400	COA	P2A-O3A	2.23	1.61	1.59
2	G	401	COA	C5A-N7A	-2.18	1.35	1.39
2	H	401	COA	P2A-O3A	2.13	1.61	1.59
2	E	401	COA	C5A-N7A	-2.13	1.35	1.39
2	F	400	COA	C5A-N7A	-2.11	1.35	1.39
2	F	400	COA	P1A-O3A	2.09	1.61	1.59
2	A	400	COA	C5A-N7A	-2.07	1.35	1.39
2	D	400	COA	C5A-N7A	-2.05	1.35	1.39
2	A	400	COA	C4A-N9A	-2.05	1.33	1.37
2	C	400	COA	C4A-N9A	-2.02	1.33	1.37
2	E	401	COA	C4A-N9A	-2.00	1.33	1.37

All (116) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	401	COA	C5A-C4A-N3A	-5.32	119.39	126.72
2	B	400	COA	C5A-C4A-N3A	-5.30	119.42	126.72
2	C	400	COA	C5A-C4A-N3A	-5.20	119.56	126.72
2	D	400	COA	C5A-C4A-N3A	-5.19	119.57	126.72
2	A	400	COA	C5A-C4A-N3A	-5.18	119.59	126.72
2	E	401	COA	C5A-C4A-N3A	-5.10	119.70	126.72
2	F	400	COA	C5A-C4A-N3A	-5.06	119.75	126.72
2	H	401	COA	C5A-C4A-N3A	-5.03	119.79	126.72
2	E	401	COA	N3A-C2A-N1A	-4.48	121.80	128.58
2	B	400	COA	O4B-C1B-N9A	4.44	116.62	108.09
2	F	400	COA	O4B-C1B-N9A	4.38	116.50	108.09
2	A	400	COA	O4B-C1B-N9A	4.35	116.44	108.09
2	B	400	COA	C4A-C5A-N7A	-4.34	105.62	110.58
2	D	400	COA	C4A-C5A-N7A	-4.32	105.65	110.58
2	E	401	COA	O4B-C1B-N9A	4.31	116.37	108.09
2	H	401	COA	O4B-C1B-N9A	4.28	116.31	108.09
2	D	400	COA	O4B-C1B-N9A	4.22	116.19	108.09
2	G	401	COA	O4B-C1B-N9A	4.22	116.19	108.09
2	H	401	COA	N3A-C2A-N1A	-4.20	122.23	128.58
2	C	400	COA	O4B-C1B-N9A	4.19	116.13	108.09
2	G	401	COA	C4A-C5A-N7A	-4.18	105.80	110.58
2	D	400	COA	N3A-C2A-N1A	-4.15	122.30	128.58
2	A	400	COA	N3A-C2A-N1A	-4.15	122.30	128.58
2	F	400	COA	C4A-C5A-N7A	-4.10	105.89	110.58
2	B	400	COA	N3A-C2A-N1A	-4.07	122.42	128.58
2	E	401	COA	C2A-N3A-C4A	4.06	121.74	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	400	COA	N3A-C2A-N1A	-4.03	122.48	128.58
2	A	400	COA	C4A-C5A-N7A	-4.03	105.97	110.58
2	C	400	COA	C4A-C5A-N7A	-4.00	106.01	110.58
2	C	400	COA	N3A-C2A-N1A	-4.00	122.53	128.58
2	G	401	COA	N3A-C2A-N1A	-3.96	122.59	128.58
2	B	400	COA	C2A-N3A-C4A	3.93	121.42	111.83
2	H	401	COA	C4A-C5A-N7A	-3.90	106.12	110.58
2	G	401	COA	C2A-N3A-C4A	3.89	121.33	111.83
2	A	400	COA	C2A-N3A-C4A	3.87	121.29	111.83
2	D	400	COA	C2A-N3A-C4A	3.86	121.25	111.83
2	H	401	COA	C2A-N3A-C4A	3.83	121.18	111.83
2	E	401	COA	C4A-C5A-N7A	-3.83	106.21	110.58
2	C	400	COA	C2A-N3A-C4A	3.81	121.14	111.83
2	F	400	COA	C2A-N3A-C4A	3.79	121.09	111.83
2	B	400	COA	N3A-C4A-N9A	3.77	133.58	127.17
2	G	401	COA	N3A-C4A-N9A	3.74	133.53	127.17
2	A	400	COA	N3A-C4A-N9A	3.74	133.53	127.17
2	F	400	COA	N3A-C4A-N9A	3.73	133.52	127.17
2	D	400	COA	N3A-C4A-N9A	3.70	133.46	127.17
2	E	401	COA	N3A-C4A-N9A	3.69	133.44	127.17
2	D	400	COA	C4B-O4B-C1B	-3.66	101.39	109.47
2	C	400	COA	N3A-C4A-N9A	3.61	133.31	127.17
2	A	400	COA	C4B-O4B-C1B	-3.59	101.54	109.47
2	H	401	COA	N3A-C4A-N9A	3.58	133.25	127.17
2	F	400	COA	C4B-O4B-C1B	-3.52	101.70	109.47
2	C	400	COA	C4B-O4B-C1B	-3.48	101.78	109.47
2	G	401	COA	C4B-O4B-C1B	-3.44	101.88	109.47
2	B	400	COA	C4B-O4B-C1B	-3.42	101.92	109.47
2	E	401	COA	C4B-O4B-C1B	-3.39	101.98	109.47
2	H	401	COA	C4B-O4B-C1B	-3.39	101.98	109.47
2	B	400	COA	C5A-N7A-C8A	3.24	108.55	103.45
2	D	400	COA	C5A-N7A-C8A	3.23	108.53	103.45
2	B	400	COA	C6A-C5A-N7A	3.12	138.11	132.09
2	D	400	COA	C6A-C5A-N7A	3.08	138.03	132.09
2	F	400	COA	C5A-N7A-C8A	3.03	108.22	103.45
2	G	401	COA	C6A-C5A-N7A	2.99	137.86	132.09
2	E	401	COA	C6A-C5A-N7A	2.99	137.86	132.09
2	G	401	COA	C5A-N7A-C8A	2.99	108.15	103.45
2	F	400	COA	C4A-N9A-C8A	2.96	108.85	105.74
2	F	400	COA	C6A-C5A-N7A	2.96	137.79	132.09
3	E	402	GOL	O3-C3-C2	2.95	123.65	110.38
2	A	400	COA	C5A-N7A-C8A	2.93	108.06	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	401	COA	C5A-N7A-C8A	2.93	108.06	103.45
2	A	400	COA	C6A-C5A-N7A	2.90	137.68	132.09
2	H	401	COA	C6A-C5A-N7A	2.89	137.66	132.09
2	C	400	COA	C5A-N7A-C8A	2.88	107.98	103.45
2	A	400	COA	C4A-N9A-C8A	2.87	108.75	105.74
2	C	400	COA	C6A-C5A-N7A	2.86	137.60	132.09
2	D	400	COA	C4A-N9A-C8A	2.85	108.73	105.74
2	E	401	COA	C5A-N7A-C8A	2.84	107.92	103.45
2	B	400	COA	C4A-N9A-C8A	2.72	108.59	105.74
2	E	401	COA	C4A-N9A-C8A	2.70	108.57	105.74
2	D	400	COA	N9A-C8A-N7A	-2.69	110.11	113.94
2	E	401	COA	O5P-C5P-C6P	-2.67	117.19	122.02
2	G	401	COA	C4A-N9A-C8A	2.66	108.53	105.74
2	B	400	COA	N9A-C8A-N7A	-2.62	110.22	113.94
3	G	402	GOL	O3-C3-C2	2.57	121.95	110.38
2	F	400	COA	N9A-C8A-N7A	-2.57	110.30	113.94
2	C	400	COA	C4A-N9A-C8A	2.56	108.42	105.74
2	A	400	COA	N9A-C8A-N7A	-2.55	110.33	113.94
2	H	401	COA	C4A-N9A-C8A	2.48	108.35	105.74
2	D	400	COA	C2A-N1A-C6A	2.47	122.79	118.73
2	E	401	COA	N9A-C8A-N7A	-2.47	110.43	113.94
2	G	401	COA	N9A-C8A-N7A	-2.47	110.44	113.94
2	C	400	COA	N9A-C8A-N7A	-2.45	110.45	113.94
2	H	401	COA	C2A-N1A-C6A	2.42	122.70	118.73
2	H	401	COA	N9A-C8A-N7A	-2.42	110.51	113.94
2	E	401	COA	C2A-N1A-C6A	2.39	122.65	118.73
2	F	400	COA	C2A-N1A-C6A	2.39	122.65	118.73
2	A	400	COA	C2A-N1A-C6A	2.37	122.62	118.73
2	B	400	COA	C4A-N9A-C1B	-2.36	121.11	126.63
2	F	400	COA	C4A-N9A-C1B	-2.34	121.15	126.63
2	D	400	COA	C4A-N9A-C1B	-2.34	121.17	126.63
2	C	400	COA	C2A-N1A-C6A	2.33	122.56	118.73
2	G	401	COA	C4A-N9A-C1B	-2.31	121.23	126.63
2	B	400	COA	O5P-C5P-C6P	-2.30	117.85	122.02
2	E	401	COA	C4A-N9A-C1B	-2.28	121.29	126.63
2	A	400	COA	O5P-C5P-C6P	-2.27	117.91	122.02
2	G	401	COA	OAP-CAP-CBP	-2.26	104.95	110.18
2	H	401	COA	C4A-N9A-C1B	-2.25	121.38	126.63
2	C	400	COA	C4A-N9A-C1B	-2.24	121.39	126.63
2	A	400	COA	C4A-N9A-C1B	-2.24	121.41	126.63
2	G	401	COA	O5P-C5P-C6P	-2.22	118.00	122.02
3	G	403	GOL	O3-C3-C2	2.21	120.34	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	401	COA	C2A-N1A-C6A	2.15	122.26	118.73
2	B	400	COA	C2A-N1A-C6A	2.12	122.21	118.73
3	H	402	GOL	O1-C1-C2	2.11	119.90	110.38
2	E	401	COA	O5P-C5P-N4P	2.11	127.17	123.03
2	D	400	COA	O6A-CCP-CBP	2.08	113.89	110.55
3	E	402	GOL	O1-C1-C2	2.05	119.59	110.38

