



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:05 PM UTC

PDB ID : 8ECM / pdb_00008ecm
Title : Crystal Structure Analysis of Acetyl-CoA acetyltransferase from Firmicutes bacterium
Authors : Seo, H.S.; Dhe-Paganon, S.
Deposited on : 2022-09-02
Resolution : 1.89 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
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

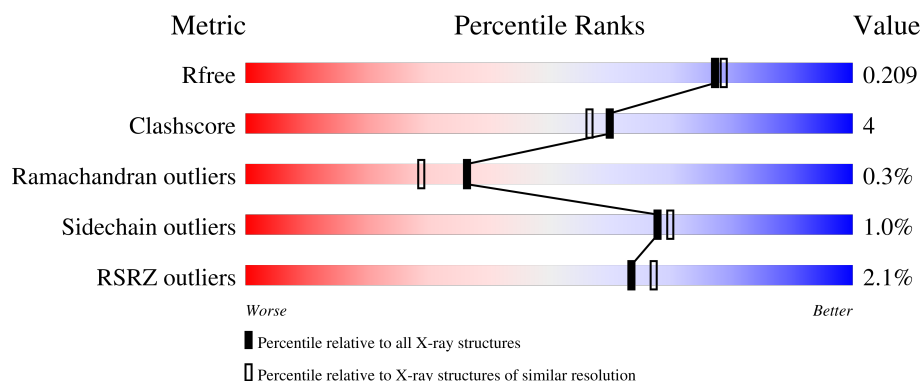
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	428	2% 87% 10% ..
1	B	428	% 88% 10% .
1	C	428	% 88% 10% .
1	D	428	% 91% 7% .
1	E	428	% 89% 9% .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	428	% 91% 7%
1	G	428	% 91% 7%
1	H	428	9% 82% 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 25714 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	419	Total	C	N	O	S	0	0	0
			3044	1904	530	584	26			
1	B	420	Total	C	N	O	S	0	0	0
			3056	1911	532	587	26			
1	C	421	Total	C	N	O	S	0	0	0
			3065	1917	534	588	26			
1	D	423	Total	C	N	O	S	0	0	0
			3085	1929	540	590	26			
1	E	421	Total	C	N	O	S	0	0	0
			3077	1925	536	590	26			
1	F	423	Total	C	N	O	S	0	0	0
			3092	1934	540	592	26			
1	G	420	Total	C	N	O	S	0	0	0
			3063	1917	533	587	26			
1	H	417	Total	C	N	O	S	0	0	0
			3039	1902	530	582	25			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP R6CZ24
A	2	ALA	-	expression tag	UNP R6CZ24
A	421	LEU	-	expression tag	UNP R6CZ24
A	422	GLU	-	expression tag	UNP R6CZ24
A	423	HIS	-	expression tag	UNP R6CZ24
A	424	HIS	-	expression tag	UNP R6CZ24
A	425	HIS	-	expression tag	UNP R6CZ24
A	426	HIS	-	expression tag	UNP R6CZ24
A	427	HIS	-	expression tag	UNP R6CZ24
A	428	HIS	-	expression tag	UNP R6CZ24
B	1	MET	-	initiating methionine	UNP R6CZ24
B	2	ALA	-	expression tag	UNP R6CZ24
B	421	LEU	-	expression tag	UNP R6CZ24

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	422	GLU	-	expression tag	UNP R6CZ24
B	423	HIS	-	expression tag	UNP R6CZ24
B	424	HIS	-	expression tag	UNP R6CZ24
B	425	HIS	-	expression tag	UNP R6CZ24
B	426	HIS	-	expression tag	UNP R6CZ24
B	427	HIS	-	expression tag	UNP R6CZ24
B	428	HIS	-	expression tag	UNP R6CZ24
C	1	MET	-	initiating methionine	UNP R6CZ24
C	2	ALA	-	expression tag	UNP R6CZ24
C	421	LEU	-	expression tag	UNP R6CZ24
C	422	GLU	-	expression tag	UNP R6CZ24
C	423	HIS	-	expression tag	UNP R6CZ24
C	424	HIS	-	expression tag	UNP R6CZ24
C	425	HIS	-	expression tag	UNP R6CZ24
C	426	HIS	-	expression tag	UNP R6CZ24
C	427	HIS	-	expression tag	UNP R6CZ24
C	428	HIS	-	expression tag	UNP R6CZ24
D	1	MET	-	initiating methionine	UNP R6CZ24
D	2	ALA	-	expression tag	UNP R6CZ24
D	421	LEU	-	expression tag	UNP R6CZ24
D	422	GLU	-	expression tag	UNP R6CZ24
D	423	HIS	-	expression tag	UNP R6CZ24
D	424	HIS	-	expression tag	UNP R6CZ24
D	425	HIS	-	expression tag	UNP R6CZ24
D	426	HIS	-	expression tag	UNP R6CZ24
D	427	HIS	-	expression tag	UNP R6CZ24
D	428	HIS	-	expression tag	UNP R6CZ24
E	1	MET	-	initiating methionine	UNP R6CZ24
E	2	ALA	-	expression tag	UNP R6CZ24
E	421	LEU	-	expression tag	UNP R6CZ24
E	422	GLU	-	expression tag	UNP R6CZ24
E	423	HIS	-	expression tag	UNP R6CZ24
E	424	HIS	-	expression tag	UNP R6CZ24
E	425	HIS	-	expression tag	UNP R6CZ24
E	426	HIS	-	expression tag	UNP R6CZ24
E	427	HIS	-	expression tag	UNP R6CZ24
E	428	HIS	-	expression tag	UNP R6CZ24
F	1	MET	-	initiating methionine	UNP R6CZ24
F	2	ALA	-	expression tag	UNP R6CZ24
F	421	LEU	-	expression tag	UNP R6CZ24
F	422	GLU	-	expression tag	UNP R6CZ24
F	423	HIS	-	expression tag	UNP R6CZ24

