



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:52 AM UTC

PDB ID : 2F2S / pdb_00002f2s
Title : Human mitochondrial acetoacetyl-CoA thiolase
Authors : Min, J.R.; Dombrowski, L.; Antoshenko, T.; Wu, H.; Loppnau, P.; Weigelt, J.; Sundstrom, M.; Arrowsmith, C.H.; Edwards, A.M.; Bochkarev, A.; Plotnikov, A.N.; Structural Genomics Consortium (SGC)
Deposited on : 2005-11-17
Resolution : 2.00 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

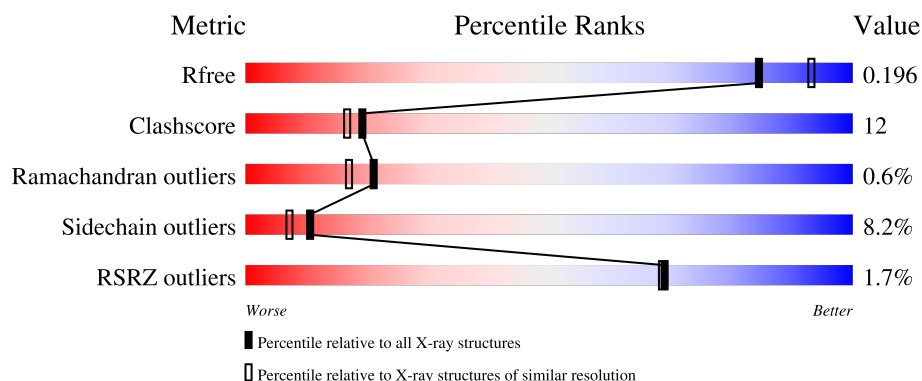
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	406	4% 69% 23% ...
1	B	406	75% 17% 5%
1	C	406	2% 71% 21% ...
1	D	406	% 70% 21% 6%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	SCY	A	126	-	-	X	-
1	SCY	B	126	-	-	X	-
1	SCY	C	126	-	-	X	-
1	SCY	D	126	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12020 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, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	0	0
			2834	1783	486	544	21			
1	B	386	Total	C	N	O	S	0	0	0
			2818	1773	484	540	21			
1	C	388	Total	C	N	O	S	0	0	0
			2841	1789	487	544	21			
1	D	382	Total	C	N	O	S	0	0	0
			2801	1765	478	537	21			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	MET	-	cloning artifact	UNP P24752
A	23	GLY	-	cloning artifact	UNP P24752
A	24	SER	-	cloning artifact	UNP P24752
A	25	SER	-	cloning artifact	UNP P24752
A	26	HIS	-	expression tag	UNP P24752
A	27	HIS	-	expression tag	UNP P24752
A	28	HIS	-	expression tag	UNP P24752
A	29	HIS	-	expression tag	UNP P24752
A	30	HIS	-	expression tag	UNP P24752
A	31	HIS	-	expression tag	UNP P24752
A	32	SER	-	cloning artifact	UNP P24752
A	33	SER	-	cloning artifact	UNP P24752
A	34	GLY	-	cloning artifact	UNP P24752
A	35	LEU	-	cloning artifact	UNP P24752
A	36	VAL	-	cloning artifact	UNP P24752
A	37	PRO	-	cloning artifact	UNP P24752
A	38	ARG	-	cloning artifact	UNP P24752
A	39	GLY	-	cloning artifact	UNP P24752
A	40	SER	-	cloning artifact	UNP P24752
A	126	SCY	CYS	modified residue	UNP P24752
B	22	MET	-	cloning artifact	UNP P24752

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Chain	Residue	Modelled	Actual	Comment	Reference
B	23	GLY	-	cloning artifact	UNP P24752
B	24	SER	-	cloning artifact	UNP P24752
B	25	SER	-	cloning artifact	UNP P24752
B	26	HIS	-	expression tag	UNP P24752
B	27	HIS	-	expression tag	UNP P24752
B	28	HIS	-	expression tag	UNP P24752
B	29	HIS	-	expression tag	UNP P24752
B	30	HIS	-	expression tag	UNP P24752
B	31	HIS	-	expression tag	UNP P24752
B	32	SER	-	cloning artifact	UNP P24752
B	33	SER	-	cloning artifact	UNP P24752
B	34	GLY	-	cloning artifact	UNP P24752
B	35	LEU	-	cloning artifact	UNP P24752
B	36	VAL	-	cloning artifact	UNP P24752
B	37	PRO	-	cloning artifact	UNP P24752
B	38	ARG	-	cloning artifact	UNP P24752
B	39	GLY	-	cloning artifact	UNP P24752
B	40	SER	-	cloning artifact	UNP P24752
B	126	SCY	CYS	modified residue	UNP P24752
C	22	MET	-	cloning artifact	UNP P24752
C	23	GLY	-	cloning artifact	UNP P24752
C	24	SER	-	cloning artifact	UNP P24752
C	25	SER	-	cloning artifact	UNP P24752
C	26	HIS	-	expression tag	UNP P24752
C	27	HIS	-	expression tag	UNP P24752
C	28	HIS	-	expression tag	UNP P24752
C	29	HIS	-	expression tag	UNP P24752
C	30	HIS	-	expression tag	UNP P24752
C	31	HIS	-	expression tag	UNP P24752
C	32	SER	-	cloning artifact	UNP P24752
C	33	SER	-	cloning artifact	UNP P24752
C	34	GLY	-	cloning artifact	UNP P24752
C	35	LEU	-	cloning artifact	UNP P24752
C	36	VAL	-	cloning artifact	UNP P24752
C	37	PRO	-	cloning artifact	UNP P24752
C	38	ARG	-	cloning artifact	UNP P24752
C	39	GLY	-	cloning artifact	UNP P24752
C	40	SER	-	cloning artifact	UNP P24752
C	126	SCY	CYS	modified residue	UNP P24752
D	22	MET	-	cloning artifact	UNP P24752
D	23	GLY	-	cloning artifact	UNP P24752
D	24	SER	-	cloning artifact	UNP P24752

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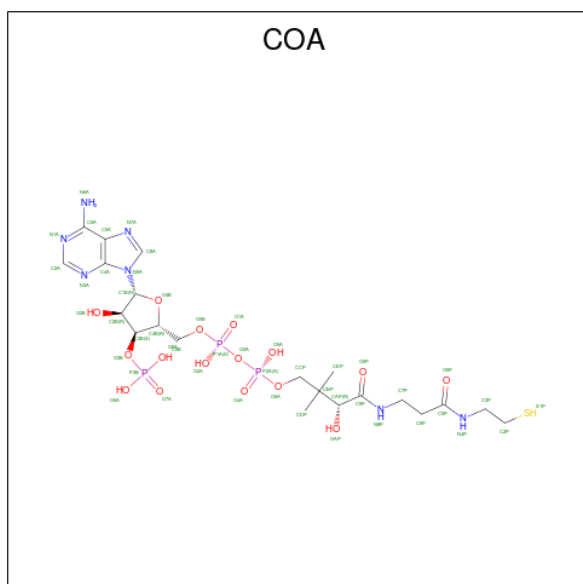
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Chain	Residue	Modelled	Actual	Comment	Reference
D	25	SER	-	cloning artifact	UNP P24752
D	26	HIS	-	expression tag	UNP P24752
D	27	HIS	-	expression tag	UNP P24752
D	28	HIS	-	expression tag	UNP P24752
D	29	HIS	-	expression tag	UNP P24752
D	30	HIS	-	expression tag	UNP P24752
D	31	HIS	-	expression tag	UNP P24752
D	32	SER	-	cloning artifact	UNP P24752
D	33	SER	-	cloning artifact	UNP P24752
D	34	GLY	-	cloning artifact	UNP P24752
D	35	LEU	-	cloning artifact	UNP P24752
D	36	VAL	-	cloning artifact	UNP P24752
D	37	PRO	-	cloning artifact	UNP P24752
D	38	ARG	-	cloning artifact	UNP P24752
D	39	GLY	-	cloning artifact	UNP P24752
D	40	SER	-	cloning artifact	UNP P24752
D	126	SCY	CYS	modified residue	UNP P24752

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0
2	B	2	Total Cl 2 2	0	0

- Molecule 3 is COENZYME A (CCD ID: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S).