There are no chirality outliers.

All (127) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	400	COA	CCP-O6A-P2A-O3A
2	A	400	COA	CCP-O6A-P2A-O4A
2	A	400	COA	CDP-CBP-CCP-O6A
2	A	400	COA	CEP-CBP-CCP-O6A
2	A	400	COA	CAP-CBP-CCP-O6A
2	B	400	COA	CCP-O6A-P2A-O3A
2	B	400	COA	CCP-O6A-P2A-O4A
2	B	400	COA	CDP-CBP-CCP-O6A
2	B	400	COA	CEP-CBP-CCP-O6A
2	B	400	COA	CAP-CBP-CCP-O6A
2	B	400	COA	S1P-C2P-C3P-N4P
2	C	400	COA	CCP-O6A-P2A-O3A
2	C	400	COA	CCP-O6A-P2A-O4A
2	C	400	COA	CDP-CBP-CCP-O6A
2	C	400	COA	CEP-CBP-CCP-O6A
2	C	400	COA	CAP-CBP-CCP-O6A
2	C	400	COA	S1P-C2P-C3P-N4P
2	D	400	COA	CCP-O6A-P2A-O3A
2	D	400	COA	CCP-O6A-P2A-O4A
2	D	400	COA	CDP-CBP-CCP-O6A
2	D	400	COA	CEP-CBP-CCP-O6A
2	D	400	COA	CAP-CBP-CCP-O6A
2	D	400	COA	S1P-C2P-C3P-N4P
2	E	401	COA	CCP-O6A-P2A-O3A
2	E	401	COA	CCP-O6A-P2A-O4A
2	E	401	COA	CDP-CBP-CCP-O6A
2	E	401	COA	CEP-CBP-CCP-O6A
2	E	401	COA	CAP-CBP-CCP-O6A
2	E	401	COA	S1P-C2P-C3P-N4P
2	F	400	COA	CCP-O6A-P2A-O3A
2	F	400	COA	CCP-O6A-P2A-O4A