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
F	424	HIS	-	expression tag	UNP R6CZ24
F	425	HIS	-	expression tag	UNP R6CZ24
F	426	HIS	-	expression tag	UNP R6CZ24
F	427	HIS	-	expression tag	UNP R6CZ24
F	428	HIS	-	expression tag	UNP R6CZ24
G	1	MET	-	initiating methionine	UNP R6CZ24
G	2	ALA	-	expression tag	UNP R6CZ24
G	421	LEU	-	expression tag	UNP R6CZ24
G	422	GLU	-	expression tag	UNP R6CZ24
G	423	HIS	-	expression tag	UNP R6CZ24
G	424	HIS	-	expression tag	UNP R6CZ24
G	425	HIS	-	expression tag	UNP R6CZ24
G	426	HIS	-	expression tag	UNP R6CZ24
G	427	HIS	-	expression tag	UNP R6CZ24
G	428	HIS	-	expression tag	UNP R6CZ24
H	1	MET	-	initiating methionine	UNP R6CZ24
H	2	ALA	-	expression tag	UNP R6CZ24
H	421	LEU	-	expression tag	UNP R6CZ24
H	422	GLU	-	expression tag	UNP R6CZ24
H	423	HIS	-	expression tag	UNP R6CZ24
H	424	HIS	-	expression tag	UNP R6CZ24
H	425	HIS	-	expression tag	UNP R6CZ24
H	426	HIS	-	expression tag	UNP R6CZ24
H	427	HIS	-	expression tag	UNP R6CZ24
H	428	HIS	-	expression tag	UNP R6CZ24

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	132	Total O 132 132	0	0
2	B	150	Total O 150 150	0	0
2	C	161	Total O 161 161	0	0
2	D	170	Total O 170 170	0	0
2	E	170	Total O 170 170	0	0
2	F	181	Total O 181 181	0	0
2	G	156	Total O 156 156	0	0

Continued on next page...

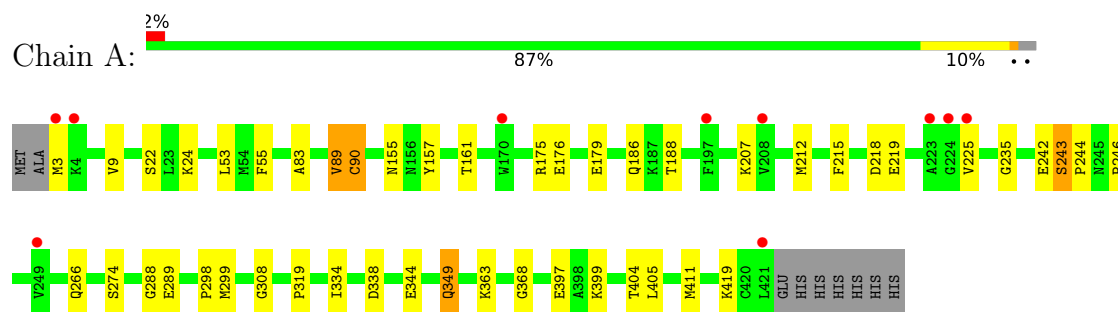
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	H	73	Total	O	0	0
			73	73		

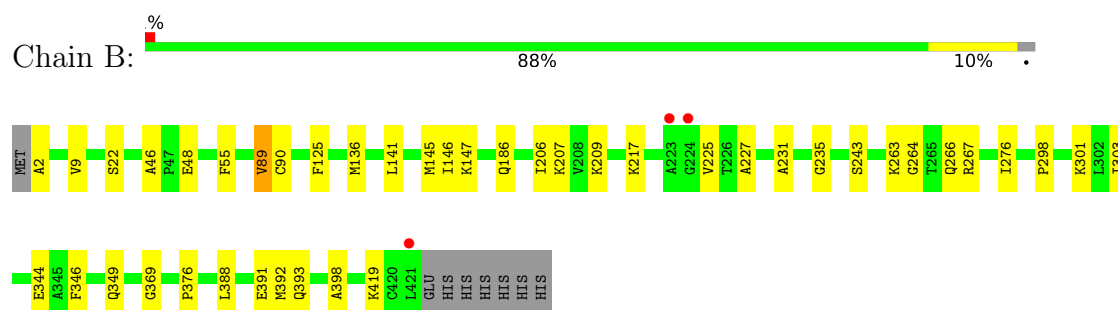
3 Residue-property plots

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.

