



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 05:59 AM GMT

PDB ID : 2VAT
Title : CRYSTAL STRUCTURE OF DEACETYLCEPHALOSPORIN C ACETYL-TRANSFERASE IN COMPLEX WITH COENZYME A
Authors : Lejon, S.; Ellis, J.; Valegard, K.
Deposited on : 2007-09-04
Resolution : 2.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

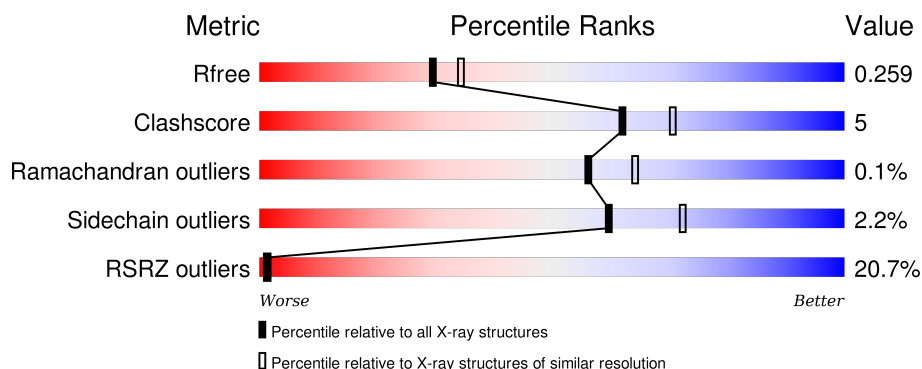
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	444	<div> <div>3%</div> <div>68% 10% 22%</div> </div>
1	B	444	<div> <div>8%</div> <div>71% 7% 22%</div> </div>
1	C	444	<div> <div>11%</div> <div>69% 7% 22%</div> </div>
1	D	444	<div> <div>8%</div> <div>70% 6% 22%</div> </div>
1	E	444	<div> <div>3%</div> <div>70% 7% 22%</div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	444	
1	G	444	
1	H	444	
1	I	444	
1	J	444	
1	K	444	
1	L	444	

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
2	COA	B	1383	-	-	-	X
2	COA	C	1383	-	-	-	X
2	COA	D	1383	-	-	-	X
2	COA	F	1383	-	-	-	X
2	COA	G	1383	-	-	-	X
2	COA	L	1383	-	-	-	X
3	ACT	A	1384	-	-	-	X
3	ACT	C	1384	-	-	-	X
3	ACT	D	1384	-	-	-	X
3	ACT	F	1384	-	-	-	X
3	ACT	H	1384	-	-	X	-
3	ACT	I	1385	-	-	-	X
3	ACT	K	1386	-	-	X	-
4	GOL	B	1384	-	-	-	X
4	GOL	I	1384	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 35013 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--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	347	Total	C	N	O	S	0	4	0
			2744	1731	485	509	19			
1	B	347	Total	C	N	O	S	0	3	0
			2744	1731	488	506	19			
1	C	345	Total	C	N	O	S	0	4	0
			2745	1730	488	508	19			
1	D	345	Total	C	N	O	S	0	4	0
			2748	1732	488	509	19			
1	E	346	Total	C	N	O	S	0	5	0
			2744	1732	487	506	19			
1	F	342	Total	C	N	O	S	0	2	0
			2694	1700	474	501	19			
1	G	342	Total	C	N	O	S	0	6	0
			2733	1724	484	506	19			
1	H	345	Total	C	N	O	S	0	8	0
			2760	1741	492	508	19			
1	I	347	Total	C	N	O	S	0	3	0
			2739	1728	485	507	19			
1	J	344	Total	C	N	O	S	0	4	0
			2724	1719	482	504	19			
1	K	344	Total	C	N	O	S	0	5	0
			2732	1724	485	504	19			
1	L	341	Total	C	N	O	S	0	3	0
			2697	1702	476	500	19			

- Molecule 2 is COENZYME A (three-letter code: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S).



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

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	F	1	Total	C	O	0	0
			4	2	2		
3	G	1	Total	C	O	0	0
			4	2	2		
3	H	1	Total	C	O	0	0
			4	2	2		
3	I	1	Total	C	O	0	0
			4	2	2		
3	J	1	Total	C	O	0	0
			4	2	2		
3	K	1	Total	C	O	0	0
			4	2	2		
3	L	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	I	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	165	Total	O	0	0
			165	165		
5	B	171	Total	O	0	0
			171	171		
5	C	137	Total	O	0	0
			137	137		
5	D	123	Total	O	0	0
			123	123		
5	E	165	Total	O	0	0
			165	165		
5	F	102	Total	O	0	0
			102	102		
5	G	140	Total	O	0	0
			140	140		
5	H	144	Total	O	0	0
			144	144		
5	I	164	Total	O	0	0
			164	164		
5	J	80	Total	O	0	0
			80	80		

Continued on next page...