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

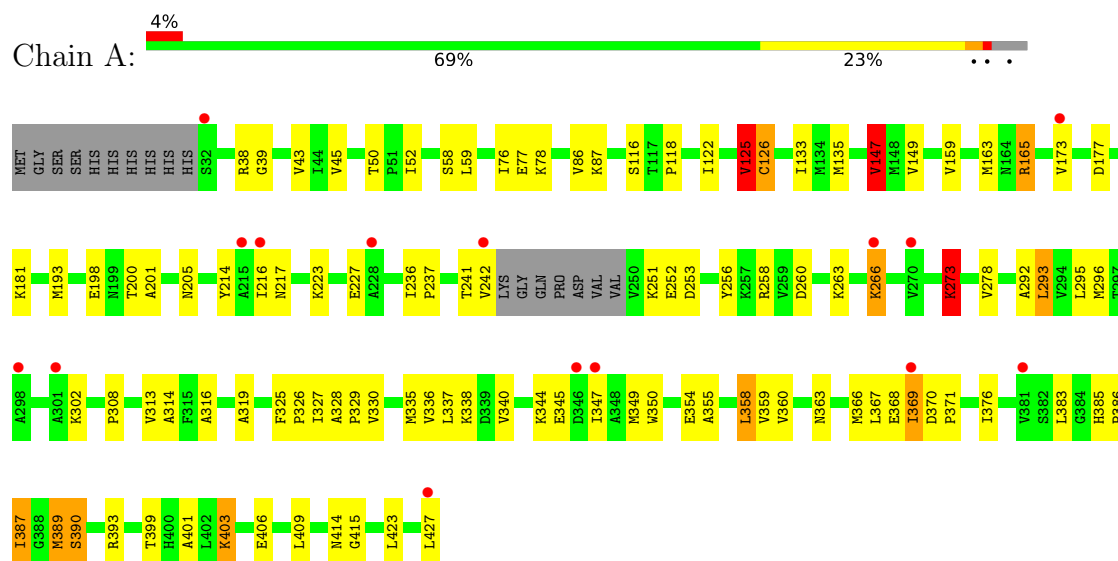
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	99	Total 99	O 99	0	0
4	B	172	Total 172	O 172	0	0
4	C	137	Total 137	O 137	0	0
4	D	123	Total 123	O 123	0	0

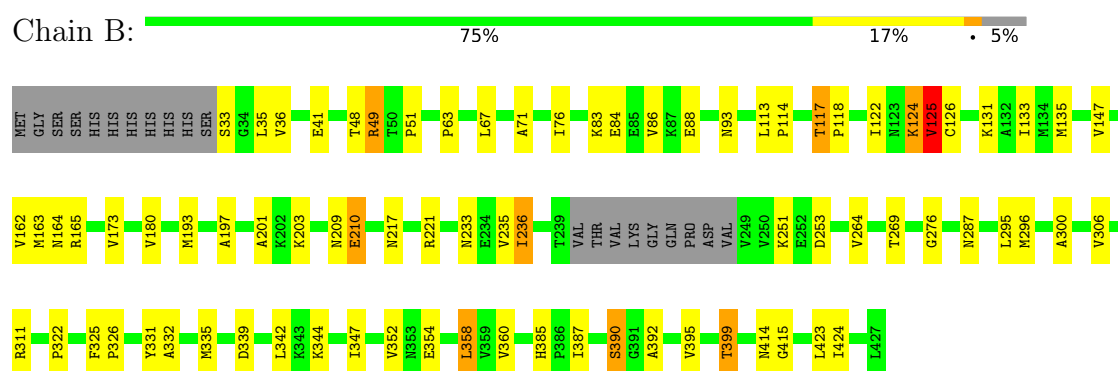
3 Residue-property plots [i](#)

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.

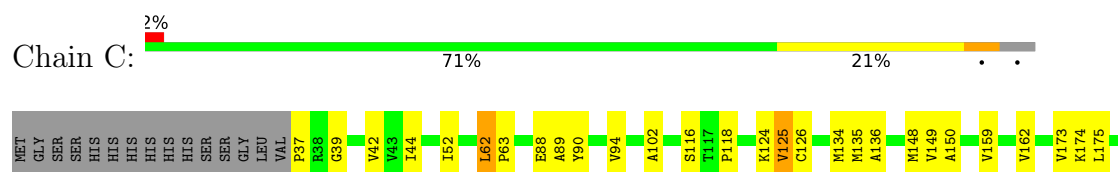
- Molecule 1: Acetyl-CoA acetyltransferase, mitochondrial

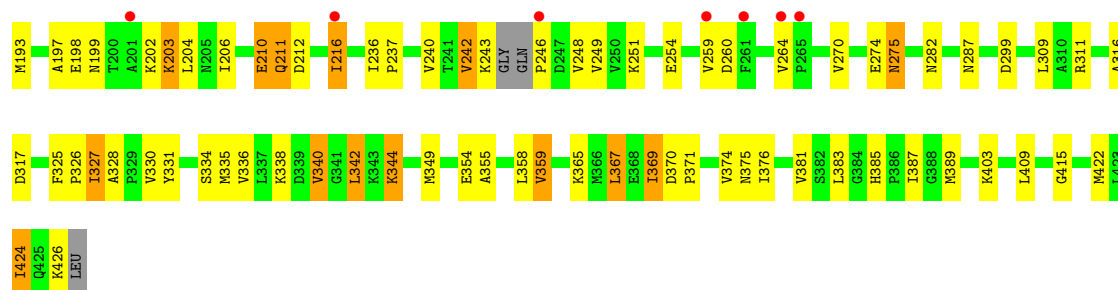


- Molecule 1: Acetyl-CoA acetyltransferase, mitochondrial

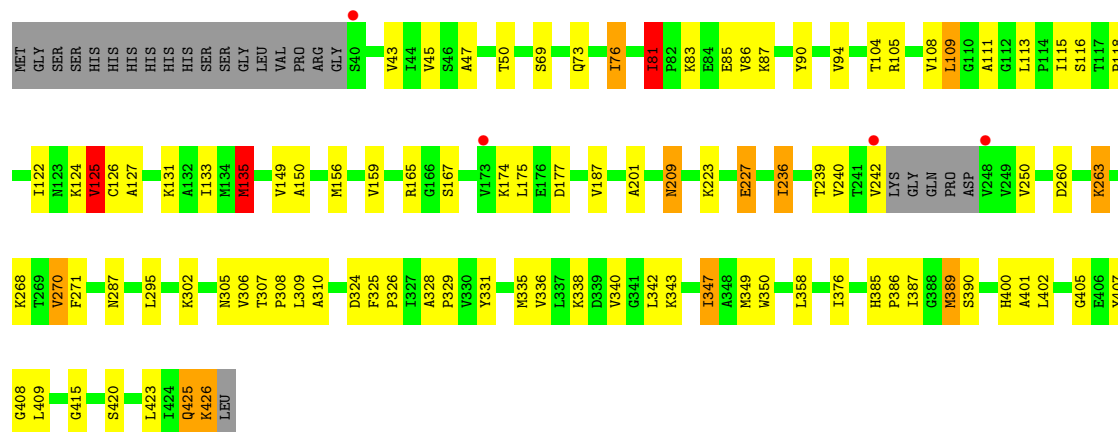


- Molecule 1: Acetyl-CoA acetyltransferase, mitochondrial





- Molecule 1: Acetyl-CoA acetyltransferase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.99Å 126.64Å 111.86Å 90.00° 98.64° 90.00°	Depositor
Resolution (Å)	33.80 – 2.00 33.80 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.9 (33.80-2.00) 98.9 (33.80-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.199 , 0.257 0.202 , 0.196	Depositor DCC
R_{free} test set	5269 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12020	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.50% 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: SCY, 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	1.00	1/2864 (0.0%)	1.18	6/3876 (0.2%)
1	B	1.26	9/2848 (0.3%)	1.25	14/3851 (0.4%)
1	C	1.17	6/2871 (0.2%)	1.22	10/3883 (0.3%)
1	D	1.05	1/2830 (0.0%)	1.18	11/3828 (0.3%)
All	All	1.12	17/11413 (0.1%)	1.21	41/15438 (0.3%)

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	D	0	1
All	All	0	3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	236	ILE	CA-CB	8.92	1.63	1.53
1	B	71	ALA	CA-CB	7.43	1.64	1.53
1	B	264	VAL	CA-CB	7.39	1.57	1.54
1	D	81	ILE	CA-CB	7.31	1.61	1.55
1	A	147	VAL	CA-CB	6.99	1.64	1.54
1	B	347	ILE	CA-CB	5.79	1.60	1.54
1	C	359	VAL	CA-CB	5.68	1.61	1.54
1	C	52	ILE	CA-CB	5.55	1.60	1.54
1	B	387	ILE	CA-CB	5.36	1.60	1.54
1	C	102	ALA	CA-C	5.34	1.57	1.53
1	C	102	ALA	C-O	-5.34	1.21	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	352	VAL	CA-CB	5.25	1.60	1.53
1	C	42	VAL	CA-CB	5.13	1.60	1.54
1	B	392	ALA	CA-CB	-5.13	1.45	1.53
1	C	136	ALA	CA-CB	5.10	1.61	1.53
1	B	88	GLU	CA-C	5.00	1.58	1.52
1	B	332	ALA	CA-CB	5.00	1.61	1.53