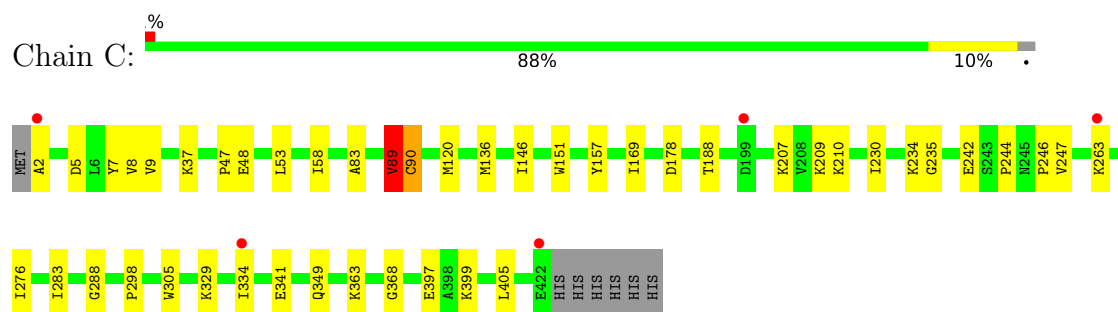
• Molecule 1: Acetyl-CoA acetyltransferase



• Molecule 1: Acetyl-CoA acetyltransferase

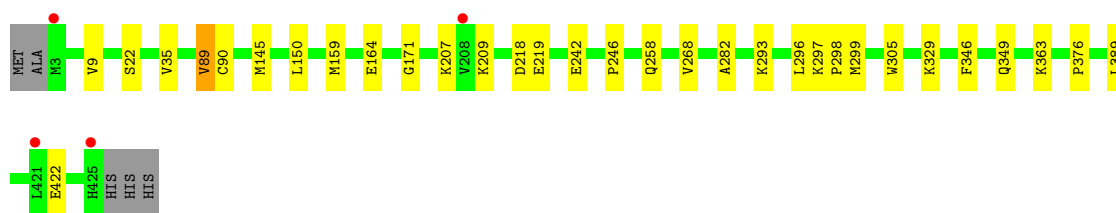


• Molecule 1: Acetyl-CoA acetyltransferase

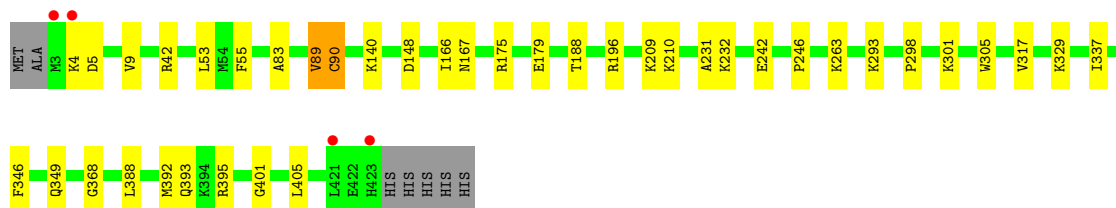
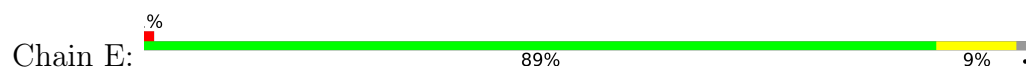


• Molecule 1: Acetyl-CoA acetyltransferase

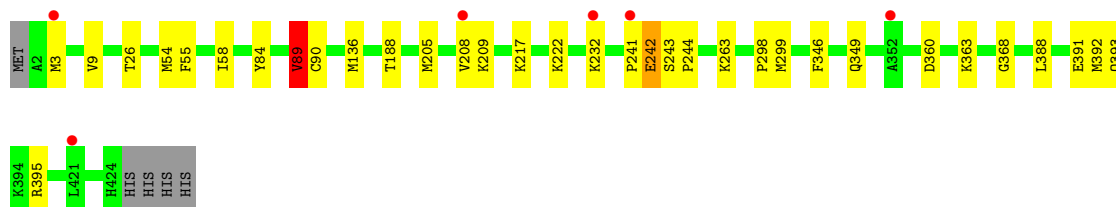




- Molecule 1: Acetyl-CoA acetyltransferase



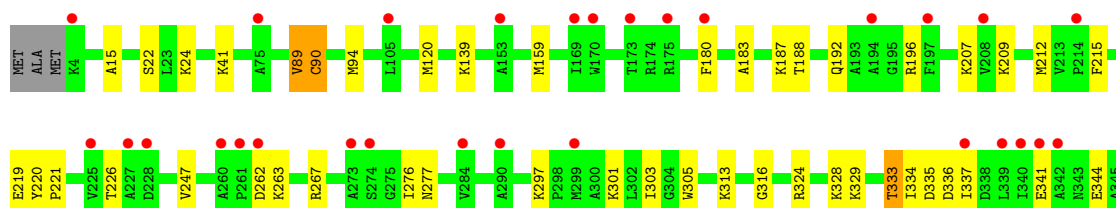
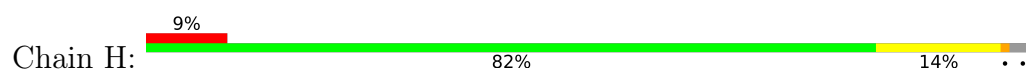
- Molecule 1: Acetyl-CoA acetyltransferase