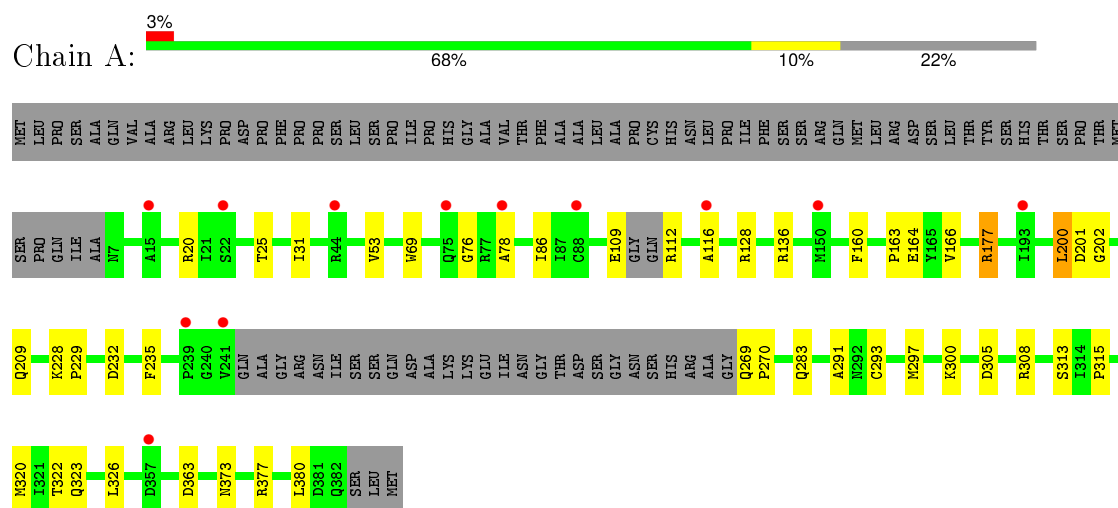
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	K	78	Total 78	O 78	0	0
5	L	104	Total 104	O 104	0	0

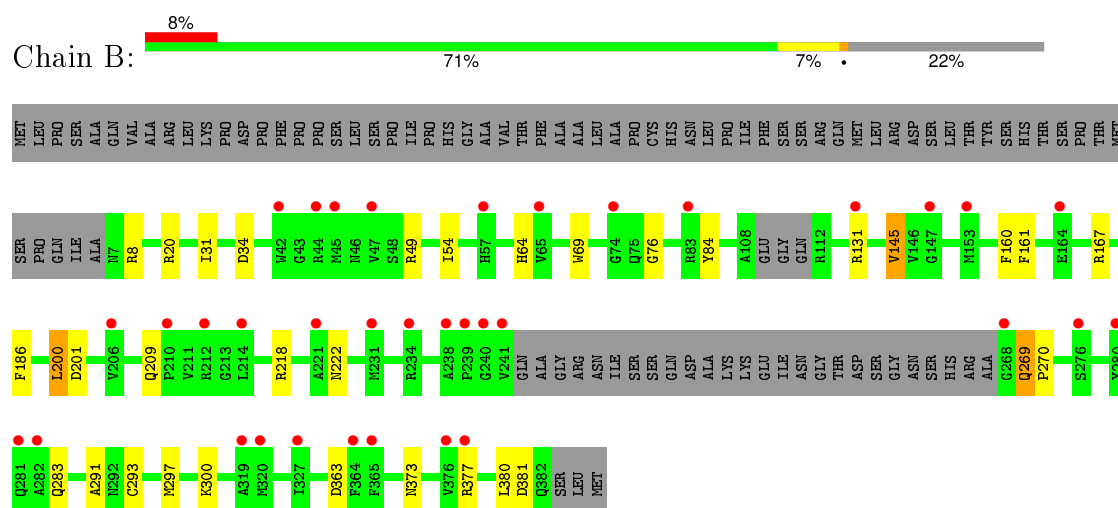
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

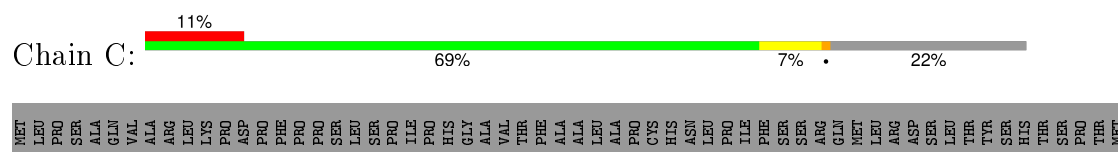
• Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