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Mol	Chain	Res	Type	Atoms
2	F	400	COA	CDP-CBP-CCP-O6A
2	F	400	COA	CEP-CBP-CCP-O6A
2	F	400	COA	CAP-CBP-CCP-O6A
2	F	400	COA	S1P-C2P-C3P-N4P
2	G	401	COA	CCP-O6A-P2A-O3A
2	G	401	COA	CCP-O6A-P2A-O4A
2	G	401	COA	CDP-CBP-CCP-O6A
2	G	401	COA	CEP-CBP-CCP-O6A
2	G	401	COA	CAP-CBP-CCP-O6A
2	G	401	COA	S1P-C2P-C3P-N4P
2	H	401	COA	CCP-O6A-P2A-O3A
2	H	401	COA	CCP-O6A-P2A-O4A
2	H	401	COA	CDP-CBP-CCP-O6A
2	H	401	COA	CEP-CBP-CCP-O6A
2	H	401	COA	CAP-CBP-CCP-O6A
2	H	401	COA	S1P-C2P-C3P-N4P
3	G	402	GOL	O1-C1-C2-C3
3	G	402	GOL	C1-C2-C3-O3
3	G	403	GOL	C1-C2-C3-O3
3	H	402	GOL	C1-C2-C3-O3
2	E	401	COA	C6P-C7P-N8P-C9P
2	G	401	COA	C6P-C7P-N8P-C9P
2	A	400	COA	C6P-C7P-N8P-C9P
2	B	400	COA	C6P-C7P-N8P-C9P
2	C	400	COA	C6P-C7P-N8P-C9P
2	D	400	COA	C6P-C7P-N8P-C9P
2	F	400	COA	C6P-C7P-N8P-C9P
2	H	401	COA	C6P-C7P-N8P-C9P
2	B	400	COA	O4B-C4B-C5B-O5B
2	A	400	COA	C2B-C3B-O3B-P3B
2	B	400	COA	C2B-C3B-O3B-P3B
2	C	400	COA	C2B-C3B-O3B-P3B
2	D	400	COA	C2B-C3B-O3B-P3B
2	E	401	COA	C2B-C3B-O3B-P3B
2	F	400	COA	C2B-C3B-O3B-P3B
2	G	401	COA	C2B-C3B-O3B-P3B
2	A	400	COA	O4B-C4B-C5B-O5B
2	D	400	COA	O4B-C4B-C5B-O5B
2	E	401	COA	O4B-C4B-C5B-O5B
2	F	400	COA	O4B-C4B-C5B-O5B
2	G	401	COA	O4B-C4B-C5B-O5B
2	H	401	COA	O4B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
3	E	402	GOL	C1-C2-C3-O3
3	H	402	GOL	O1-C1-C2-C3
3	G	402	GOL	O1-C1-C2-O2
3	G	403	GOL	O2-C2-C3-O3
3	H	402	GOL	O1-C1-C2-O2
2	C	400	COA	O4B-C4B-C5B-O5B
2	H	401	COA	C2B-C3B-O3B-P3B
3	H	402	GOL	O2-C2-C3-O3
2	D	400	COA	C4B-C3B-O3B-P3B
2	A	400	COA	C3B-C4B-C5B-O5B
2	B	400	COA	C3B-C4B-C5B-O5B
2	D	400	COA	C3B-C4B-C5B-O5B
2	H	401	COA	C3B-C4B-C5B-O5B
3	G	402	GOL	O2-C2-C3-O3
2	E	401	COA	C3B-C4B-C5B-O5B
2	F	400	COA	C3B-C4B-C5B-O5B
2	G	401	COA	C3B-C4B-C5B-O5B
2	C	400	COA	C3B-C4B-C5B-O5B
2	G	401	COA	C4B-C3B-O3B-P3B
2	C	400	COA	C4B-C3B-O3B-P3B
2	E	401	COA	C4B-C3B-O3B-P3B
2	A	400	COA	S1P-C2P-C3P-N4P
2	A	400	COA	CCP-O6A-P2A-O5A
2	B	400	COA	CCP-O6A-P2A-O5A
2	C	400	COA	CCP-O6A-P2A-O5A
2	D	400	COA	CCP-O6A-P2A-O5A
2	E	401	COA	CCP-O6A-P2A-O5A
2	F	400	COA	CCP-O6A-P2A-O5A
2	G	401	COA	CCP-O6A-P2A-O5A
2	H	401	COA	CCP-O6A-P2A-O5A
2	A	400	COA	C4B-C3B-O3B-P3B
2	B	400	COA	C4B-C3B-O3B-P3B
2	F	400	COA	C4B-C3B-O3B-P3B
2	H	401	COA	C4B-C3B-O3B-P3B
2	A	400	COA	O5P-C5P-C6P-C7P
2	E	401	COA	O5P-C5P-C6P-C7P
2	B	400	COA	O5P-C5P-C6P-C7P
2	C	400	COA	O5P-C5P-C6P-C7P
2	D	400	COA	O5P-C5P-C6P-C7P
2	F	400	COA	O5P-C5P-C6P-C7P
2	H	401	COA	O5P-C5P-C6P-C7P
2	G	401	COA	O5P-C5P-C6P-C7P