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	246	PRO	N-CA-CB	7.07	110.78	103.00
1	B	113	LEU	CA-C-N	-7.01	112.54	119.76
1	B	113	LEU	C-N-CA	-7.01	112.54	119.76
1	A	58	SER	N-CA-C	6.87	119.61	111.71
1	A	50	THR	CA-C-N	-6.83	113.27	120.03
1	A	50	THR	C-N-CA	-6.83	113.27	120.03
1	B	49	ARG	NE-CZ-NH2	-6.63	113.23	119.20
1	D	307	THR	CA-C-N	6.47	126.43	119.76
1	D	307	THR	C-N-CA	6.47	126.43	119.76
1	C	264	VAL	CA-C-N	-6.40	114.30	120.83
1	C	264	VAL	C-N-CA	-6.40	114.30	120.83
1	B	387	ILE	N-CA-C	6.39	117.07	110.36
1	D	135	MET	N-CA-C	6.38	118.24	111.28
1	B	390	SER	N-CA-C	6.29	118.94	111.33
1	A	147	VAL	N-CA-C	6.28	116.63	108.35
1	B	180	VAL	N-CA-C	6.16	116.33	110.42
1	A	165	ARG	NE-CZ-NH2	-6.15	113.66	119.20
1	D	236	ILE	CA-C-N	6.06	126.02	119.78
1	D	236	ILE	C-N-CA	6.06	126.02	119.78
1	B	117	THR	CA-C-N	5.72	125.65	119.76
1	B	117	THR	C-N-CA	5.72	125.65	119.76
1	D	187	VAL	N-CA-C	5.70	116.43	110.62
1	C	325	PHE	CA-C-N	-5.66	114.28	119.82
1	C	325	PHE	C-N-CA	-5.66	114.28	119.82
1	C	211	GLN	CA-C-N	-5.62	113.26	122.54
1	C	211	GLN	C-N-CA	-5.62	113.26	122.54
1	B	49	ARG	CD-NE-CZ	5.61	132.26	124.40
1	C	328	ALA	CA-C-N	-5.56	113.89	119.56
1	C	328	ALA	C-N-CA	-5.56	113.89	119.56
1	D	50	THR	CA-C-N	-5.56	113.97	119.92
1	D	50	THR	C-N-CA	-5.56	113.97	119.92
1	A	390	SER	N-CA-C	5.53	116.99	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	287	ASN	N-CA-C	5.49	116.60	108.86
1	B	235	VAL	N-CA-C	5.34	116.40	108.23
1	B	124	LYS	N-CA-C	-5.31	104.06	110.44
1	B	49	ARG	NE-CZ-NH1	5.24	126.74	121.50
1	B	358	LEU	N-CA-C	5.24	117.67	111.33
1	D	425	GLN	CA-C-N	-5.17	112.39	121.70
1	D	425	GLN	C-N-CA	-5.17	112.39	121.70
1	D	81	ILE	N-CA-CB	5.14	115.38	111.83
1	C	355	ALA	N-CA-C	-5.08	105.74	111.28

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	386	PRO	Peptide
1	B	125	VAL	Mainchain
1	D	425	GLN	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	2834	0	2887	70	0
1	B	2818	0	2882	63	0
1	C	2841	0	2910	65	0
1	D	2801	0	2869	82	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
3	A	48	0	32	2	0
3	B	48	0	32	5	0
3	C	48	0	32	2	0
3	D	48	0	32	0	0
4	A	99	0	0	0	0
4	B	172	0	0	7	0
4	C	137	0	0	5	0
4	D	123	0	0	6	0
All	All	12020	0	11676	267	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 12.