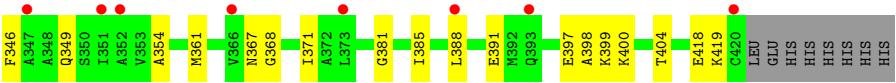


- Molecule 1: Acetyl-CoA acetyltransferase



- Molecule 1: Acetyl-CoA acetyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	75.23Å 81.46Å 160.19Å 95.61° 100.45° 108.27°	Depositor
Resolution (Å)	73.46 – 1.89 73.46 – 1.89	Depositor EDS
% Data completeness (in resolution range)	94.0 (73.46-1.89) 94.1 (73.46-1.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.176 , 0.209 0.176 , 0.209	Depositor DCC
R_{free} test set	13250 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	39.0	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	25714	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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

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	0.48	0/3082	0.64	0/4169
1	B	0.55	0/3094	0.69	2/4184 (0.0%)
1	C	0.57	0/3103	0.68	0/4195
1	D	0.56	0/3126	0.70	1/4229 (0.0%)
1	E	0.63	0/3116	0.74	0/4212
1	F	0.61	0/3132	0.70	0/4234
1	G	0.56	0/3101	0.67	0/4192
1	H	0.47	0/3077	0.66	0/4160
All	All	0.56	0/24831	0.68	3/33575 (0.0%)

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	B	0	1
1	D	0	1
1	E	0	1
All	All	0	3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	145	MET	CA-CB-CG	5.42	124.93	114.10
1	B	2	ALA	CA-C-N	-5.06	110.69	121.32
1	B	2	ALA	C-N-CA	-5.06	110.69	121.32

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	376	PRO	Peptide
1	D	376	PRO	Peptide
1	E	196	ARG	Sidechain

5.2 Too-close contacts [i](#)

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	3044	0	3052	32	0
1	B	3056	0	3070	23	0
1	C	3065	0	3083	30	1
1	D	3085	0	3086	17	0
1	E	3077	0	3098	26	1
1	F	3092	0	3110	22	0
1	G	3063	0	3090	19	0
1	H	3039	0	3063	44	0
2	A	132	0	0	3	0
2	B	150	0	0	5	0
2	C	161	0	0	4	0
2	D	170	0	0	3	0
2	E	170	0	0	5	0
2	F	181	0	0	5	0
2	G	156	0	0	5	0
2	H	73	0	0	2	0
All	All	25714	0	24652	209	1

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 4.

All (209) 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:E:392:MET:HE1	1:E:401:GLY:HA3	1.39	1.02
1:E:167:ASN:O	1:E:263:LYS:NZ	1.92	1.02
1:G:242:GLU:HG3	1:G:246:PRO:HA	1.49	0.93