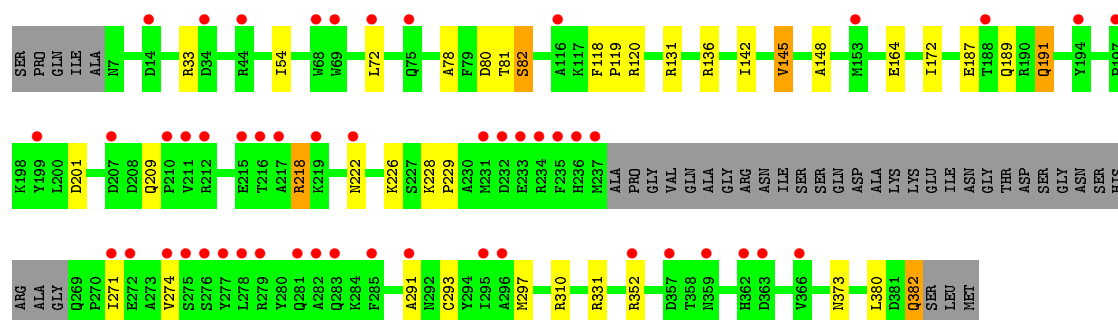


• Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

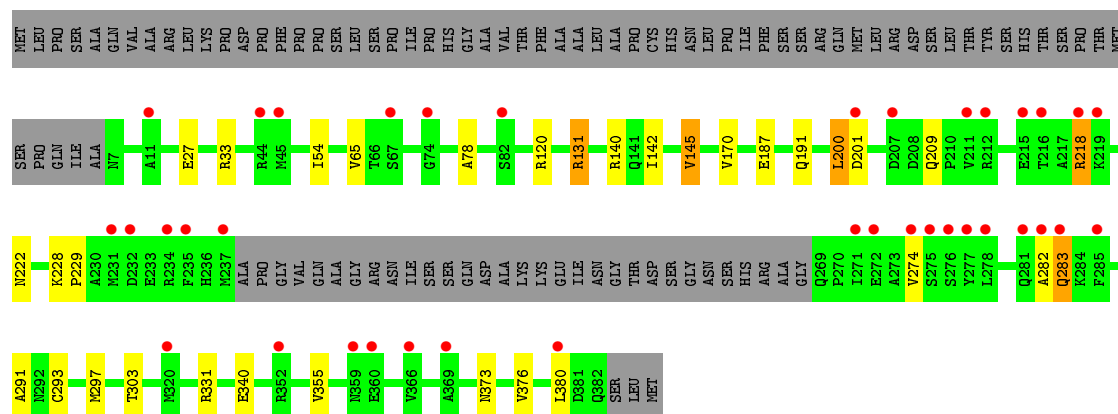


• Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

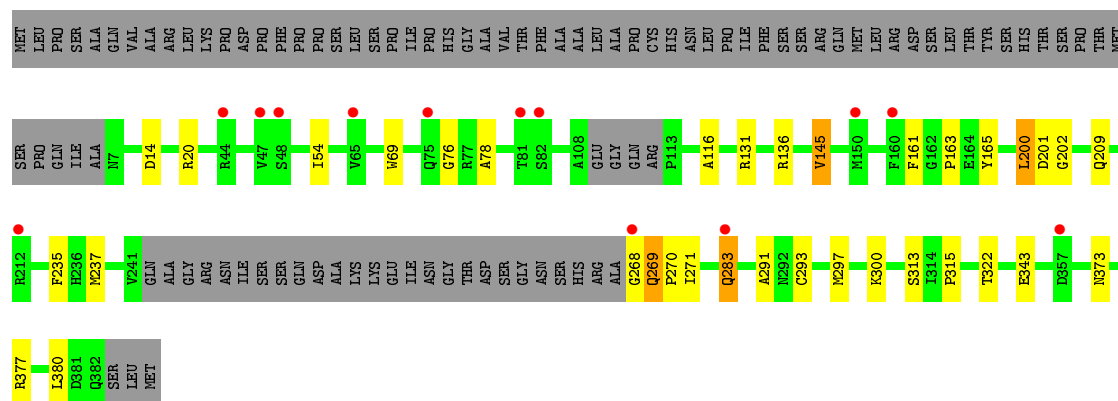




• Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

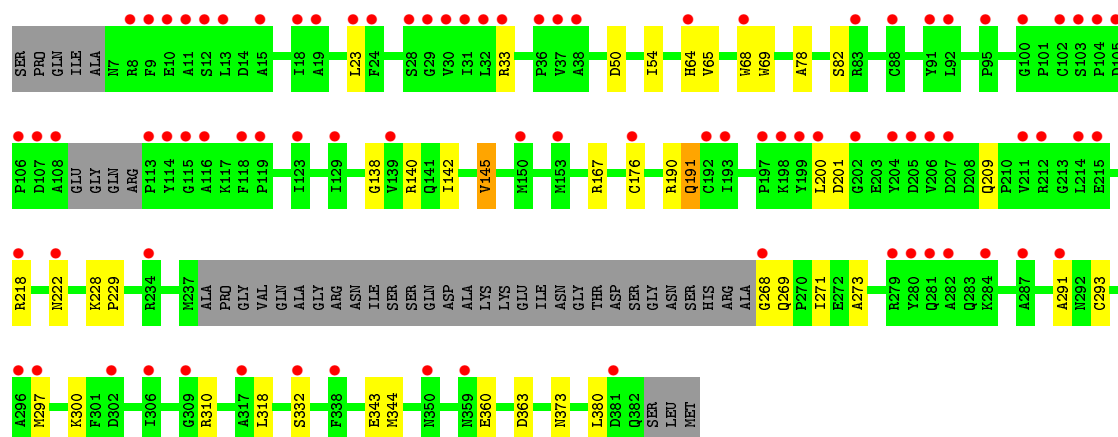


• Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

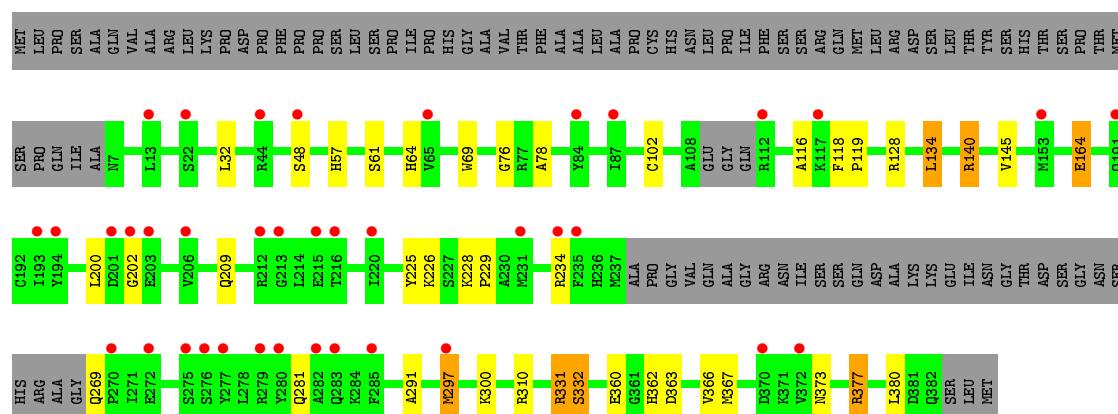


• Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

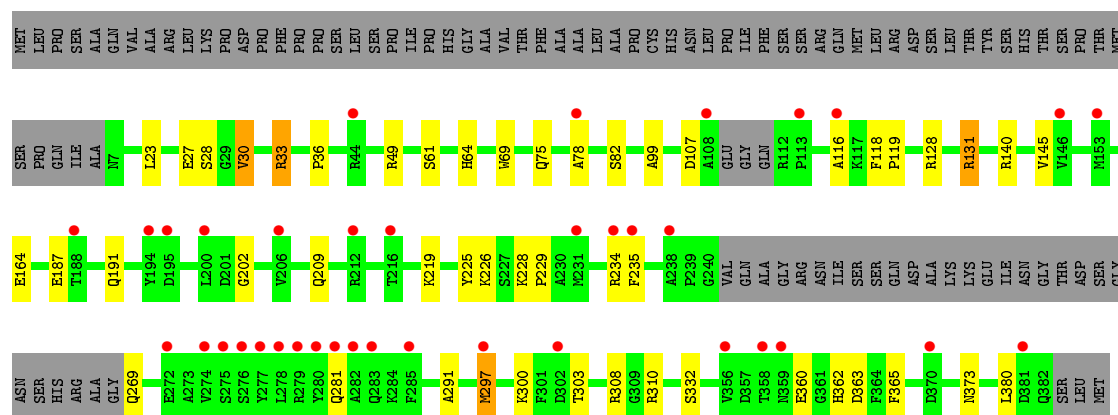




● Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

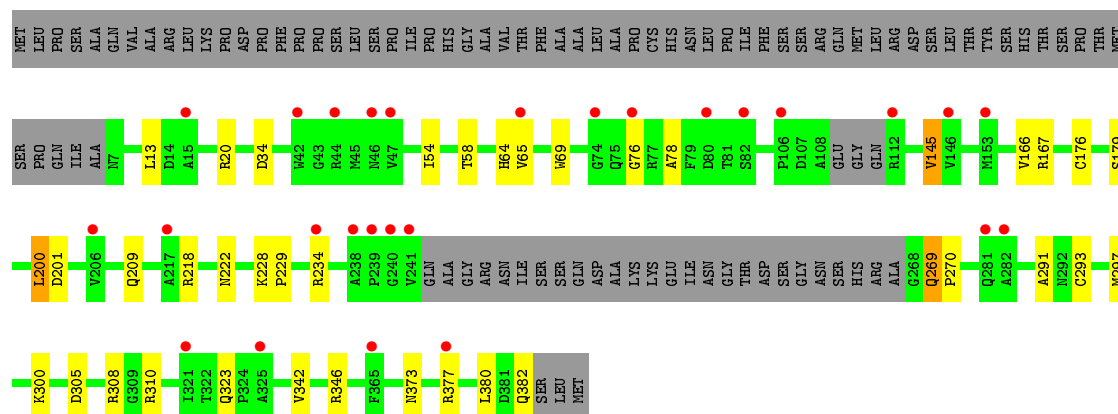


● Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

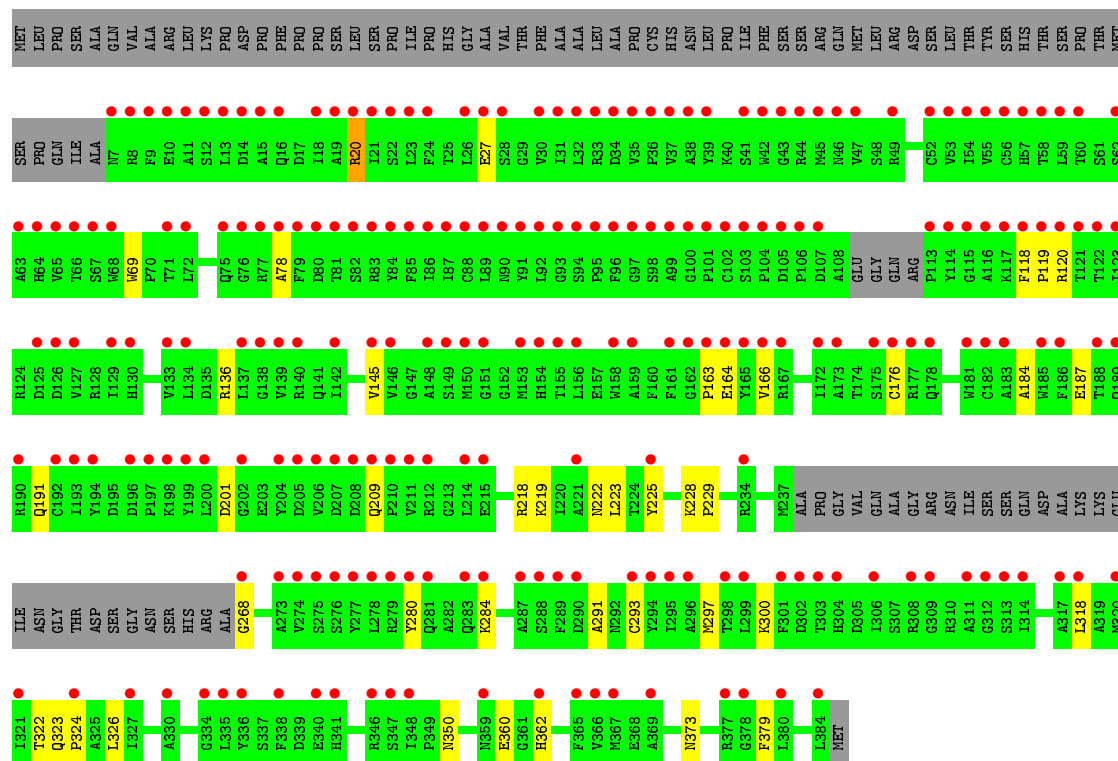


● Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

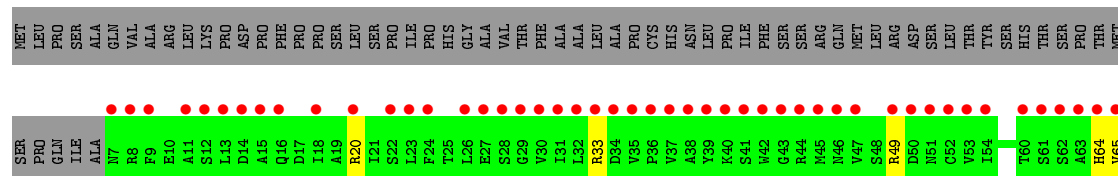


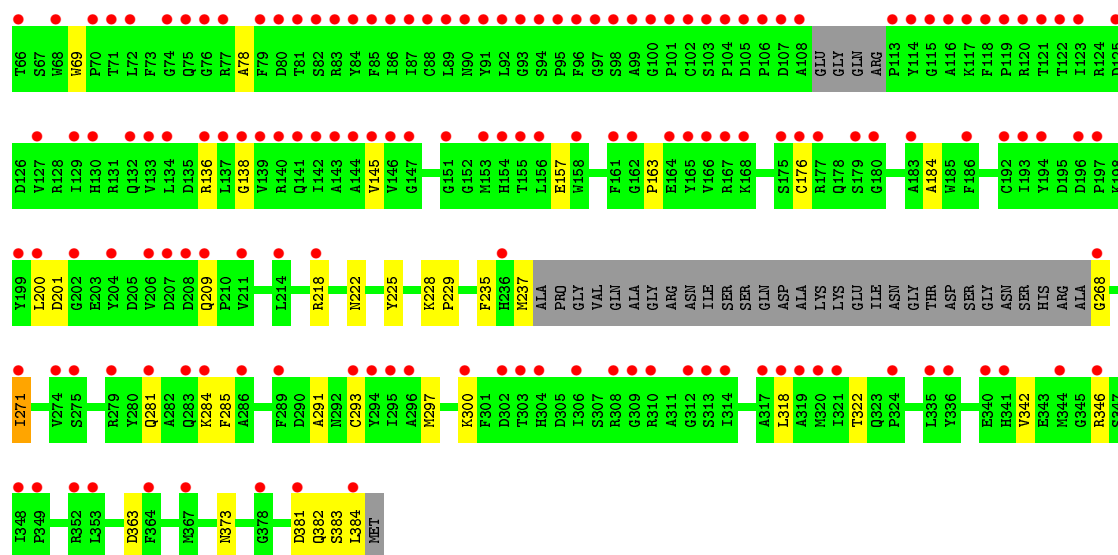


● Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE

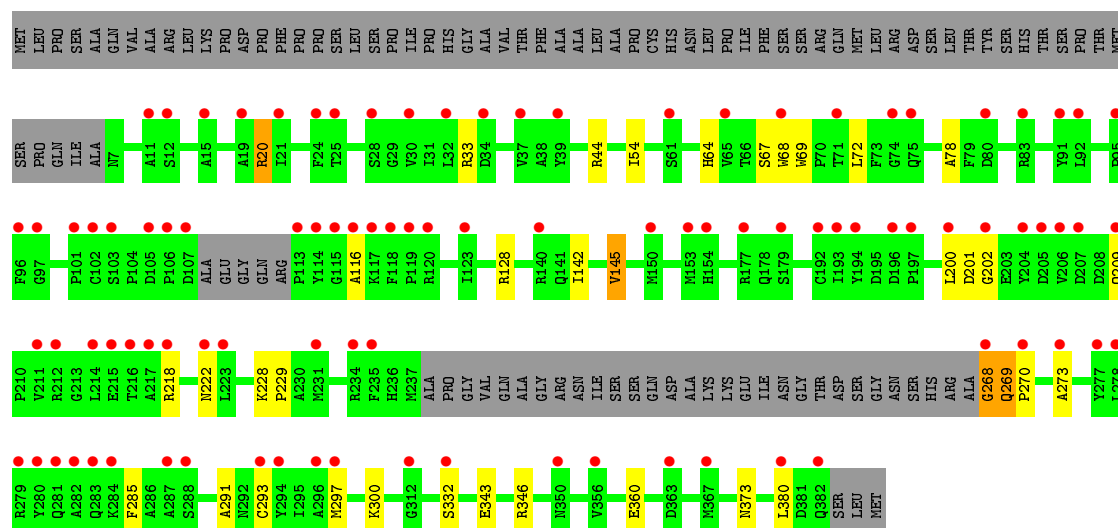


● Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE





● Molecule 1: ACETYL-COA--DEACETYLCEPHALOSPORIN C ACETYLTRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	121.96Å 109.28Å 197.00Å 90.00° 90.23° 90.00°	Depositor
Resolution (Å)	122.17 – 2.20 27.45 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.5 (122.17-2.20) 98.9 (27.45-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.200 , 0.230 0.263 , 0.259	Depositor DCC
R_{free} test set	10106 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 16.2	EDS
Estimated twinning fraction	0.479 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 200784 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	35013	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.89	3/2816 (0.1%)	0.75	4/3817 (0.1%)
1	B	0.99	10/2816 (0.4%)	0.76	3/3816 (0.1%)
1	C	0.89	6/2820 (0.2%)	0.71	1/3821 (0.0%)
1	D	0.94	5/2826 (0.2%)	0.74	2/3829 (0.1%)
1	E	0.85	1/2822 (0.0%)	0.70	0/3823
1	F	0.89	5/2759 (0.2%)	0.75	6/3739 (0.2%)
1	G	0.87	3/2813 (0.1%)	0.73	3/3811 (0.1%)
1	H	0.84	0/2844	0.73	4/3852 (0.1%)
1	I	0.87	2/2811 (0.1%)	0.73	3/3810 (0.1%)
1	J	0.89	2/2795 (0.1%)	0.72	1/3786 (0.0%)
1	K	0.86	1/2806 (0.0%)	0.69	0/3800
1	L	0.86	1/2765 (0.0%)	0.71	0/3746
All	All	0.89	39/33693 (0.1%)	0.73	27/45650 (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	L	0	1