All (267) 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:125:VAL:C	1:A:126:SCY:H2	1.21	1.49
1:B:131:LYS:HG3	1:B:135:MET:CE	1.86	1.06
1:B:131:LYS:CG	1:B:135:MET:HE2	1.87	1.04
1:D:426:LYS:O	1:D:426:LYS:HD3	1.58	1.03
1:D:122:ILE:HG13	1:D:135:MET:HE1	1.41	1.01
1:D:309:LEU:O	1:D:426:LYS:HG2	1.66	0.96
1:A:38:ARG:HB3	1:B:36:VAL:HG13	1.46	0.94
1:A:133:ILE:HG23	1:A:295:LEU:HD11	1.49	0.93
1:C:94:VAL:HG21	1:C:387:ILE:HG22	1.51	0.92
1:B:122:ILE:CG2	1:B:135:MET:HE1	2.03	0.89
1:D:309:LEU:O	1:D:426:LYS:CG	2.22	0.87
1:D:126:SCY:HE1	1:D:415:GLY:H	1.45	0.82
1:C:202:LYS:O	1:C:203:LYS:HB2	1.77	0.81
1:D:223:LYS:O	1:D:227:GLU:HG2	1.81	0.81
1:A:125:VAL:C	1:A:126:SCY:H	1.86	0.80
1:D:125:VAL:CG1	1:D:126:SCY:HE2	2.13	0.79
1:D:347:ILE:HD11	1:D:350:TRP:CZ2	2.18	0.79
1:B:122:ILE:HG23	1:B:135:MET:HE1	1.65	0.78
1:D:125:VAL:HG12	1:D:126:SCY:HE2	1.64	0.78
1:C:126:SCY:CE	1:C:415:GLY:H	1.97	0.77
1:B:131:LYS:HG3	1:B:135:MET:HE2	0.91	0.77
1:B:126:SCY:HE1	1:B:415:GLY:H	1.48	0.77
1:C:126:SCY:HE1	1:C:415:GLY:H	1.49	0.76
1:A:201:ALA:HB2	1:A:358:LEU:CD1	2.18	0.74
1:B:48:THR:HG22	4:B:2104:HOH:O	1.86	0.74
1:A:252:GLU:OE2	1:A:256:TYR:OH	2.05	0.74
1:B:197:ALA:O	1:B:358:LEU:HD21	1.89	0.73
1:B:122:ILE:HG21	1:B:135:MET:HE1	1.69	0.73
1:C:274:GLU:HG2	1:C:275:ASN:ND2	2.03	0.73
1:D:126:SCY:C	1:D:127:ALA:CA	2.67	0.73
1:A:125:VAL:HG22	1:A:126:SCY:CE	2.19	0.72
1:A:223:LYS:O	1:A:227:GLU:HG2	1.90	0.72
1:D:426:LYS:O	1:D:426:LYS:CD	2.37	0.71
1:D:122:ILE:HG13	1:D:135:MET:CE	2.19	0.70
1:D:125:VAL:HG12	1:D:126:SCY:CE	2.21	0.70
1:B:126:SCY:HE1	1:B:415:GLY:N	2.06	0.70
1:C:116:SER:O	1:C:118:PRO:HD3	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:197:ALA:O	1:C:358:LEU:HD11	1.90	0.70
1:A:201:ALA:HB2	1:A:358:LEU:HD13	1.74	0.69
1:D:125:VAL:CG1	1:D:126:SCY:CE	2.69	0.69
1:D:131:LYS:O	1:D:135:MET:HE2	1.91	0.69
1:B:126:SCY:CE	1:B:415:GLY:H	2.05	0.69
1:D:385:HIS:HD2	1:D:390:SER:OG	1.74	0.69
1:B:49:ARG:NH2	1:B:253:ASP:OD1	2.25	0.68
1:A:126:SCY:HE1	1:A:415:GLY:H	1.59	0.67
1:C:212:ASP:O	1:C:216:ILE:HG23	1.93	0.67
1:B:67:LEU:HD12	1:B:67:LEU:H	1.59	0.66
1:B:124:LYS:NZ	1:B:414:ASN:HD21	1.94	0.66
1:C:210:GLU:HB3	4:C:1137:HOH:O	1.95	0.66
1:B:124:LYS:HZ3	1:B:414:ASN:HD21	1.44	0.66
1:D:209:ASN:H	1:D:209:ASN:HD22	1.41	0.66
1:C:274:GLU:HG2	1:C:275:ASN:HD22	1.60	0.66
1:B:122:ILE:HG21	1:B:135:MET:CE	2.26	0.65
1:B:197:ALA:HB1	1:B:358:LEU:HD23	1.77	0.65
1:A:125:VAL:HG22	1:A:126:SCY:HE2	1.77	0.65
1:A:163:MET:HE1	1:D:175:LEU:HD21	1.79	0.65
1:C:259:VAL:HG22	1:C:260:ASP:H	1.60	0.65
1:B:126:SCY:HE2	1:B:126:SCY:H	1.62	0.64
1:C:287:ASN:ND2	1:C:385:HIS:H	1.96	0.64
1:D:94:VAL:HG21	1:D:387:ILE:HG22	1.79	0.64
1:B:233:ASN:HB2	4:B:2062:HOH:O	1.97	0.64
1:D:47:ALA:O	1:D:400:HIS:HE1	1.79	0.64
1:B:164:ASN:ND2	4:B:2149:HOH:O	2.29	0.63
1:C:126:SCY:HE1	1:C:415:GLY:N	2.12	0.63
1:C:344:LYS:HB2	1:C:369:ILE:HD11	1.80	0.63
1:C:193:MET:HE1	1:C:415:GLY:HA2	1.80	0.63
1:C:330:VAL:HG22	1:C:367:LEU:CD2	2.29	0.62
1:D:385:HIS:CD2	1:D:390:SER:OG	2.53	0.62
1:D:309:LEU:O	1:D:426:LYS:HG3	1.99	0.61
1:D:76:ILE:HD13	1:D:113:LEU:HD11	1.81	0.61
1:D:86:VAL:HG11	1:D:149:VAL:HG12	1.81	0.61
1:D:122:ILE:CG1	1:D:135:MET:HE1	2.25	0.61
1:A:38:ARG:HB3	1:B:36:VAL:CG1	2.28	0.61
1:A:163:MET:CE	1:D:175:LEU:HD21	2.31	0.61
1:C:354:GLU:HB3	1:C:381:VAL:HG23	1.82	0.60
3:B:1002:COA:H52A	4:B:2111:HOH:O	2.00	0.60
1:C:199:ASN:HB3	4:C:1065:HOH:O	2.01	0.60
1:D:402:LEU:HD11	1:D:408:GLY:HA3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:LYS:HE2	1:C:317:ASP:OD1	2.02	0.60
1:D:126:SCY:HE2	1:D:126:SCY:H	1.65	0.60
1:D:349:MET:HE3	4:D:1045:HOH:O	2.01	0.59
1:C:39:GLY:HA2	1:C:299:ASP:OD2	2.00	0.59
1:B:395:VAL:O	1:B:399:THR:HG23	2.03	0.59
1:D:426:LYS:CB	4:D:1126:HOH:O	2.51	0.59
1:D:125:VAL:CB	1:D:126:SCY:HE2	2.34	0.58
1:A:330:VAL:HG21	1:A:366:MET:HE2	1.86	0.58
1:A:314:ALA:HB3	1:A:336:VAL:HG13	1.85	0.57
1:C:358:LEU:C	1:C:358:LEU:HD23	2.28	0.57
1:D:405:GLY:H	1:D:426:LYS:C	2.12	0.57
1:A:253:ASP:OD2	1:A:383:LEU:HD22	2.04	0.57
1:A:403:LYS:N	1:A:406:GLU:OE1	2.33	0.57
1:A:86:VAL:HG11	1:A:149:VAL:HG23	1.86	0.57
1:B:114:PRO:HG2	1:B:117:THR:HG23	1.87	0.57
1:C:204:LEU:HD12	1:C:365:LYS:HD2	1.86	0.56
1:A:349:MET:HE1	1:A:401:ALA:HB1	1.88	0.56
1:D:81:ILE:HG22	1:D:306:VAL:HG21	1.87	0.56
1:D:328:ALA:N	1:D:329:PRO:CD	2.69	0.56
1:A:337:LEU:HD21	1:A:347:ILE:HD11	1.87	0.56
1:B:217:ASN:O	1:B:221:ARG:HG3	2.04	0.56
1:C:426:LYS:C	4:C:1128:HOH:O	2.47	0.56
1:C:44:ILE:HD12	1:C:424:ILE:HD12	1.86	0.56
1:C:259:VAL:HG22	1:C:260:ASP:N	2.21	0.56
1:C:259:VAL:HG23	3:C:1003:COA:H61A	1.71	0.55
1:C:287:ASN:HD22	1:C:385:HIS:H	1.54	0.55
1:A:198:GLU:OE2	1:A:278:VAL:HG12	2.06	0.55
1:C:309:LEU:O	1:C:426:LYS:HD2	2.06	0.55
1:A:147:VAL:HB	1:A:296:MET:HG3	1.89	0.55
1:C:326:PRO:HB3	1:C:359:VAL:HG22	1.89	0.55
1:B:83:LYS:HE2	4:B:2088:HOH:O	2.07	0.54
1:A:126:SCY:CE	1:A:415:GLY:H	2.19	0.54
1:A:326:PRO:HB3	1:A:359:VAL:HG22	1.89	0.54
1:C:94:VAL:HG21	1:C:387:ILE:CG2	2.32	0.54
1:D:45:VAL:CG1	1:D:306:VAL:HG11	2.38	0.53
1:A:389:MET:HE3	1:A:390:SER:HA	1.88	0.53
1:C:126:SCY:HE2	1:C:126:SCY:H	1.73	0.53
1:D:69:SER:O	1:D:73:GLN:HG3	2.07	0.53
1:A:214:TYR:O	1:A:217:ASN:HB3	2.09	0.53
1:A:86:VAL:CG1	1:A:149:VAL:HG23	2.39	0.52
1:A:329:PRO:HG2	1:A:363:ASN:HD21	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:LEU:HD12	1:B:67:LEU:N	2.23	0.52
1:C:89:ALA:CB	1:C:149:VAL:HG23	2.40	0.52
1:C:249:VAL:HG12	1:C:251:LYS:HG2	1.91	0.52
1:B:125:VAL:HG12	1:B:126:SCY:CE	2.39	0.51
1:C:126:SCY:OCD	3:C:1003:COA:S1P	2.67	0.51
1:A:349:MET:HE1	1:A:401:ALA:CB	2.40	0.51
1:D:125:VAL:HB	1:D:126:SCY:HE2	1.91	0.51
1:D:260:ASP:OD2	1:D:263:LYS:HB2	2.10	0.51
1:A:260:ASP:CG	1:A:263:LYS:HG3	2.35	0.51
1:D:94:VAL:HG21	1:D:387:ILE:CG2	2.41	0.51
1:C:212:ASP:HB3	1:C:282:ASN:HB3	1.93	0.51
1:B:342:LEU:HD12	1:B:423:LEU:HD21	1.93	0.50
1:B:51:PRO:HD3	1:B:236:ILE:O	2.12	0.50
1:A:125:VAL:HG22	1:A:126:SCY:HE3	1.93	0.50
1:C:197:ALA:O	1:C:358:LEU:CD1	2.58	0.50
1:C:44:ILE:CD1	1:C:424:ILE:HD12	2.41	0.50
1:C:349:MET:HG2	1:C:376:ILE:HD12	1.92	0.50
1:A:126:SCY:OCD	3:A:1001:COA:S1P	2.69	0.49
1:B:41:GLU:OE2	1:B:311:ARG:NH2	2.37	0.49
1:D:386:PRO:HD2	1:D:389:MET:HE2	1.94	0.49
1:A:39:GLY:O	1:B:36:VAL:HG11	2.12	0.49
1:C:331:TYR:O	1:C:335:MET:HG3	2.13	0.49
1:C:316:ALA:HB1	1:C:335:MET:HE2	1.94	0.49
1:A:87:LYS:O	1:A:118:PRO:HD2	2.12	0.49
1:A:116:SER:O	1:A:118:PRO:HD3	2.12	0.49
1:D:209:ASN:H	1:D:209:ASN:ND2	2.09	0.49
1:D:45:VAL:HG11	1:D:306:VAL:HG11	1.94	0.49
1:A:126:SCY:SG	1:A:387:ILE:HG12	2.53	0.48
1:A:358:LEU:O	1:A:358:LEU:HG	2.13	0.48
1:B:210:GLU:HG3	4:B:2051:HOH:O	2.13	0.48
1:D:325:PHE:N	1:D:326:PRO:CD	2.76	0.48
1:B:385:HIS:HD2	1:B:390:SER:OG	1.96	0.48
1:A:135:MET:HE1	1:D:122:ILE:HD11	1.96	0.48
1:B:193:MET:HE3	3:B:1002:COA:S1P	2.53	0.48
1:D:126:SCY:HA	1:D:387:ILE:HG23	1.95	0.48
1:B:296:MET:HG2	1:B:300:ALA:HB3	1.96	0.48
1:A:313:VAL:HG11	1:A:423:LEU:HD23	1.95	0.48
1:D:43:VAL:HB	1:D:308:PRO:HB3	1.96	0.48
1:D:124:LYS:HE2	1:D:420:SER:OG	2.14	0.48
1:D:358:LEU:O	1:D:358:LEU:HG	2.04	0.47
1:B:399:THR:HG22	1:B:424:ILE:HD13	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:197:ALA:HB1	1:C:358:LEU:HD13	1.96	0.47
1:A:52:ILE:HD11	1:A:393:ARG:HD3	1.96	0.47
1:A:177:ASP:OD2	1:D:165:ARG:NH2	2.47	0.47
1:A:266:LYS:HB2	1:A:266:LYS:HE3	1.52	0.47
1:B:354:GLU:HG2	1:B:360:VAL:HG21	1.97	0.47
1:C:37:PRO:HG3	1:C:311:ARG:HD2	1.96	0.47
1:A:43:VAL:HB	1:A:308:PRO:HB3	1.96	0.47
1:B:165:ARG:HD2	1:C:159:VAL:O	2.14	0.47
1:C:242:VAL:O	1:C:243:LYS:HB2	2.15	0.47
1:A:354:GLU:HG2	1:A:360:VAL:HG21	1.97	0.46
1:B:114:PRO:HG2	1:B:117:THR:CG2	2.45	0.46
1:A:355:ALA:HB1	3:A:1001:COA:C2P	2.45	0.46
1:D:76:ILE:CD1	1:D:113:LEU:HD11	2.46	0.46
1:D:310:ALA:HA	1:D:426:LYS:HG2	1.97	0.46
1:D:347:ILE:HG22	1:D:407:TYR:HB3	1.98	0.46
1:B:122:ILE:HD11	1:C:135:MET:HE1	1.97	0.46
1:B:63:PRO:O	1:B:67:LEU:CD1	2.64	0.46
1:C:243:LYS:N	4:C:1104:HOH:O	2.48	0.46
1:A:316:ALA:HB1	1:A:335:MET:HE2	1.98	0.45
1:B:125:VAL:CG1	1:B:126:SCY:CE	2.94	0.45
1:B:269:THR:HG21	1:B:276:GLY:C	2.41	0.45
1:C:126:SCY:SG	1:C:387:ILE:HD12	2.56	0.45
1:B:193:MET:HG3	3:B:1002:COA:H22	1.98	0.45
1:D:201:ALA:HB2	1:D:358:LEU:HD13	1.97	0.45
1:C:126:SCY:HE2	1:C:415:GLY:H	1.79	0.45
1:B:124:LYS:NZ	1:B:414:ASN:ND2	2.63	0.45
1:A:292:ALA:C	1:A:293:LEU:HD23	2.42	0.45
1:B:162:VAL:HG12	1:C:162:VAL:HG12	1.98	0.45
1:A:328:ALA:N	1:A:329:PRO:HD2	2.32	0.45
1:B:126:SCY:OCD	3:B:1002:COA:S1P	2.70	0.45
1:B:125:VAL:HG12	1:B:126:SCY:HE3	1.98	0.45
1:C:242:VAL:HG12	1:C:243:LYS:N	2.32	0.44
1:D:116:SER:O	1:D:118:PRO:HD3	2.16	0.44
1:A:177:ASP:CG	1:D:165:ARG:HH22	2.24	0.44
1:B:331:TYR:O	1:B:335:MET:HE3	2.18	0.44
1:C:88:GLU:HG2	1:C:148:MET:HE3	1.99	0.44
1:D:133:ILE:HG23	1:D:295:LEU:HD11	1.99	0.44
1:D:331:TYR:O	1:D:335:MET:HG3	2.18	0.44
1:A:313:VAL:CG1	1:A:423:LEU:HD23	2.48	0.44
1:B:84:GLU:HA	1:B:114:PRO:HG3	1.99	0.44
1:B:193:MET:CE	3:B:1002:COA:S1P	3.06	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:426:LYS:HB2	4:D:1126:HOH:O	2.15	0.44
1:A:76:ILE:CD1	1:A:86:VAL:HG21	2.48	0.43
1:A:159:VAL:O	1:D:165:ARG:NH1	2.50	0.43
1:A:125:VAL:HG13	1:A:414:ASN:HB2	1.99	0.43
1:D:306:VAL:HG13	4:D:1112:HOH:O	2.18	0.43
1:D:83:LYS:HG2	1:D:113:LEU:HD12	1.98	0.43
1:C:89:ALA:HB2	1:C:149:VAL:HG23	1.99	0.43
1:A:193:MET:HE2	1:A:193:MET:HA	2.00	0.43
1:B:133:ILE:HG23	1:B:295:LEU:HD11	1.99	0.43
1:D:270:VAL:HG23	1:D:271:PHE:CD2	2.54	0.43
1:D:268:LYS:HG2	4:D:1099:HOH:O	2.17	0.43
1:B:33:SER:N	4:B:2076:HOH:O	2.51	0.43
1:C:340:VAL:HG12	1:C:342:LEU:HB2	2.00	0.43
1:C:327:ILE:O	1:C:330:VAL:HB	2.19	0.43
1:B:49:ARG:HH22	1:B:253:ASP:CG	2.24	0.42
1:B:325:PHE:N	1:B:326:PRO:CD	2.82	0.42
1:A:198:GLU:CD	1:A:278:VAL:HG12	2.44	0.42
1:D:105:ARG:HG3	1:D:109:LEU:HD22	2.01	0.42
1:D:126:SCY:CA	1:D:127:ALA:N	2.71	0.42
1:C:330:VAL:HG22	1:C:367:LEU:HD22	2.01	0.42
1:B:201:ALA:HB2	1:B:358:LEU:CD2	2.49	0.42
1:C:134:MET:HG3	1:C:422:MET:CE	2.49	0.42
1:D:287:ASN:ND2	1:D:385:HIS:H	2.18	0.42
1:A:214:TYR:CD2	1:A:214:TYR:C	2.98	0.41
1:A:385:HIS:HD2	1:A:390:SER:OG	2.03	0.41
1:D:73:GLN:HG2	1:D:111:ALA:HB1	2.01	0.41
1:A:214:TYR:OH	1:A:360:VAL:HG11	2.20	0.41
1:C:275:ASN:HD22	1:C:275:ASN:N	2.17	0.41
1:D:86:VAL:CG1	1:D:149:VAL:HG12	2.48	0.41
1:A:126:SCY:HE1	1:A:415:GLY:N	2.31	0.41
1:B:147:VAL:HG22	1:B:296:MET:HG3	2.02	0.41
1:A:325:PHE:N	1:A:326:PRO:CD	2.83	0.41
1:A:370:ASP:HA	1:A:371:PRO:HD2	1.93	0.41
1:D:45:VAL:HG13	1:D:306:VAL:CG1	2.50	0.41
1:A:273:LYS:H	1:A:273:LYS:HD2	1.86	0.41
1:B:125:VAL:HG12	1:B:126:SCY:HE2	2.01	0.41
1:C:236:ILE:HA	1:C:237:PRO:HD3	1.94	0.41
1:C:374:VAL:O	1:C:375:ASN:C	2.62	0.41
1:A:319:ALA:O	1:D:116:SER:HA	2.21	0.41
1:B:117:THR:HA	1:B:118:PRO:HD3	1.84	0.41
1:D:90:TYR:O	1:D:150:ALA:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:340:VAL:HG23	1:D:342:LEU:HG	2.02	0.41
1:D:104:THR:O	1:D:108:VAL:HG23	2.20	0.41
1:D:156:MET:HA	1:D:159:VAL:HG23	2.02	0.41
1:D:349:MET:HE1	1:D:401:ALA:O	2.20	0.41
1:A:237:PRO:HG3	1:A:251:LYS:HD3	2.03	0.41
1:B:76:ILE:HD11	1:B:86:VAL:HG21	2.03	0.41
1:C:62:LEU:HA	1:C:63:PRO:HD2	1.87	0.41
1:C:90:TYR:O	1:C:150:ALA:HA	2.21	0.41
1:C:370:ASP:HA	1:C:371:PRO:HD3	1.82	0.41
1:D:324:ASP:C	1:D:326:PRO:HD2	2.45	0.41
1:D:328:ALA:N	1:D:329:PRO:HD3	2.36	0.41
1:A:337:LEU:O	1:A:338:LYS:C	2.64	0.40
1:A:350:TRP:CZ2	1:A:369:ILE:HG21	2.55	0.40
1:B:67:LEU:H	1:B:67:LEU:CD1	2.32	0.40
1:B:163:MET:HE2	1:B:163:MET:HB3	1.89	0.40
1:C:210:GLU:CB	4:C:1137:HOH:O	2.64	0.40
1:D:426:LYS:HB3	4:D:1126:HOH:O	2.19	0.40
1:A:165:ARG:HH22	1:D:177:ASP:CG	2.28	0.40
1:A:251:LYS:HD2	1:A:251:LYS:O	2.20	0.40
1:A:345:GLU:CD	1:A:345:GLU:H	2.28	0.40
1:C:89:ALA:HB2	1:C:149:VAL:CG2	2.51	0.40
1:C:206:ILE:HG22	1:C:211:GLN:HG3	2.04	0.40
1:D:85:GLU:O	1:D:87:LYS:HD3	2.22	0.40
1:D:336:VAL:HG11	1:D:423:LEU:HB2	2.04	0.40
1:D:347:ILE:HG23	1:D:347:ILE:HD13	1.79	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	384/406 (95%)	370 (96%)	11 (3%)	3 (1%)	16 11