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:363:LYS:NZ	2:D:501:HOH:O	2.02	0.91
1:H:354:ALA:HB1	1:H:361:MET:HE1	1.50	0.91
1:A:176:GLU:OE1	2:A:501:HOH:O	1.90	0.88
1:E:242:GLU:OE2	2:E:501:HOH:O	1.94	0.84
1:A:207:LYS:HG3	1:A:212:MET:HE1	1.65	0.79
1:E:392:MET:CE	1:E:401:GLY:HA3	2.12	0.79
1:C:207:LYS:HE2	1:C:210:LYS:HD2	1.65	0.79
1:E:4:LYS:N	2:E:502:HOH:O	2.16	0.78
1:H:346:PHE:H	1:H:349:GLN:HE21	1.32	0.77
1:C:334:ILE:HD12	1:C:334:ILE:H	1.49	0.77
1:C:341:GLU:OE1	2:C:501:HOH:O	2.02	0.77
1:G:189:GLU:O	2:G:501:HOH:O	2.03	0.75
1:C:234:LYS:HD2	1:C:235:GLY:N	2.01	0.75
1:A:308:GLY:O	1:A:411:MET:HB3	1.90	0.72
1:C:234:LYS:HD2	1:C:235:GLY:H	1.56	0.71
1:A:3:MET:HE3	1:A:289:GLU:HG3	1.73	0.70
1:H:196:ARG:NH2	1:H:391:GLU:OE2	2.23	0.70
1:D:346:PHE:H	1:D:349:GLN:HE21	1.40	0.70
1:F:26:THR:OG1	2:F:501:HOH:O	2.09	0.70
1:B:393:GLN:NE2	2:B:501:HOH:O	2.10	0.70
1:F:244:PRO:HB2	1:F:263:LYS:HG2	1.74	0.70
1:E:392:MET:HE1	1:E:401:GLY:CA	2.18	0.69
1:C:363:LYS:NZ	2:C:502:HOH:O	2.26	0.68
1:H:333:THR:HG23	1:H:335:ASP:H	1.59	0.67
1:C:244:PRO:HB2	1:C:263:LYS:HZ3	1.59	0.67
1:H:212:MET:HE3	1:H:212:MET:HA	1.77	0.67
1:H:367:ASN:ND2	1:H:391:GLU:OE1	2.21	0.67
1:B:346:PHE:H	1:B:349:GLN:HE21	1.43	0.66
1:F:393:GLN:HG2	2:F:541:HOH:O	1.94	0.66
1:H:297:LYS:HD2	1:H:297:LYS:O	1.97	0.65
1:A:242:GLU:HB3	1:A:246:PRO:HA	1.77	0.65
1:A:24:LYS:NZ	2:A:504:HOH:O	2.29	0.64
1:B:46:ALA:HB1	1:B:48:GLU:OE1	1.98	0.64
1:E:346:PHE:H	1:E:349:GLN:HE21	1.46	0.64
1:A:243:SER:HB2	1:A:244:PRO:HD3	1.79	0.64
1:C:37:LYS:NZ	2:C:504:HOH:O	2.30	0.63
1:A:338:ASP:O	1:A:363:LYS:HG2	1.99	0.62
1:H:183:ALA:O	1:H:187:LYS:HG3	2.01	0.61
1:A:22:SER:OG	1:A:219:GLU:OE2	2.09	0.61
1:H:344:GLU:HG3	1:H:371:ILE:HG13	1.83	0.60
1:H:180:PHE:HE1	1:H:344:GLU:OE1	1.84	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:263:LYS:HA	1:H:267:ARG:NH1	2.17	0.59
1:H:305:TRP:O	1:H:329:LYS:HE3	2.02	0.59
1:D:207:LYS:NZ	1:D:209:LYS:O	2.32	0.59
1:D:293:LYS:NZ	2:D:503:HOH:O	2.36	0.59
1:C:329:LYS:HG3	2:C:530:HOH:O	2.02	0.58
1:H:22:SER:OG	1:H:219:GLU:OE2	2.15	0.58
1:H:120:MET:HG3	1:H:277:ASN:O	2.03	0.58
1:H:398:ALA:O	1:H:419:LYS:HE2	2.04	0.58
1:F:388:LEU:O	1:F:392:MET:HG3	2.04	0.58
1:D:22:SER:OG	1:D:219:GLU:OE2	2.11	0.57
1:D:299:MET:HE2	1:D:389:LEU:HB3	1.87	0.56
1:F:346:PHE:H	1:F:349:GLN:HE21	1.53	0.56
1:G:4:LYS:NZ	1:G:104:ILE:O	2.36	0.56
1:A:175:ARG:O	1:A:179:GLU:HG3	2.05	0.56
1:D:242:GLU:HB2	1:D:246:PRO:HA	1.87	0.56
1:H:400:LYS:NZ	1:H:418:GLU:OE2	2.38	0.55
1:H:301:LYS:HD3	1:H:303:ILE:HG22	1.89	0.55
1:H:354:ALA:CB	1:H:361:MET:HE1	2.29	0.55
1:G:135:ARG:HH11	1:G:135:ARG:HG2	1.72	0.55
1:H:94:MET:HG3	1:H:385:ILE:HD11	1.88	0.55
1:C:2:ALA:HA	1:C:5:ASP:OD2	2.08	0.54
1:G:346:PHE:H	1:G:349:GLN:HE21	1.55	0.54
1:F:217:LYS:NZ	2:F:504:HOH:O	2.41	0.53
1:F:136:MET:HG2	1:G:146:ILE:HD11	1.90	0.53
1:C:334:ILE:H	1:C:334:ILE:CD1	2.19	0.53
1:C:305:TRP:O	1:C:329:LYS:HD3	2.08	0.53
1:E:393:GLN:HG3	2:E:623:HOH:O	2.08	0.53
1:D:299:MET:CE	1:D:389:LEU:HB3	2.39	0.52