The worst 5 of 39 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	186	PHE	CE1-CZ	-11.74	1.15	1.37
1	B	186	PHE	CG-CD2	-11.20	1.22	1.38
1	D	340	GLU	CD-OE1	-10.64	1.14	1.25
1	D	340	GLU	CD-OE2	-10.50	1.14	1.25
1	B	186	PHE	CE2-CZ	-9.88	1.18	1.37

The worst 5 of 27 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	177	ARG	NE-CZ-NH1	11.47	126.04	120.30
1	B	8	ARG	NE-CZ-NH2	10.40	125.50	120.30
1	D	340	GLU	OE1-CD-OE2	-10.00	111.30	123.30
1	D	27	GLU	OE1-CD-OE2	-9.88	111.44	123.30
1	J	27	GLU	OE1-CD-OE2	-9.78	111.56	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	268	GLY	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	2744	0	2656	25	0
1	B	2744	0	2663	18	0
1	C	2745	0	2662	30	0
1	D	2748	0	2669	29	0
1	E	2744	0	2669	26	0
1	F	2694	0	2599	30	0
1	G	2733	0	2652	27	0
1	H	2760	0	2685	44	0
1	I	2739	0	2655	33	0
1	J	2724	0	2641	33	0
1	K	2732	0	2653	44	0
1	L	2697	0	2607	33	0
2	A	48	0	32	2	0
2	B	48	0	32	2	0
2	C	48	0	32	0	0
2	D	48	0	32	0	0
2	E	48	0	31	1	0
2	F	48	0	32	3	0
2	G	48	0	32	5	0
2	H	48	0	32	6	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	I	48	0	32	1	0
2	J	48	0	32	3	0
2	K	48	0	32	4	0
2	L	48	0	32	1	0
3	A	4	0	3	0	0
3	B	4	0	3	1	0
3	C	4	0	3	0	0
3	D	4	0	3	0	0
3	E	4	0	3	0	0
3	F	4	0	3	0	0
3	G	4	0	3	1	0
3	H	4	0	3	3	0
3	I	4	0	3	0	0
3	J	4	0	3	0	0
3	K	4	0	3	2	0
3	L	4	0	3	1	0
4	B	6	0	8	0	0
4	I	6	0	8	0	0
5	A	165	0	0	0	0
5	B	171	0	0	6	0
5	C	137	0	0	8	0
5	D	123	0	0	5	0
5	E	165	0	0	3	0
5	F	102	0	0	7	0
5	G	140	0	0	3	0
5	H	144	0	0	5	0
5	I	164	0	0	3	0
5	J	80	0	0	3	0
5	K	78	0	0	5	0
5	L	104	0	0	7	0
All	All	35013	0	32246	357	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

The worst 5 of 357 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:K:218[A]:ARG:NE	1:K:222[A]:ASN:HD21	1.04	1.42
1:K:218[A]:ARG:NE	1:K:222[A]:ASN:ND2	1.78	1.29
1:J:218:ARG:CZ	1:J:222[A]:ASN:OD1	1.98	1.11

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:363[A]:ASP:OD1	2:H:1383:COA:C6P	1.99	1.11
1:E:14:ASP:OD1	5:E:2005:HOH:O	1.72	1.08

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	343/444 (77%)	332 (97%)	11 (3%)	0	100	100
1	B	343/444 (77%)	332 (97%)	10 (3%)	1 (0%)	46	50
1	C	344/444 (78%)	334 (97%)	10 (3%)	0	100	100
1	D	345/444 (78%)	335 (97%)	10 (3%)	0	100	100
1	E	344/444 (78%)	333 (97%)	10 (3%)	1 (0%)	46	50
1	F	336/444 (76%)	325 (97%)	11 (3%)	0	100	100
1	G	341/444 (77%)	332 (97%)	9 (3%)	0	100	100
1	H	345/444 (78%)	334 (97%)	11 (3%)	0	100	100
1	I	343/444 (77%)	333 (97%)	9 (3%)	1 (0%)	46	50
1	J	340/444 (77%)	327 (96%)	13 (4%)	0	100	100
1	K	341/444 (77%)	328 (96%)	13 (4%)	0	100	100
1	L	336/444 (76%)	327 (97%)	8 (2%)	1 (0%)	46	50
All	All	4101/5328 (77%)	3972 (97%)	125 (3%)	4 (0%)	56	64

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	269	GLN
1	E	269	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	I	269	GLN
1	L	269	GLN