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	381/406 (94%)	369 (97%)	10 (3%)	2 (0%)	24	21
1	C	383/406 (94%)	362 (94%)	18 (5%)	3 (1%)	16	11
1	D	377/406 (93%)	367 (97%)	9 (2%)	1 (0%)	36	35
All	All	1525/1624 (94%)	1468 (96%)	48 (3%)	9 (1%)	21	17

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	125	VAL
1	A	340	VAL
1	C	203	LYS
1	A	273	LYS
1	C	242	VAL
1	D	125	VAL
1	B	125	VAL
1	C	125	VAL
1	B	93	ASN

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	296/317 (93%)	263 (89%)	33 (11%)	6	3
1	B	296/317 (93%)	285 (96%)	11 (4%)	30	30
1	C	299/317 (94%)	272 (91%)	27 (9%)	9	6
1	D	295/317 (93%)	269 (91%)	26 (9%)	9	6
All	All	1186/1268 (94%)	1089 (92%)	97 (8%)	10	7

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	VAL
1	A	59	LEU

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Mol	Chain	Res	Type
1	A	77	GLU
1	A	78	LYS
1	A	122	ILE
1	A	125	VAL
1	A	147	VAL
1	A	173	VAL
1	A	181	LYS
1	A	200	THR
1	A	205	ASN
1	A	216	ILE
1	A	236	ILE
1	A	241	THR
1	A	242	VAL
1	A	258	ARG
1	A	266	LYS
1	A	273	LYS
1	A	293	LEU
1	A	302	LYS
1	A	327	ILE
1	A	344	LYS
1	A	358	LEU
1	A	367	LEU
1	A	368	GLU
1	A	369	ILE
1	A	376	ILE
1	A	387	ILE
1	A	389	MET
1	A	399	THR
1	A	403	LYS
1	A	409	LEU
1	A	427	LEU
1	B	35	LEU
1	B	173	VAL
1	B	203	LYS
1	B	209	ASN
1	B	210	GLU
1	B	251	LYS
1	B	306	VAL
1	B	322	PRO
1	B	339	ASP
1	B	344	LYS
1	B	399	THR

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Mol	Chain	Res	Type
1	C	62	LEU
1	C	125	VAL
1	C	173	VAL
1	C	174	LYS
1	C	175	LEU
1	C	198	GLU
1	C	210	GLU
1	C	216	ILE
1	C	240	VAL
1	C	248	VAL
1	C	254	GLU
1	C	270	VAL
1	C	275	ASN
1	C	327	ILE
1	C	334	SER
1	C	336	VAL
1	C	338	LYS
1	C	340	VAL
1	C	342	LEU
1	C	344	LYS
1	C	367	LEU
1	C	369	ILE
1	C	383	LEU
1	C	389	MET
1	C	403	LYS
1	C	409	LEU
1	C	424	ILE
1	D	76	ILE
1	D	81	ILE
1	D	109	LEU
1	D	115	ILE
1	D	125	VAL
1	D	135	MET
1	D	167	SER
1	D	174	LYS
1	D	209	ASN
1	D	227	GLU
1	D	236	ILE
1	D	239	THR
1	D	240	VAL
1	D	242	VAL
1	D	250	VAL