1:F:54:MET:HG2	1:F:84:TYR:CE1	2.45	0.52
1:E:5:ASP:CG	1:E:301:LYS:HE3	2.35	0.51
1:E:242:GLU:HB2	1:E:246:PRO:HA	1.93	0.51
1:A:9:VAL:HA	1:A:298:PRO:HA	1.91	0.51
1:H:263:LYS:HA	1:H:267:ARG:HH12	1.73	0.51
1:E:140:LYS:HD3	1:H:139:LYS:O	2.10	0.51
1:F:391:GLU:OE2	1:F:395:ARG:NE	2.40	0.50
1:E:175:ARG:O	1:E:179:GLU:HG3	2.11	0.50
1:H:90:SCY:O	2:H:501:HOH:O	2.19	0.50
1:F:9:VAL:O	1:F:299:MET:HG3	2.11	0.50
1:B:146:ILE:HD11	1:C:136:MET:HG2	1.92	0.49
1:E:305:TRP:H	1:E:329:LYS:HD2	1.77	0.49
1:C:188:THR:OG1	1:C:368:GLY:HA3	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48:GLU:H	1:B:48:GLU:CD	2.20	0.49
1:B:231:ALA:HB2	2:B:502:HOH:O	2.13	0.49
1:H:94:MET:HB2	1:H:381:GLY:O	2.12	0.49
1:H:333:THR:N	1:H:336:ASP:OD2	2.41	0.49
1:H:344:GLU:HG3	1:H:371:ILE:H	1.77	0.49
1:C:48:GLU:CD	1:C:48:GLU:H	2.21	0.48
1:A:207:LYS:HG3	1:A:212:MET:CE	2.40	0.48
1:A:3:MET:HA	1:A:288:GLY:HA3	1.96	0.48
1:A:9:VAL:O	1:A:299:MET:HG3	2.13	0.48
1:F:222:LYS:HE2	2:F:513:HOH:O	2.12	0.48
1:F:9:VAL:HA	1:F:298:PRO:HA	1.96	0.48
1:H:15:ALA:N	2:H:502:HOH:O	2.22	0.48
1:A:90:SCY:O	1:A:405:LEU:HD11	2.14	0.48
1:H:324:ARG:O	1:H:328:LYS:HG3	2.14	0.48
1:B:388:LEU:O	1:B:392:MET:HG3	2.14	0.47
1:A:235:GLY:HA3	1:A:266:GLN:HB3	1.97	0.47
1:B:227:ALA:O	2:B:502:HOH:O	2.20	0.47
1:H:159:MET:HA	1:H:159:MET:HE2	1.95	0.47
1:D:305:TRP:H	1:D:329:LYS:HD2	1.80	0.47
1:G:237:PHE:O	1:G:266:GLN:HG2	2.14	0.47
1:G:360:ASP:HB3	1:G:363:LYS:HD2	1.95	0.47
1:G:394:LYS:NZ	2:G:512:HOH:O	2.42	0.47
1:F:360:ASP:HB3	1:F:363:LYS:HD2	1.97	0.47
1:A:188:THR:OG1	1:A:368:GLY:HA3	2.15	0.46
1:B:391:GLU:OE2	2:B:503:HOH:O	2.20	0.46
1:G:9:VAL:HA	1:G:298:PRO:HA	1.97	0.46
1:G:176:GLU:OE1	2:G:502:HOH:O	2.20	0.46
1:G:209:LYS:NZ	2:G:520:HOH:O	2.48	0.46
1:F:208:VAL:HG12	1:F:209:LYS:HD2	1.97	0.46
1:H:247:VAL:O	1:H:313:LYS:NZ	2.34	0.46
1:C:120:MET:HE2	1:C:276:ILE:HG23	1.99	0.45
1:E:188:THR:OG1	1:E:368:GLY:HA3	2.16	0.45
1:A:186:GLN:NE2	1:A:225:VAL:O	2.48	0.45
1:C:9:VAL:HA	1:C:298:PRO:HA	1.97	0.45
1:B:263:LYS:HA	1:B:267:ARG:NH1	2.31	0.45
1:G:166:ILE:HD13	1:G:317:VAL:HG13	1.99	0.45
1:A:218:ASP:HB3	2:A:583:HOH:O	2.15	0.45
1:C:58:ILE:HD13	1:C:89:VAL:HA	1.98	0.45
1:F:58:ILE:HD13	1:F:89:VAL:HA	1.98	0.45
1:A:243:SER:CB	1:A:244:PRO:HD3	2.46	0.45
1:F:299:MET:HE3	1:F:393:GLN:HG3	1.99	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:47:PRO:HD2	1:C:48:GLU:OE2	2.17	0.45
1:B:186:GLN:NE2	1:B:225:VAL:O	2.45	0.44
1:H:94:MET:HE1	1:H:404:THR:HA	1.98	0.44
1:B:145:MET:HE1	1:B:276:ILE:HD13	2.00	0.44
1:E:232:LYS:HD3	1:E:232:LYS:HA	1.58	0.44
1:H:120:MET:HE2	1:H:276:ILE:HG23	1.99	0.44
1:G:305:TRP:H	1:G:329:LYS:HD2	1.83	0.44
1:B:136:MET:HG3	1:C:146:ILE:HD11	1.99	0.44
1:B:147:LYS:NZ	2:B:523:HOH:O	2.51	0.44
1:E:231:ALA:HB3	1:E:232:LYS:HE2	1.99	0.44
1:H:397:GLU:O	1:H:399:LYS:HG2	2.18	0.44
1:H:188:THR:OG1	1:H:368:GLY:HA3	2.19	0.43
1:H:192:GLN:HE22	1:H:221:PRO:HG2	1.83	0.43
1:H:397:GLU:HG2	1:H:398:ALA:N	2.33	0.43
1:B:344:GLU:CD	1:B:369:GLY:HA3	2.43	0.43
1:A:215:PHE:HE1	1:A:219:GLU:HG3	1.83	0.43