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	293/372 (79%)	286 (98%)	7 (2%)	57	69
1	B	292/372 (78%)	286 (98%)	6 (2%)	61	74
1	C	293/372 (79%)	286 (98%)	7 (2%)	57	69
1	D	294/372 (79%)	286 (97%)	8 (3%)	52	64
1	E	293/372 (79%)	287 (98%)	6 (2%)	63	76
1	F	287/372 (77%)	280 (98%)	7 (2%)	57	69
1	G	293/372 (79%)	283 (97%)	10 (3%)	44	54
1	H	295/372 (79%)	286 (97%)	9 (3%)	47	59
1	I	292/372 (78%)	287 (98%)	5 (2%)	68	81
1	J	291/372 (78%)	287 (99%)	4 (1%)	74	85
1	K	292/372 (78%)	286 (98%)	6 (2%)	61	74
1	L	288/372 (77%)	281 (98%)	7 (2%)	57	69
All	All	3503/4464 (78%)	3421 (98%)	82 (2%)	60	71

5 of 82 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	201	ASP
1	G	145	VAL
1	L	20	ARG
1	F	269	GLN
1	G	69	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 41 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	269	GLN
1	G	269	GLN
1	K	373	ASN
1	F	373	ASN
1	G	64	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

26 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	A	1383	-	40,50,50	2.19	14 (35%)	50,75,75	2.17	9 (18%)
3	ACT	A	1384	-	1,3,3	1.27	0	0,3,3	0.00	-
2	COA	B	1383	-	40,50,50	1.69	3 (7%)	50,75,75	2.13	6 (12%)
4	GOL	B	1384	-	5,5,5	0.30	0	5,5,5	0.36	0
3	ACT	B	1385	-	1,3,3	0.68	0	0,3,3	0.00	-
2	COA	C	1383	-	40,50,50	1.85	4 (10%)	50,75,75	2.07	4 (8%)
3	ACT	C	1384	-	1,3,3	0.91	0	0,3,3	0.00	-
2	COA	D	1383	-	40,50,50	1.85	4 (10%)	50,75,75	2.03	5 (10%)
3	ACT	D	1384	-	1,3,3	0.28	0	0,3,3	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	COA	E	1383	-	40,50,50	2.13	12 (30%)	50,75,75	2.03	9 (18%)
3	ACT	E	1384	-	1,3,3	1.31	0	0,3,3	0.00	-
2	COA	F	1383	-	40,50,50	1.86	3 (7%)	50,75,75	2.03	3 (6%)
3	ACT	F	1384	-	1,3,3	1.74	0	0,3,3	0.00	-
2	COA	G	1383	-	40,50,50	1.79	3 (7%)	50,75,75	1.92	5 (10%)
3	ACT	G	1384	-	1,3,3	1.11	0	0,3,3	0.00	-
2	COA	H	1383	-	40,50,50	2.03	9 (22%)	50,75,75	2.33	11 (22%)
3	ACT	H	1384	-	1,3,3	1.87	0	0,3,3	0.00	-
2	COA	I	1383	-	40,50,50	1.70	3 (7%)	50,75,75	2.15	7 (14%)
4	GOL	I	1384	-	5,5,5	0.35	0	5,5,5	0.21	0
3	ACT	I	1385	-	1,3,3	1.44	0	0,3,3	0.00	-
2	COA	J	1385	-	40,50,50	1.80	4 (10%)	50,75,75	1.95	3 (6%)
3	ACT	J	1386	-	1,3,3	1.71	0	0,3,3	0.00	-
2	COA	K	1385	-	40,50,50	1.81	4 (10%)	50,75,75	2.12	5 (10%)
3	ACT	K	1386	-	1,3,3	1.75	0	0,3,3	0.00	-
2	COA	L	1383	-	40,50,50	1.80	3 (7%)	50,75,75	2.15	3 (6%)
3	ACT	L	1384	-	1,3,3	1.77	0	0,3,3	0.00	-

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	COA	A	1383	-	-	2/44/64/64	0/3/3/3
3	ACT	A	1384	-	-	0/0/0/0	0/0/0/0
2	COA	B	1383	-	-	1/44/64/64	0/3/3/3
4	GOL	B	1384	-	-	0/4/4/4	0/0/0/0
3	ACT	B	1385	-	-	0/0/0/0	0/0/0/0
2	COA	C	1383	-	-	0/44/64/64	0/3/3/3
3	ACT	C	1384	-	-	0/0/0/0	0/0/0/0
2	COA	D	1383	-	-	0/44/64/64	0/3/3/3
3	ACT	D	1384	-	-	0/0/0/0	0/0/0/0
2	COA	E	1383	-	-	1/44/64/64	0/3/3/3
3	ACT	E	1384	-	-	0/0/0/0	0/0/0/0
2	COA	F	1383	-	-	1/44/64/64	0/3/3/3
3	ACT	F	1384	-	-	0/0/0/0	0/0/0/0
2	COA	G	1383	-	-	0/44/64/64	0/3/3/3
3	ACT	G	1384	-	-	0/0/0/0	0/0/0/0
2	COA	H	1383	-	-	0/44/64/64	0/3/3/3
3	ACT	H	1384	-	-	0/0/0/0	0/0/0/0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	COA	I	1383	-	-	1/44/64/64	0/3/3/3
4	GOL	I	1384	-	-	0/4/4/4	0/0/0/0
3	ACT	I	1385	-	-	0/0/0/0	0/0/0/0
2	COA	J	1385	-	-	0/44/64/64	0/3/3/3
3	ACT	J	1386	-	-	0/0/0/0	0/0/0/0
2	COA	K	1385	-	-	0/44/64/64	0/3/3/3
3	ACT	K	1386	-	-	0/0/0/0	0/0/0/0
2	COA	L	1383	-	-	0/44/64/64	0/3/3/3
3	ACT	L	1384	-	-	0/0/0/0	0/0/0/0