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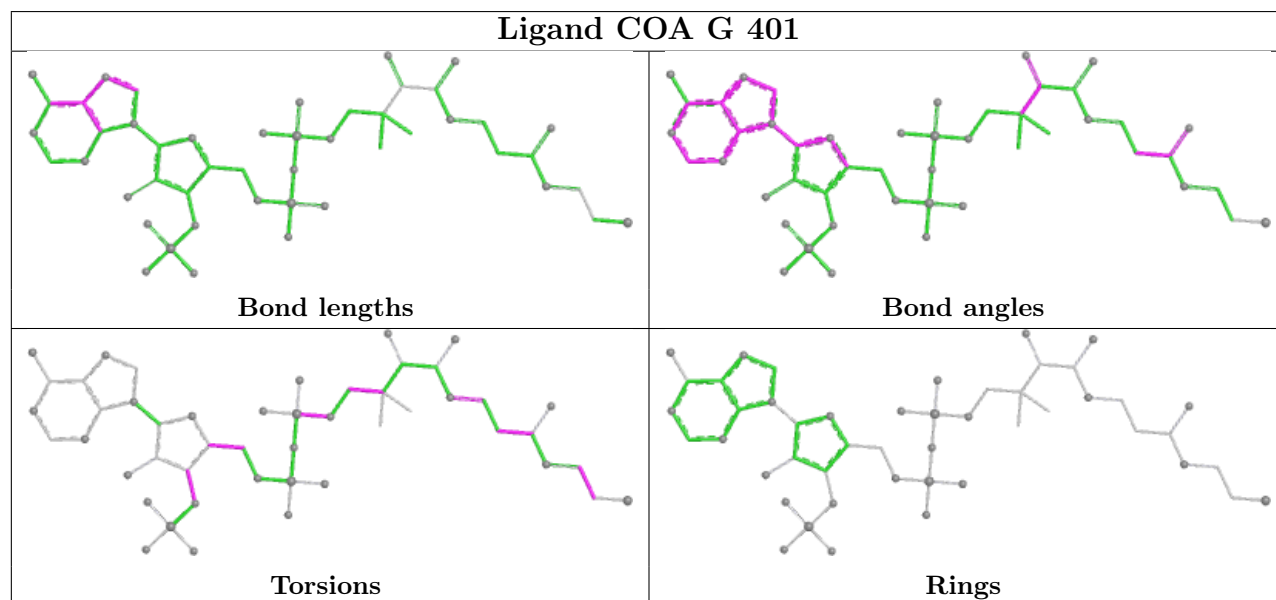
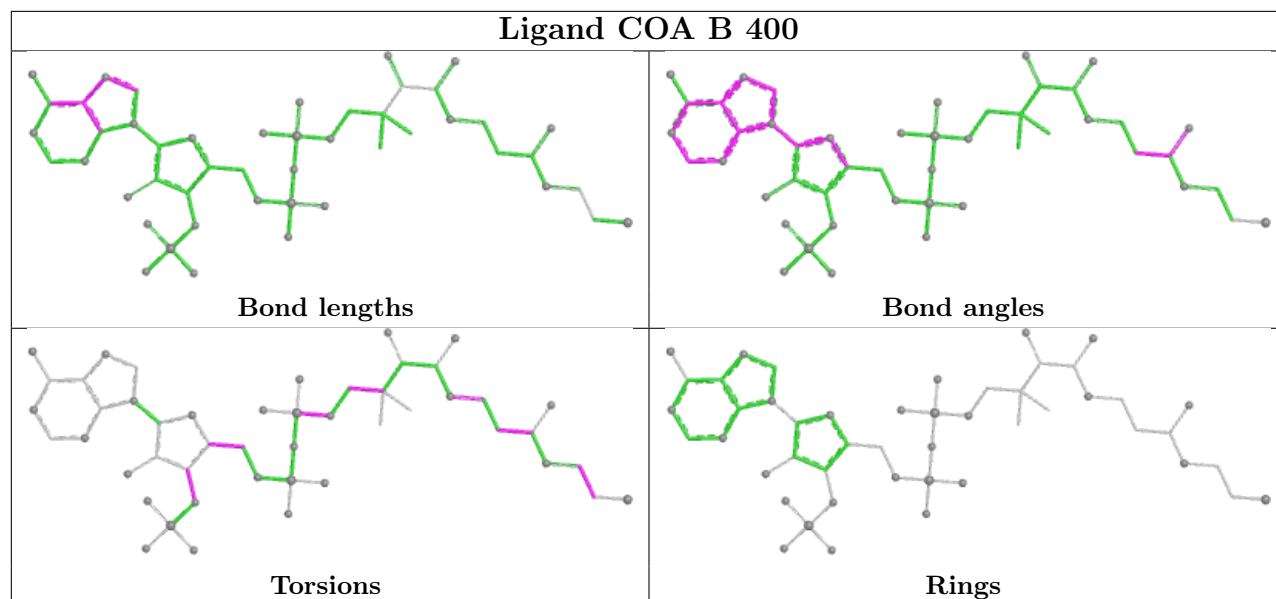
Mol	Chain	Res	Type	Atoms
2	A	400	COA	N4P-C5P-C6P-C7P
2	C	400	COA	N4P-C5P-C6P-C7P
2	E	401	COA	N4P-C5P-C6P-C7P
2	H	401	COA	N4P-C5P-C6P-C7P
2	B	400	COA	N4P-C5P-C6P-C7P
2	D	400	COA	N4P-C5P-C6P-C7P
2	F	400	COA	N4P-C5P-C6P-C7P
2	G	401	COA	N4P-C5P-C6P-C7P
2	A	400	COA	C3B-O3B-P3B-O8A
2	E	401	COA	C3B-O3B-P3B-O8A
2	F	400	COA	C3B-O3B-P3B-O8A
2	H	401	COA	C3B-O3B-P3B-O8A