1:B:22:SER:HB3	1:B:206:ILE:HD13	1.99	0.43
1:E:53:LEU:O	1:E:83:ALA:HA	2.18	0.43
1:H:262:ASP:O	1:H:267:ARG:NH1	2.44	0.43
1:A:274:SER:HB3	1:A:344:GLU:O	2.18	0.43
1:B:9:VAL:HA	1:B:298:PRO:HA	1.99	0.43
1:H:41:LYS:HD3	1:H:41:LYS:HA	1.92	0.43
1:H:180:PHE:CD1	1:H:180:PHE:C	2.97	0.43
1:D:35:VAL:HG23	1:D:282:ALA:HB2	2.01	0.43
1:C:8:VAL:HG13	1:C:283:ILE:HG23	2.00	0.43
1:G:135:ARG:HG2	1:G:135:ARG:NH1	2.34	0.43
1:H:215:PHE:HE1	1:H:219:GLU:HG3	1.84	0.43
1:H:341:GLU:HG3	1:H:388:LEU:HB2	2.01	0.43
1:C:2:ALA:O	1:C:288:GLY:HA3	2.18	0.43
1:G:140:LYS:NZ	2:G:509:HOH:O	2.36	0.43
1:H:220:TYR:N	1:H:221:PRO:HD2	2.34	0.43
1:G:53:LEU:O	1:G:83:ALA:HA	2.19	0.42
1:H:334:ILE:HA	1:H:337:ILE:CD1	2.49	0.42
1:B:398:ALA:O	1:B:419:LYS:HE2	2.19	0.42
1:F:188:THR:OG1	1:F:368:GLY:HA3	2.19	0.42
1:B:207:LYS:HE2	1:B:209:LYS:O	2.20	0.42
1:C:120:MET:HE1	1:C:276:ILE:HD12	2.00	0.42
1:E:9:VAL:HA	1:E:298:PRO:HA	2.02	0.42
1:E:148:ASP:OD2	2:E:503:HOH:O	2.20	0.42
1:A:155:ASN:HB3	1:A:157:TYR:CE2	2.55	0.42
1:A:349:GLN:HE21	1:A:349:GLN:HB2	1.69	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:PHE:HB3	1:B:141:LEU:HD22	2.02	0.42
1:B:301:LYS:HD2	1:B:303:ILE:HG22	2.02	0.42
1:C:2:ALA:HB1	1:C:7:TYR:OH	2.20	0.42
1:C:90:SCY:O	1:C:405:LEU:HD11	2.20	0.42
1:C:151:TRP:CZ2	1:C:157:TYR:HA	2.55	0.42
1:C:169:ILE:HD11	1:C:246:PRO:HB2	2.02	0.42
1:C:242:GLU:OE2	1:C:247:VAL:HG12	2.20	0.41
1:D:164:GLU:CD	1:D:268:VAL:HG23	2.45	0.41
1:E:90:SCY:O	1:E:405:LEU:HD11	2.19	0.41
1:F:3:MET:HA	2:F:553:HOH:O	2.19	0.41
1:G:242:GLU:H	1:G:242:GLU:HG2	1.65	0.41
1:D:218:ASP:HB3	2:D:543:HOH:O	2.20	0.41
1:D:422:GLU:H	1:D:422:GLU:HG3	1.71	0.41
1:F:205:MET:HE3	1:F:205:MET:HB3	1.79	0.41
1:F:241:PRO:O	1:F:242:GLU:HG3	2.20	0.41
1:A:53:LEU:O	1:A:83:ALA:HA	2.20	0.41
1:B:235:GLY:HA3	1:B:266:GLN:HB3	2.02	0.41
1:C:178:ASP:HB3	1:C:230:ILE:HD11	2.02	0.41
1:E:42:ARG:NH1	2:E:511:HOH:O	2.47	0.41
1:H:207:LYS:HE2	1:H:209:LYS:O	2.21	0.41
1:A:243:SER:CB	1:A:244:PRO:CD	2.99	0.41
1:A:319:PRO:HB3	1:A:404:THR:OG1	2.20	0.41
1:A:243:SER:HB2	1:A:244:PRO:CD	2.48	0.41
1:B:243:SER:OG	1:B:264:GLY:HA2	2.21	0.41
1:A:157:TYR:CD1	1:A:161:THR:HB	2.56	0.41
1:H:316:GLY:HA3	1:H:349:GLN:HB3	2.02	0.41
1:C:53:LEU:O	1:C:83:ALA:HA	2.20	0.41
1:D:150:LEU:O	1:D:159:MET:HG2	2.21	0.41
1:D:9:VAL:HA	1:D:298:PRO:HA	2.03	0.41
1:E:337:ILE:HG23	1:E:401:GLY:HA2	2.03	0.41
1:E:388:LEU:HD11	1:E:392:MET:HE3	2.03	0.41
1:A:242:GLU:CB	1:A:246:PRO:HA	2.50	0.40
1:D:296:LEU:C	1:D:297:LYS:HD3	2.46	0.40
1:F:241:PRO:O	1:F:243:SER:N	2.54	0.40
1:G:173:THR:HG22	1:G:261:PRO:HD2	2.02	0.40
1:D:171:GLY:HA3	1:D:258:GLN:O	2.22	0.40
1:E:166:ILE:HD13	1:E:317:VAL:HG13	2.03	0.40
1:A:397:GLU:O	1:A:399:LYS:HG2	2.21	0.40
1:E:209:LYS:HB3	1:E:209:LYS:HE2	1.99	0.40
1:E:329:LYS:HB2	1:E:329:LYS:HE3	1.78	0.40
1:A:419:LYS:HB3	1:A:419:LYS:HE2	1.82	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:241:PRO:C	1:F:243:SER:H	2.29	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:397:GLU:OE2	1:E:395:ARG:NH2[1_565]	2.12	0.08