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Mol	Chain	Res	Type
1	D	263	LYS
1	D	270	VAL
1	D	302	LYS
1	D	305	ASN
1	D	338	LYS
1	D	343	LYS
1	D	347	ILE
1	D	376	ILE
1	D	389	MET
1	D	409	LEU
1	D	426	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	189	ASN
1	A	287	ASN
1	A	363	ASN
1	A	377	ASN
1	A	385	HIS
1	B	73	GLN
1	B	211	GLN
1	B	385	HIS
1	B	414	ASN
1	C	73	GLN
1	C	158	ASN
1	C	217	ASN
1	C	275	ASN
1	C	287	ASN
1	C	305	ASN
1	C	377	ASN
1	C	425	GLN
1	D	209	ASN
1	D	272	GLN
1	D	287	ASN
1	D	385	HIS
1	D	400	HIS
1	D	414	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	SCY	A	126	1	7,8,9	1.26	1 (14%)	4,9,11	0.41	0
1	SCY	C	126	1	7,8,9	0.97	0	4,9,11	1.17	0
1	SCY	B	126	1	7,8,9	0.99	0	4,9,11	0.82	0
1	SCY	D	126	1	7,8,9	0.92	0	4,9,11	0.77	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
1	SCY	A	126	1	-	3/5/7/9	-
1	SCY	C	126	1	-	2/5/7/9	-
1	SCY	B	126	1	-	3/5/7/9	-
1	SCY	D	126	1	-	2/5/7/9	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	SCY	CB-SG	-2.35	1.76	1.81

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	126	SCY	OCD-CD-SG-CB
1	A	126	SCY	CE-CD-SG-CB
1	B	126	SCY	OCD-CD-SG-CB
1	B	126	SCY	CE-CD-SG-CB
1	C	126	SCY	OCD-CD-SG-CB
1	C	126	SCY	CE-CD-SG-CB
1	D	126	SCY	OCD-CD-SG-CB
1	D	126	SCY	CE-CD-SG-CB
1	A	126	SCY	CA-CB-SG-CD
1	B	126	SCY	CA-CB-SG-CD

There are no ring outliers.

4 monomers are involved in 37 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	126	SCY	10	0
1	C	126	SCY	7	0
1	B	126	SCY	9	0
1	D	126	SCY	11	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 3 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	$\# Z > 2$
3	COA	C	1003	-	47,50,50	1.80	6 (12%)	69,75,75	1.52	9 (13%)
3	COA	D	1004	-	47,50,50	1.72	6 (12%)	69,75,75	1.52	9 (13%)
3	COA	A	1001	-	47,50,50	1.73	6 (12%)	69,75,75	1.61	10 (14%)
3	COA	B	1002	-	47,50,50	1.80	6 (12%)	69,75,75	1.68	12 (17%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	COA	C	1003	-	-	9/48/64/64	0/3/3/3
3	COA	D	1004	-	-	10/48/64/64	0/3/3/3
3	COA	A	1001	-	-	14/48/64/64	0/3/3/3
3	COA	B	1002	-	-	6/48/64/64	0/3/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1001	COA	O9P-C9P	9.32	1.41	1.23
3	C	1003	COA	O9P-C9P	9.12	1.40	1.23
3	B	1002	COA	O9P-C9P	9.02	1.40	1.23
3	D	1004	COA	O9P-C9P	9.00	1.40	1.23
3	B	1002	COA	P1A-O3A	4.42	1.64	1.59
3	C	1003	COA	P2A-O3A	4.21	1.64	1.59
3	B	1002	COA	P2A-O3A	3.85	1.63	1.59
3	C	1003	COA	P1A-O3A	3.31	1.63	1.59
3	D	1004	COA	P2A-O3A	3.00	1.62	1.59
3	C	1003	COA	C2A-N1A	2.96	1.39	1.33
3	D	1004	COA	C2A-N3A	2.94	1.39	1.33
3	A	1001	COA	P2A-O3A	2.91	1.62	1.59
3	D	1004	COA	C2A-N1A	2.78	1.38	1.33
3	A	1001	COA	P1A-O3A	2.75	1.62	1.59
3	B	1002	COA	C2A-N1A	2.66	1.38	1.33
3	C	1003	COA	C2A-N3A	2.65	1.38	1.33
3	B	1002	COA	C2A-N3A	2.57	1.38	1.33
3	A	1001	COA	C2A-N1A	2.55	1.38	1.33
3	A	1001	COA	C2A-N3A	2.52	1.38	1.33
3	C	1003	COA	C8A-N7A	2.34	1.36	1.31
3	A	1001	COA	C8A-N7A	2.21	1.35	1.31
3	B	1002	COA	C8A-N7A	2.19	1.35	1.31
3	D	1004	COA	P1A-O3A	2.19	1.61	1.59
3	D	1004	COA	C8A-N7A	2.19	1.35	1.31

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1003	COA	N3A-C2A-N1A	-6.03	119.46	128.58
3	D	1004	COA	N3A-C2A-N1A	-5.48	120.29	128.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1001	COA	N3A-C2A-N1A	-5.39	120.43	128.58
3	B	1002	COA	N3A-C2A-N1A	-5.26	120.62	128.58
3	A	1001	COA	C5A-C4A-N3A	-4.90	119.97	126.72
3	B	1002	COA	N9A-C8A-N7A	-4.38	107.73	113.94
3	D	1004	COA	C5A-C4A-N3A	-4.27	120.83	126.72
3	B	1002	COA	C5A-C4A-N3A	-4.18	120.96	126.72
3	B	1002	COA	C7P-C6P-C5P	-4.03	105.68	112.39
3	C	1003	COA	C5A-C4A-N3A	-3.94	121.29	126.72
3	A	1001	COA	C5A-N7A-C8A	3.93	109.63	103.45
3	A	1001	COA	N9A-C8A-N7A	-3.89	108.42	113.94
3	D	1004	COA	N9A-C8A-N7A	-3.75	108.62	113.94
3	C	1003	COA	N9A-C8A-N7A	-3.70	108.69	113.94
3	B	1002	COA	C5A-N7A-C8A	3.68	109.23	103.45
3	C	1003	COA	C2A-N3A-C4A	3.47	120.31	111.83
3	D	1004	COA	C2A-N3A-C4A	3.46	120.28	111.83
3	A	1001	COA	C2A-N3A-C4A	3.46	120.28	111.83
3	A	1001	COA	O4B-C1B-N9A	3.30	114.42	108.09
3	C	1003	COA	C5A-N7A-C8A	3.28	108.61	103.45
3	D	1004	COA	C5A-N7A-C8A	3.24	108.54	103.45
3	A	1001	COA	C4A-C5A-N7A	-3.15	106.98	110.58
3	B	1002	COA	C2A-N3A-C4A	3.11	119.42	111.83
3	A	1001	COA	N3A-C4A-N9A	3.02	132.30	127.17
3	D	1004	COA	O4B-C1B-N9A	3.00	113.86	108.09
3	B	1002	COA	O5P-C5P-C6P	-2.97	116.63	122.02
3	D	1004	COA	N3A-C4A-N9A	2.94	132.17	127.17
3	B	1002	COA	N3A-C4A-N9A	2.75	131.84	127.17
3	B	1002	COA	C4A-C5A-N7A	-2.58	107.63	110.58
3	C	1003	COA	N3A-C4A-N9A	2.43	131.29	127.17
3	C	1003	COA	C4A-N9A-C1B	-2.42	120.98	126.63
3	D	1004	COA	C4A-C5A-N7A	-2.41	107.83	110.58
3	B	1002	COA	C4A-N9A-C8A	2.35	108.21	105.74
3	B	1002	COA	C2B-C3B-C4B	2.34	107.33	103.24
3	C	1003	COA	O4B-C1B-N9A	2.31	112.52	108.09
3	C	1003	COA	C4A-C5A-N7A	-2.28	107.98	110.58
3	B	1002	COA	O3B-C3B-C2B	-2.18	103.84	111.68
3	D	1004	COA	C4A-N9A-C8A	2.16	108.01	105.74
3	A	1001	COA	O6A-CCP-CBP	2.14	113.99	110.55
3	A	1001	COA	C6A-C5A-C4A	2.09	120.04	117.18

There are no chirality outliers.

All (39) torsion outliers are listed below:

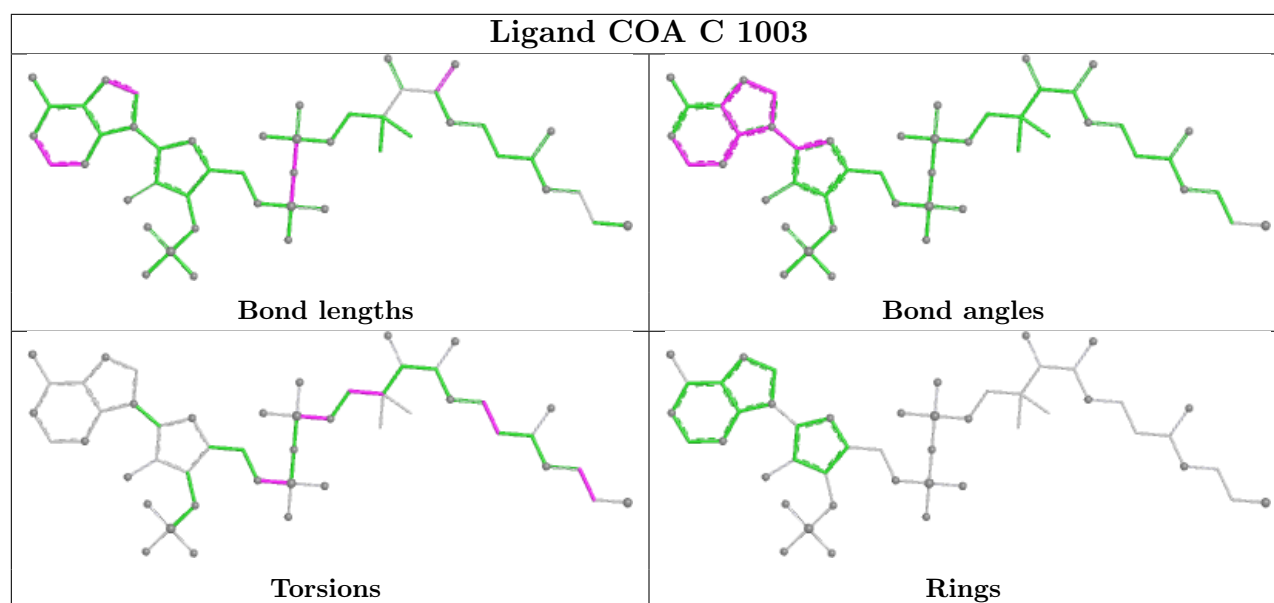
Mol	Chain	Res	Type	Atoms
3	A	1001	COA	C5B-O5B-P1A-O1A
3	A	1001	COA	C5B-O5B-P1A-O2A
3	A	1001	COA	C5B-O5B-P1A-O3A
3	A	1001	COA	CDP-CBP-CCP-O6A
3	A	1001	COA	CEP-CBP-CCP-O6A
3	A	1001	COA	CAP-CBP-CCP-O6A
3	A	1001	COA	S1P-C2P-C3P-N4P
3	B	1002	COA	C5B-O5B-P1A-O1A
3	B	1002	COA	C5B-O5B-P1A-O3A
3	B	1002	COA	C6P-C5P-N4P-C3P
3	B	1002	COA	O5P-C5P-N4P-C3P
3	C	1003	COA	C5B-O5B-P1A-O1A
3	C	1003	COA	CCP-O6A-P2A-O4A
3	C	1003	COA	CDP-CBP-CCP-O6A
3	C	1003	COA	CEP-CBP-CCP-O6A
3	C	1003	COA	CAP-CBP-CCP-O6A
3	C	1003	COA	S1P-C2P-C3P-N4P
3	D	1004	COA	C5B-O5B-P1A-O1A
3	D	1004	COA	C5B-O5B-P1A-O3A
3	D	1004	COA	CCP-O6A-P2A-O3A
3	D	1004	COA	CCP-O6A-P2A-O4A
3	D	1004	COA	CDP-CBP-CCP-O6A
3	D	1004	COA	CEP-CBP-CCP-O6A
3	D	1004	COA	CAP-CBP-CCP-O6A
3	D	1004	COA	S1P-C2P-C3P-N4P
3	A	1001	COA	O5P-C5P-N4P-C3P
3	A	1001	COA	C6P-C5P-N4P-C3P
3	A	1001	COA	O4B-C4B-C5B-O5B
3	A	1001	COA	CCP-O6A-P2A-O3A
3	A	1001	COA	CCP-O6A-P2A-O4A
3	A	1001	COA	CCP-O6A-P2A-O5A
3	B	1002	COA	C5B-O5B-P1A-O2A
3	C	1003	COA	C5B-O5B-P1A-O2A
3	C	1003	COA	C5B-O5B-P1A-O3A
3	D	1004	COA	C5B-O5B-P1A-O2A
3	C	1003	COA	C5P-C6P-C7P-N8P
3	A	1001	COA	C3B-C4B-C5B-O5B
3	D	1004	COA	C4B-C5B-O5B-P1A
3	B	1002	COA	P1A-O3A-P2A-O5A

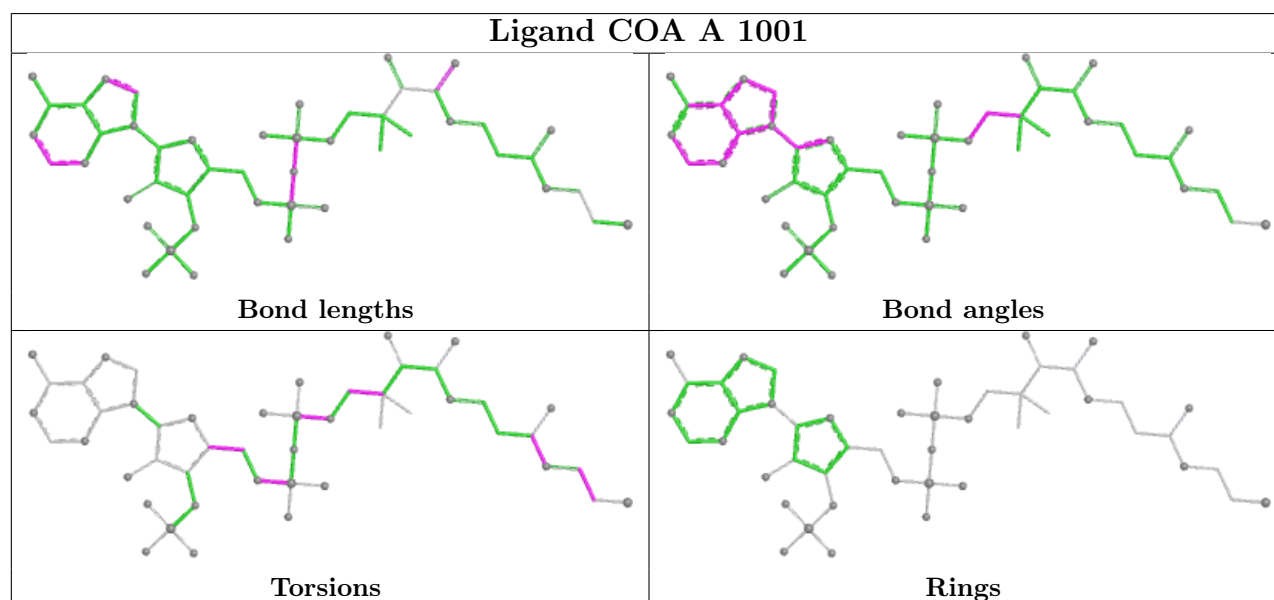
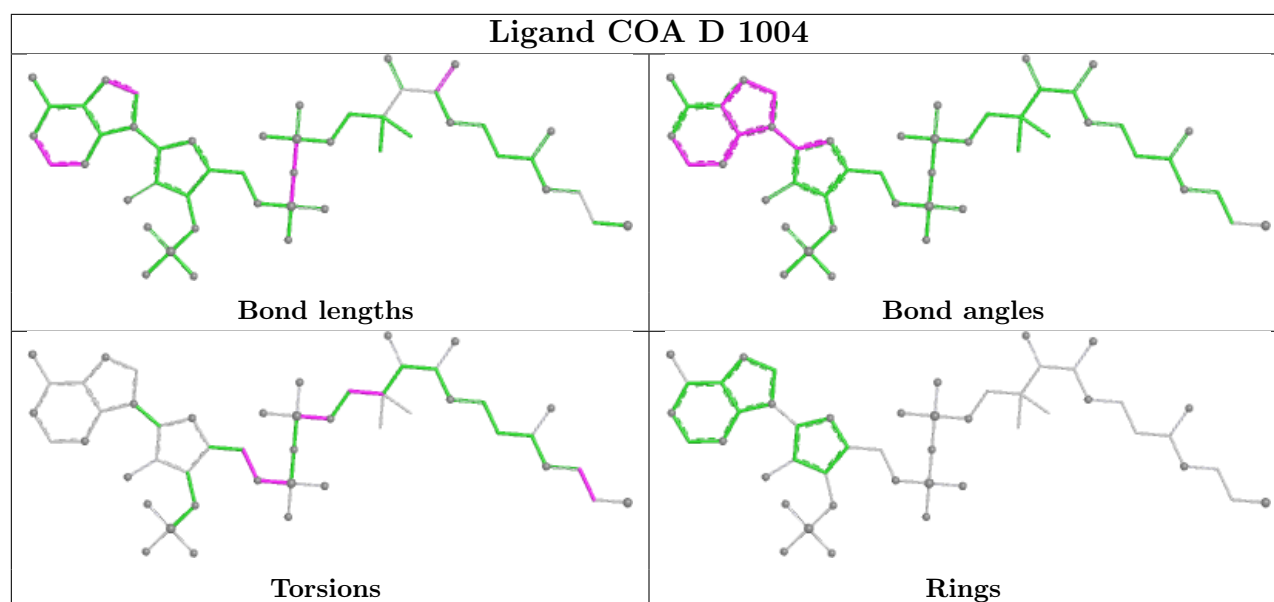
There are no ring outliers.