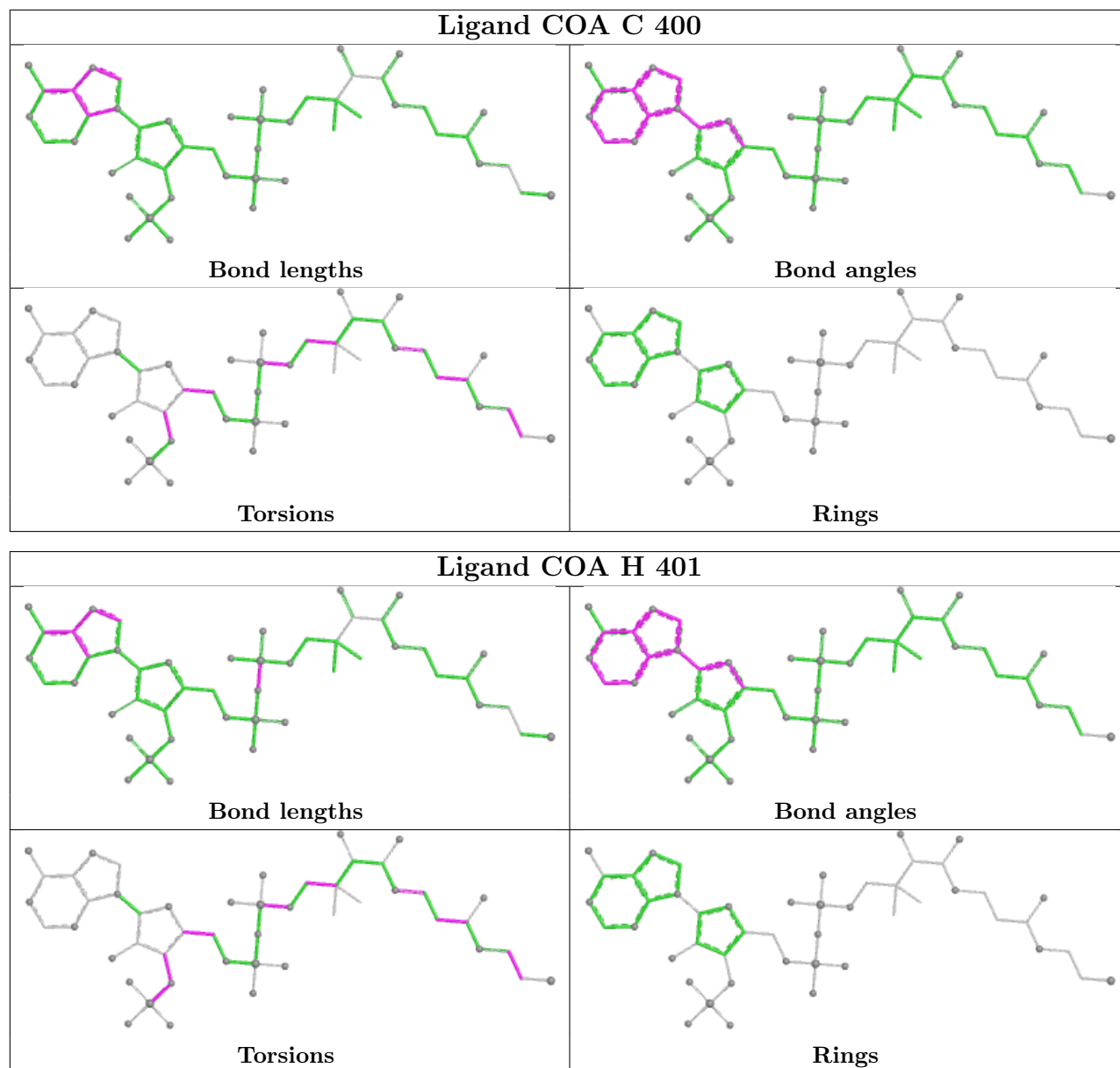
There are no ring outliers.

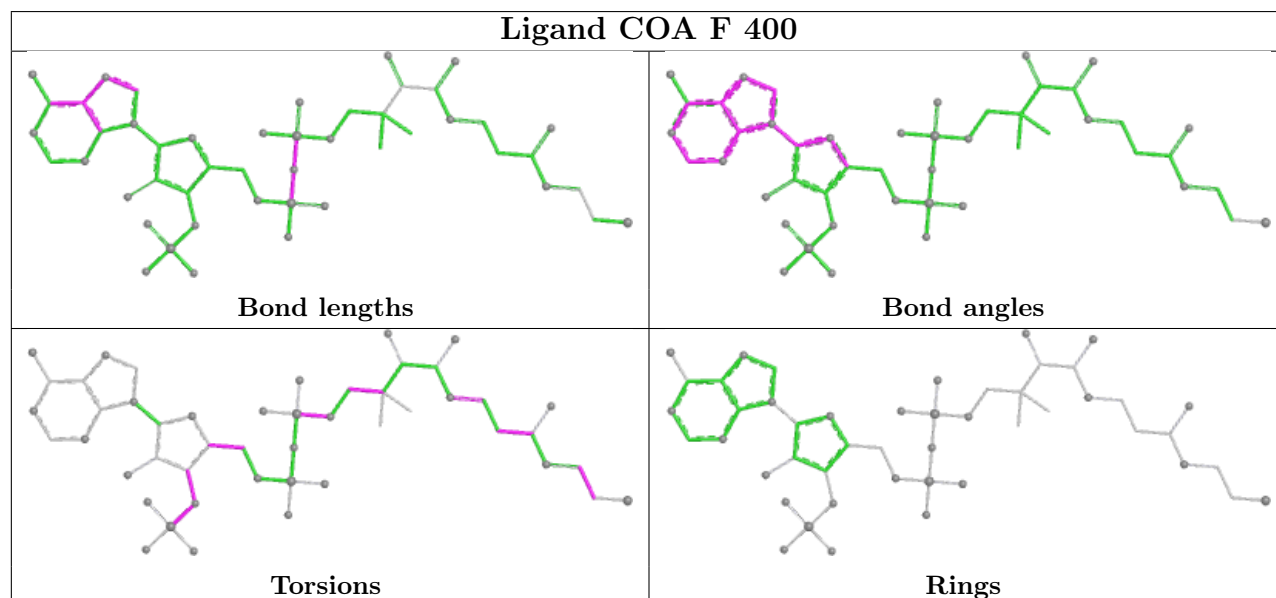
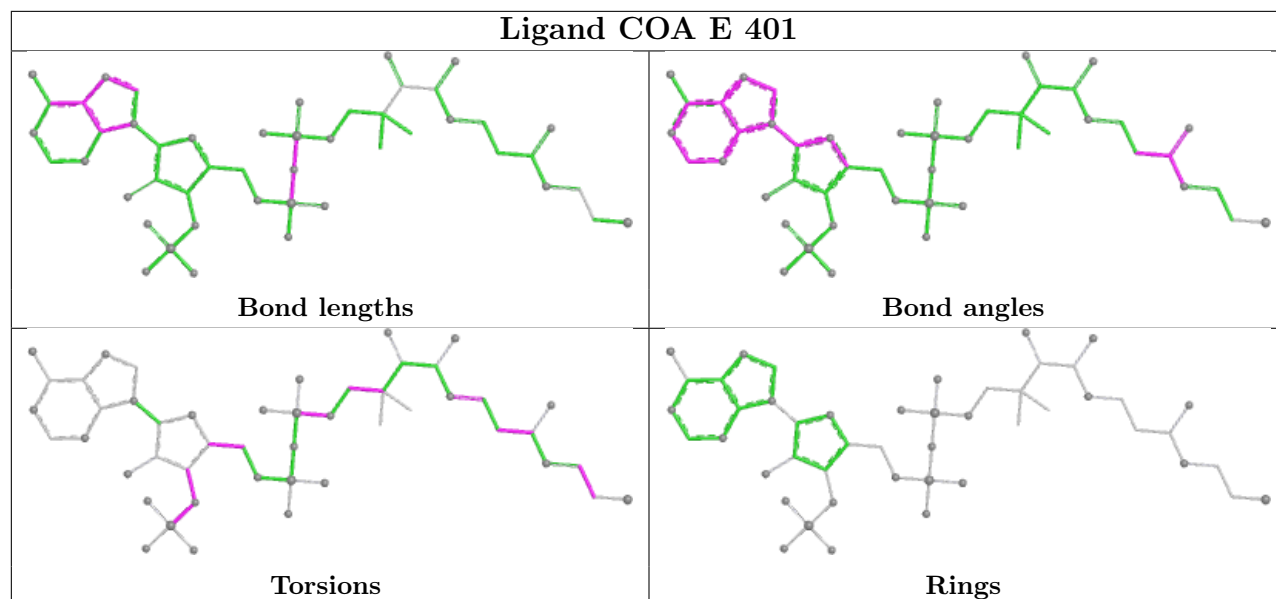
11 monomers are involved in 25 short contacts:

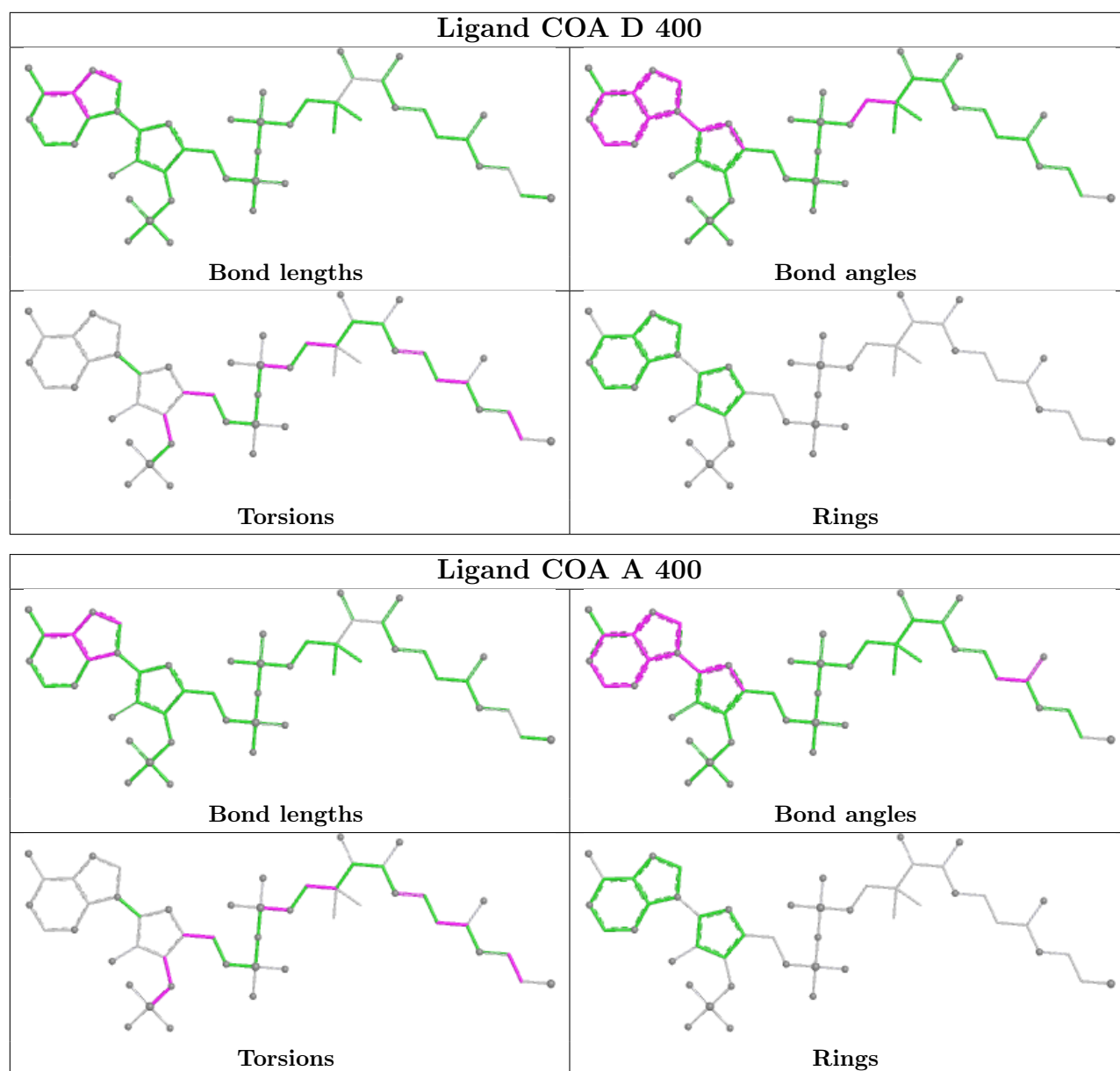
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	400	COA	2	0
2	G	401	COA	3	0
3	E	402	GOL	1	0
2	C	400	COA	3	0
2	H	401	COA	3	0
3	H	402	GOL	1	0
2	E	401	COA	4	0
3	G	403	GOL	2	0
2	F	400	COA	1	0
2	D	400	COA	2	0
2	A	400	COA	3	0