The worst 5 of 66 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	1383	COA	P3B-O8A	-3.78	1.41	1.54
2	A	1383	COA	P1A-O2A	-3.48	1.40	1.54
2	E	1383	COA	P1A-O2A	-3.47	1.40	1.54
2	A	1383	COA	O4B-C1B	-3.40	1.36	1.41
2	A	1383	COA	P3B-O8A	-3.26	1.43	1.54

The worst 5 of 70 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	1383	COA	N3A-C2A-N1A	-13.37	118.66	128.89
2	K	1385	COA	N3A-C2A-N1A	-12.92	119.00	128.89
2	I	1383	COA	N3A-C2A-N1A	-12.61	119.24	128.89
2	F	1383	COA	N3A-C2A-N1A	-12.35	119.44	128.89
2	B	1383	COA	N3A-C2A-N1A	-12.27	119.50	128.89

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	I	1383	COA	CAP-C9P-N8P-C7P
2	A	1383	COA	C6P-C5P-N4P-C3P
2	B	1383	COA	CAP-C9P-N8P-C7P
2	F	1383	COA	CAP-C9P-N8P-C7P
2	E	1383	COA	CAP-C9P-N8P-C7P

There are no ring outliers.

15 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1383	COA	2	0
2	B	1383	COA	2	0
3	B	1385	ACT	1	0
2	E	1383	COA	1	0
2	F	1383	COA	3	0
2	G	1383	COA	5	0
3	G	1384	ACT	1	0
2	H	1383	COA	6	0
3	H	1384	ACT	3	0
2	I	1383	COA	1	0
2	J	1385	COA	3	0
2	K	1385	COA	4	0
3	K	1386	ACT	2	0
2	L	1383	COA	1	0
3	L	1384	ACT	1	0

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	347/444 (78%)	0.79	12 (3%)	48	46	11, 22, 38, 57	2 (0%)
1	B	347/444 (78%)	1.10	35 (10%)	9	8	12, 22, 38, 51	1 (0%)
1	C	345/444 (77%)	1.14	50 (14%)	3	3	12, 22, 39, 59	1 (0%)
1	D	345/444 (77%)	1.04	37 (10%)	8	7	12, 22, 38, 55	0
1	E	346/444 (77%)	0.79	13 (3%)	44	43	9, 22, 36, 51	1 (0%)
1	F	342/444 (77%)	1.40	84 (24%)	1	1	12, 23, 37, 51	2 (0%)
1	G	342/444 (77%)	1.10	38 (11%)	7	7	12, 22, 37, 51	1 (0%)
1	H	345/444 (77%)	1.00	37 (10%)	8	7	12, 22, 38, 52	2 (0%)
1	I	347/444 (78%)	1.04	27 (7%)	16	15	10, 22, 38, 51	1 (0%)
1	J	344/444 (77%)	2.91	227 (65%)	0	0	13, 23, 38, 54	2 (0%)
1	K	344/444 (77%)	2.66	200 (58%)	0	0	13, 23, 38, 53	2 (0%)
1	L	341/444 (76%)	1.45	96 (28%)	1	1	12, 23, 37, 51	2 (0%)
All	All	4135/5328 (77%)	1.37	856 (20%)	1	1	9, 22, 38, 59	17 (0%)

The worst 5 of 856 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	193	ILE	11.6
1	K	106	PRO	11.5
1	J	104	PRO	10.6
1	K	104	PRO	10.5
1	K	193	ILE	10.5

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	ACT	F	1384	4/4	0.70	0.39	15.37	39,40,42,42	0
4	GOL	B	1384	6/6	0.74	0.42	11.04	33,36,39,40	0
3	ACT	I	1385	4/4	0.80	0.29	5.40	20,20,22,23	0
2	COA	D	1383	48/48	0.47	0.57	4.89	35,47,70,72	0
3	ACT	A	1384	4/4	0.91	0.28	4.79	27,27,28,29	0
3	ACT	C	1384	4/4	0.94	0.28	4.51	30,31,31,31	0
2	COA	C	1383	48/48	0.59	0.58	4.22	36,46,73,80	0
3	ACT	D	1384	4/4	0.92	0.32	3.97	28,29,29,29	0
2	COA	F	1383	48/48	0.62	0.50	3.16	39,59,69,70	0
4	GOL	I	1384	6/6	0.79	0.24	3.03	35,37,39,44	0
2	COA	L	1383	48/48	0.67	0.49	2.89	37,57,65,65	0
2	COA	G	1383	48/48	0.74	0.37	2.38	21,53,83,86	0
2	COA	B	1383	48/48	0.74	0.35	2.21	28,42,54,57	0
2	COA	H	1383	48/48	0.78	0.37	1.67	21,53,75,76	0
3	ACT	E	1384	4/4	0.94	0.24	1.67	30,31,31,31	0
3	ACT	H	1384	4/4	0.85	0.21	1.62	31,32,32,33	0
2	COA	I	1383	48/48	0.81	0.30	1.43	29,41,57,59	0
3	ACT	L	1384	4/4	0.82	0.21	0.56	46,48,48,48	0
3	ACT	K	1386	4/4	0.84	0.32	0.51	45,45,45,46	0
2	COA	A	1383	48/48	0.92	0.18	0.22	17,29,48,50	0
2	COA	J	1385	48/48	0.76	0.29	0.10	23,55,68,70	0
3	ACT	B	1385	4/4	0.84	0.20	0.02	20,22,23,24	0
2	COA	K	1385	48/48	0.78	0.24	-0.08	20,50,65,72	0
3	ACT	G	1384	4/4	0.85	0.18	-0.13	37,37,38,38	0
3	ACT	J	1386	4/4	0.86	0.24	-0.71	47,49,50,50	0
2	COA	E	1383	48/48	0.93	0.15	-1.31	14,27,49,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.