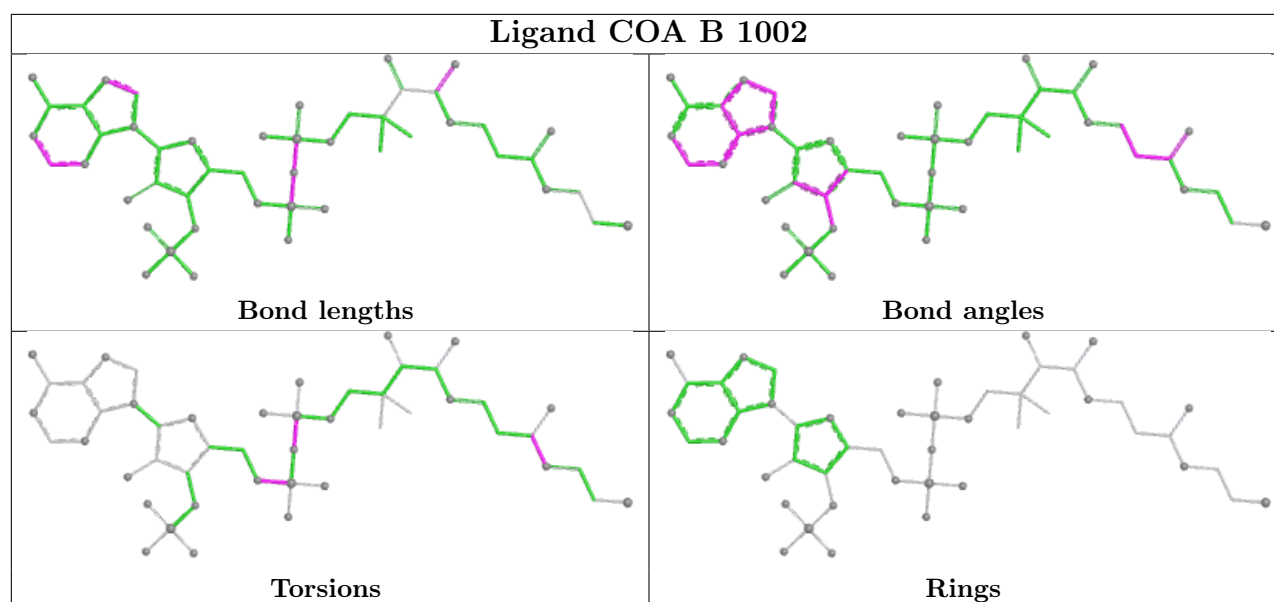
3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1003	COA	2	0
3	A	1001	COA	2	0
3	B	1002	COA	5	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1
1	D	1
1	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	125:VAL	C	126:SCY	N	1.75
1	D	126:SCY	C	127:ALA	N	1.73
1	C	126:SCY	C	127:ALA	N	1.66

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	388/406 (95%)	0.53	15 (3%) 43 42	30, 48, 69, 80	0
1	B	385/406 (94%)	-0.13	0 100 100	21, 34, 49, 55	0
1	C	387/406 (95%)	0.08	7 (1%) 67 67	23, 38, 60, 71	0
1	D	381/406 (93%)	0.06	4 (1%) 79 79	30, 41, 58, 67	0
All	All	1541/1624 (94%)	0.14	26 (1%) 69 68	21, 41, 63, 80	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	32	SER	4.2
1	A	242	VAL	3.7
1	C	261	PHE	3.5
1	A	301	ALA	3.0
1	C	265	PRO	3.0
1	D	40	SER	2.7
1	C	246	PRO	2.6
1	C	259	VAL	2.6
1	D	248	VAL	2.5
1	A	216	ILE	2.5
1	A	270	VAL	2.4
1	A	228	ALA	2.3
1	A	173	VAL	2.3
1	A	266	LYS	2.3
1	A	427	LEU	2.3
1	D	173	VAL	2.2
1	D	242	VAL	2.2
1	A	298	ALA	2.2
1	A	347	ILE	2.2
1	A	215	ALA	2.2
1	C	216	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	369	ILE	2.1
1	A	381	VAL	2.1
1	C	201	ALA	2.1
1	A	346	ASP	2.1
1	C	264	VAL	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	SCY	A	126	9/10	0.83	0.15	43,46,53,54	0
1	SCY	C	126	9/10	0.87	0.12	31,33,48,56	0
1	SCY	B	126	9/10	0.92	0.11	25,28,42,51	0
1	SCY	D	126	9/10	0.95	0.08	32,34,48,50	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.

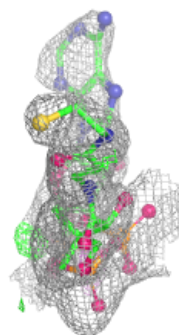
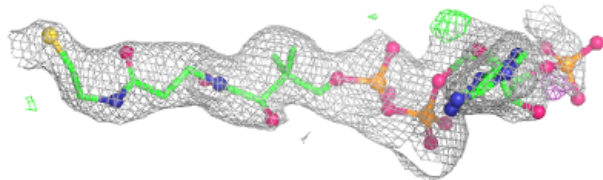
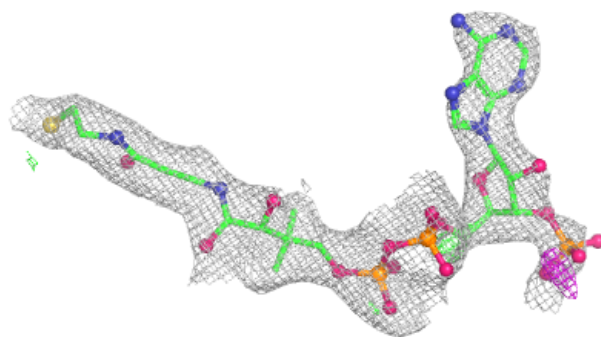
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	COA	A	1001	48/48	0.81	0.14	71,98,105,106	0
3	COA	C	1003	48/48	0.84	0.12	69,81,93,93	0
3	COA	B	1002	48/48	0.85	0.12	44,62,71,73	0
3	COA	D	1004	48/48	0.86	0.11	57,67,78,79	0
2	CL	A	2003	1/1	0.97	0.06	42,42,42,42	0
2	CL	B	2001	1/1	0.99	0.03	29,29,29,29	0
2	CL	B	2002	1/1	0.99	0.03	34,34,34,34	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.

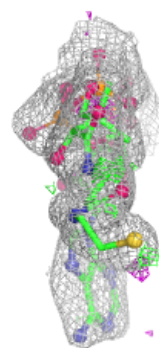
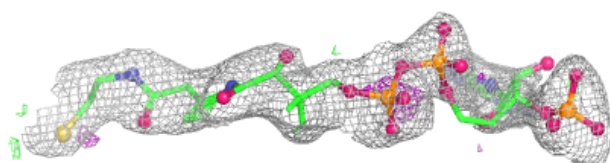
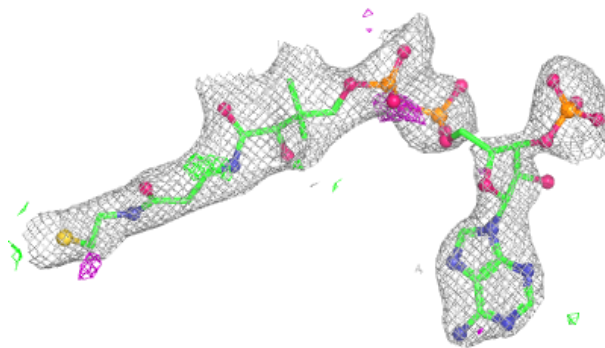
Electron density around COA A 1001:

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

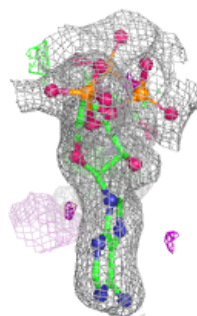
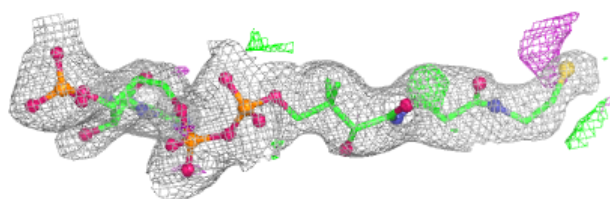
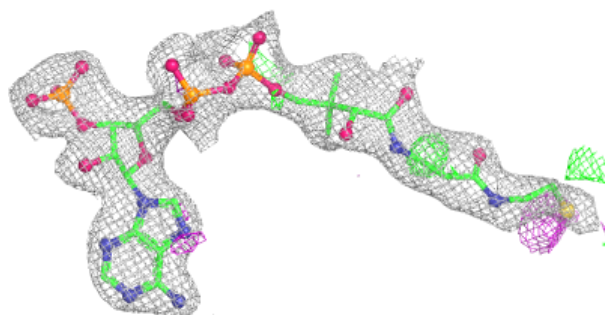


Electron density around COA C 1003:

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

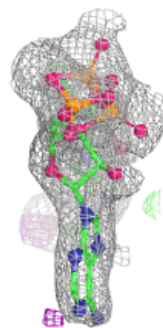
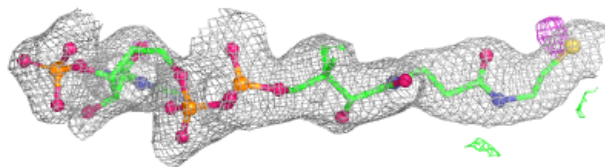
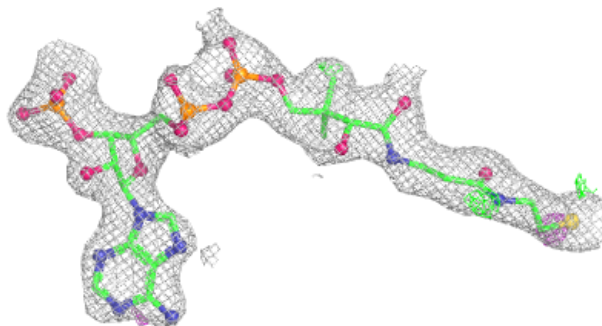
**Electron density around COA B 1002:**

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



Electron density around COA D 1004:

$2mF_o-DF_c$ (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.