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

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, 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	390/399 (97%)	-0.21	5 (1%) 75 53	20, 34, 66, 185	2 (0%)
1	B	390/399 (97%)	-0.21	4 (1%) 79 59	22, 37, 67, 147	3 (0%)
1	C	390/399 (97%)	-0.16	9 (2%) 61 38	20, 38, 67, 173	1 (0%)
1	D	390/399 (97%)	-0.23	5 (1%) 75 53	20, 36, 61, 167	2 (0%)
1	E	378/399 (94%)	-0.18	6 (1%) 70 47	21, 36, 70, 132	2 (0%)
1	F	385/399 (96%)	-0.19	6 (1%) 70 47	20, 37, 65, 157	3 (0%)
1	G	392/399 (98%)	-0.22	8 (2%) 65 41	20, 36, 70, 179	2 (0%)
1	H	393/399 (98%)	-0.03	5 (1%) 75 53	21, 45, 74, 124	3 (0%)
All	All	3108/3192 (97%)	-0.18	48 (1%) 72 49	20, 37, 70, 185	18 (0%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	222	GLU	5.3
1	E	215	PHE	4.3
1	D	1	MET	3.9
1	H	-2	HIS	3.8
1	E	216	ARG	3.7
1	E	1	MET	3.7
1	B	1	MET	3.5
1	C	215	PHE	3.4
1	B	222	GLU	3.3
1	G	209	GLY	3.2
1	D	222	GLU	3.2
1	C	222	GLU	3.1
1	E	222	GLU	3.0
1	F	1	MET	3.0
1	B	215	PHE	2.9
1	F	204	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	1	MET	2.7
1	F	215	PHE	2.7
1	F	222	GLU	2.7
1	D	215	PHE	2.6
1	G	215	PHE	2.6
1	G	204	GLN	2.6
1	G	222	GLU	2.6
1	G	212[A]	ARG	2.6
1	C	44	GLN	2.5
1	F	209	GLY	2.5
1	H	44	GLN	2.5
1	A	153	GLU	2.4
1	C	203	GLU	2.4
1	E	0	SER	2.4
1	D	205	ASN	2.4
1	D	208	THR	2.3
1	A	222	GLU	2.3
1	H	215	PHE	2.3
1	A	204	GLN	2.3
1	C	216	ARG	2.3
1	G	207	PRO	2.3
1	B	207	PRO	2.2
1	G	44	GLN	2.2
1	F	205	ASN	2.1
1	G	202	ASP	2.1
1	C	204	GLN	2.1
1	H	3[A]	TYR	2.1
1	E	204	GLN	2.1
1	C	266	HIS	2.0
1	A	142	GLN	2.0
1	A	1	MET	2.0
1	C	134	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	OAS	A	93	9/10	0.86	0.16	29,30,34,34	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	OAS	G	93	9/10	0.87	0.16	24,28,33,34	0
1	OAS	E	93	9/10	0.90	0.12	26,30,32,32	0
1	OAS	F	93	9/10	0.90	0.14	27,29,34,36	0
1	OAS	B	93	9/10	0.90	0.14	25,26,30,33	0
1	OAS	D	93	9/10	0.92	0.11	26,27,31,33	0
1	OAS	C	93	9/10	0.92	0.12	23,24,28,30	0
1	OAS	H	93	9/10	0.92	0.12	30,33,35,36	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

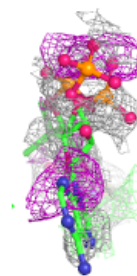
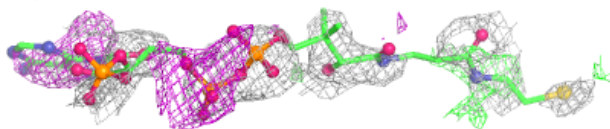
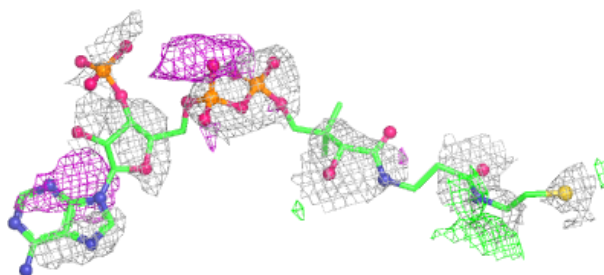
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	B	400	48/48	0.68	0.23	90,115,175,188	0
2	COA	E	401	48/48	0.71	0.20	74,113,176,183	0
2	COA	G	401	48/48	0.71	0.20	79,132,172,184	0
2	COA	H	401	48/48	0.71	0.22	70,148,201,210	0
2	COA	D	400	48/48	0.72	0.21	77,94,138,142	0
2	COA	F	400	48/48	0.72	0.18	63,106,147,159	0
3	GOL	E	402	6/6	0.74	0.23	47,57,60,61	0
2	COA	A	400	48/48	0.76	0.20	88,125,172,182	0
2	COA	C	400	48/48	0.76	0.21	82,112,197,204	0
3	GOL	H	402	6/6	0.77	0.18	70,72,75,78	0
3	GOL	G	402	6/6	0.85	0.12	57,64,67,73	0
3	GOL	G	403	6/6	0.91	0.11	58,66,70,72	0

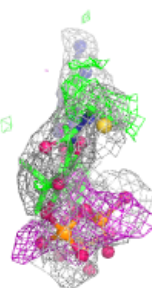
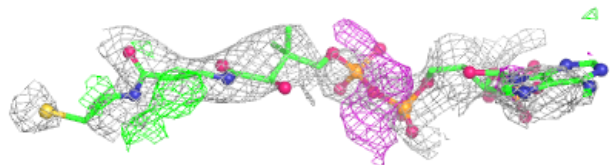
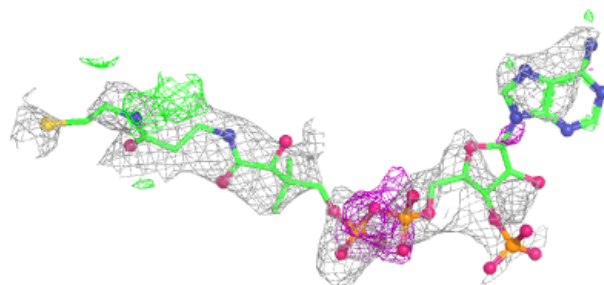
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around COA B 400:

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

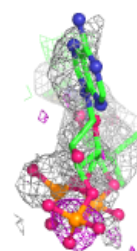
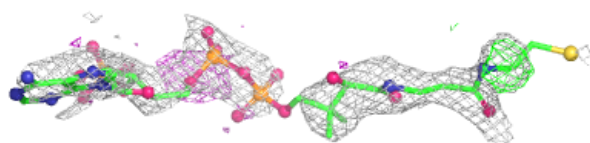
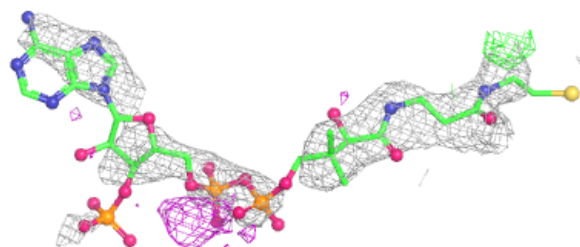
**Electron density around COA E 401:**

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

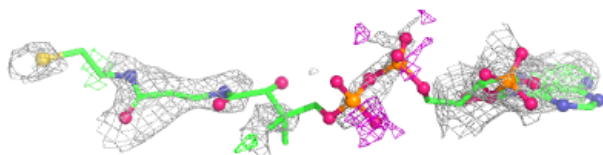
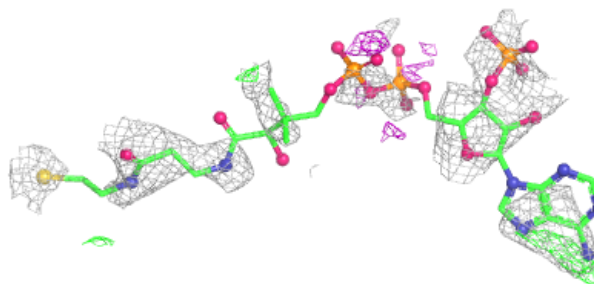


Electron density around COA G 401:

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

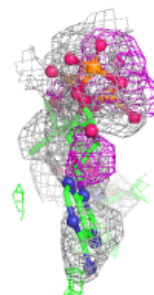
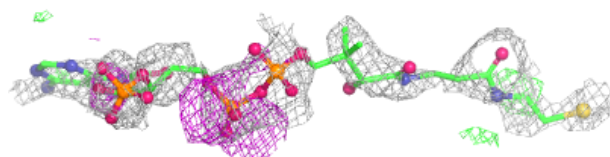
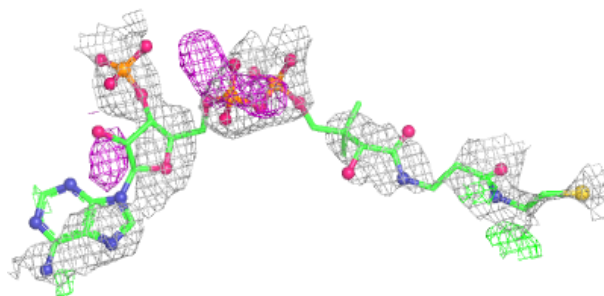
**Electron density around COA H 401:**

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

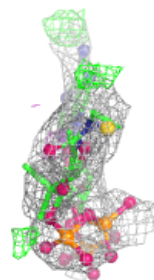
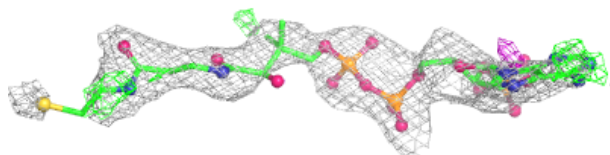
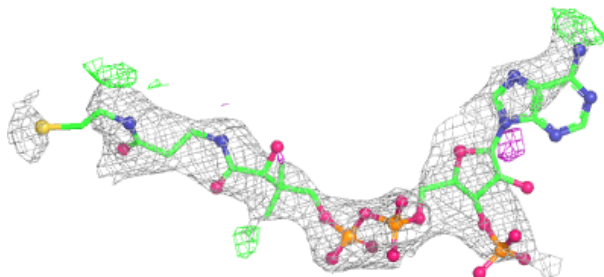


Electron density around COA D 400:

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

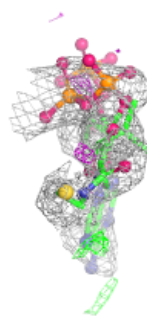
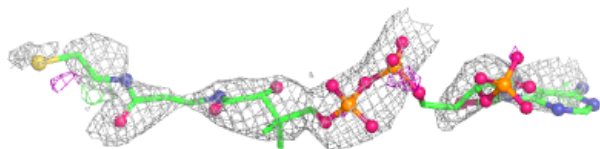
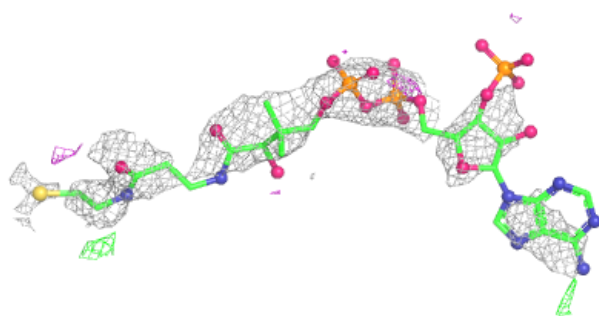
**Electron density around COA F 400:**

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

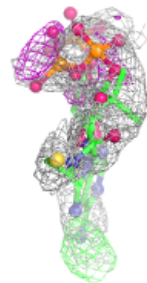
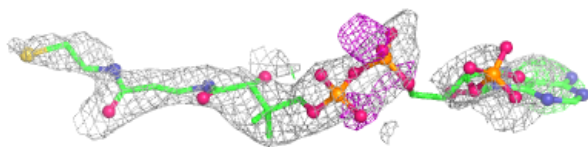
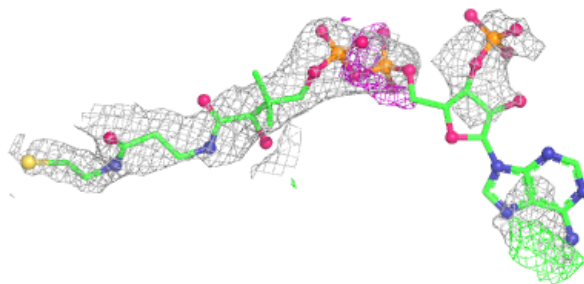


Electron density around COA A 400:

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

**Electron density around COA C 400:**

$2mF_o-DF_c$ (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.