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	416/428 (97%)	403 (97%)	11 (3%)	2 (0%)	24	16
1	B	417/428 (97%)	402 (96%)	14 (3%)	1 (0%)	43	36
1	C	418/428 (98%)	406 (97%)	11 (3%)	1 (0%)	43	36
1	D	420/428 (98%)	409 (97%)	10 (2%)	1 (0%)	43	36
1	E	418/428 (98%)	403 (96%)	14 (3%)	1 (0%)	43	36
1	F	420/428 (98%)	404 (96%)	14 (3%)	2 (0%)	24	16
1	G	417/428 (97%)	400 (96%)	16 (4%)	1 (0%)	43	36
1	H	414/428 (97%)	401 (97%)	12 (3%)	1 (0%)	43	36
All	All	3340/3424 (98%)	3228 (97%)	102 (3%)	10 (0%)	36	29

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	SER
1	A	89	VAL
1	B	89	VAL
1	C	89	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	89	VAL
1	G	89	VAL
1	H	89	VAL
1	E	89	VAL
1	F	89	VAL
1	F	242	GLU

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	309/321 (96%)	305 (99%)	4 (1%)	61	61
1	B	311/321 (97%)	308 (99%)	3 (1%)	68	70
1	C	312/321 (97%)	308 (99%)	4 (1%)	61	61
1	D	314/321 (98%)	313 (100%)	1 (0%)	86	88
1	E	315/321 (98%)	311 (99%)	4 (1%)	61	61
1	F	316/321 (98%)	313 (99%)	3 (1%)	70	73
1	G	313/321 (98%)	312 (100%)	1 (0%)	86	88
1	H	310/321 (97%)	306 (99%)	4 (1%)	61	61
All	All	2500/2568 (97%)	2476 (99%)	24 (1%)	68	70

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	PHE
1	A	89	VAL
1	A	334	ILE
1	A	349	GLN
1	B	55	PHE
1	B	89	VAL
1	B	217	LYS
1	C	89	VAL
1	C	209	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	349	GLN
1	C	399	LYS
1	D	89	VAL
1	E	55	PHE
1	E	89	VAL
1	E	210	LYS
1	E	293	LYS
1	F	55	PHE
1	F	89	VAL
1	F	232	LYS
1	G	89	VAL
1	H	24	LYS
1	H	89	VAL
1	H	226	THR
1	H	333	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	GLN
1	A	245	ASN
1	A	393	GLN
1	B	63	GLN
1	B	192	GLN
1	B	349	GLN
1	B	367	ASN
1	C	393	GLN
1	D	266	GLN
1	D	349	GLN
1	E	349	GLN
1	F	63	GLN
1	F	349	GLN
1	G	63	GLN
1	G	192	GLN
1	G	250	ASN
1	G	266	GLN
1	G	349	GLN
1	H	250	ASN
1	H	349	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	SCY	F	90	1	7,8,9	1.60	2 (28%)	4,9,11	1.74	1 (25%)
1	SCY	E	90	1	7,8,9	2.12	2 (28%)	4,9,11	2.26	1 (25%)
1	SCY	A	90	1	7,8,9	0.96	0	4,9,11	1.61	1 (25%)
1	SCY	G	90	1	7,8,9	1.74	2 (28%)	4,9,11	1.78	1 (25%)
1	SCY	B	90	1	7,8,9	1.36	2 (28%)	4,9,11	1.98	1 (25%)
1	SCY	D	90	1	7,8,9	1.49	1 (14%)	4,9,11	1.85	1 (25%)
1	SCY	H	90	1	7,8,9	1.24	1 (14%)	4,9,11	1.94	1 (25%)
1	SCY	C	90	1	7,8,9	1.66	1 (14%)	4,9,11	1.99	1 (25%)

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	F	90	1	-	0/5/7/9	-
1	SCY	E	90	1	-	2/5/7/9	-
1	SCY	A	90	1	-	0/5/7/9	-
1	SCY	G	90	1	-	2/5/7/9	-
1	SCY	B	90	1	-	2/5/7/9	-
1	SCY	D	90	1	-	2/5/7/9	-
1	SCY	H	90	1	-	2/5/7/9	-
1	SCY	C	90	1	-	2/5/7/9	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	90	SCY	CB-SG	-4.48	1.71	1.81
1	C	90	SCY	CB-SG	-3.73	1.72	1.81
1	F	90	SCY	CB-SG	-3.36	1.73	1.81
1	G	90	SCY	CB-SG	-3.35	1.73	1.81
1	D	90	SCY	CB-SG	-3.11	1.74	1.81
1	G	90	SCY	O-C	2.47	1.29	1.20
1	B	90	SCY	CB-SG	-2.44	1.75	1.81
1	E	90	SCY	CB-CA	2.27	1.59	1.53
1	H	90	SCY	CB-SG	-2.16	1.76	1.81
1	B	90	SCY	O-C	2.14	1.28	1.20
1	F	90	SCY	O-C	2.09	1.27	1.20

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	90	SCY	CB-CA-C	-3.83	100.41	110.80
1	B	90	SCY	CB-CA-C	-3.81	100.46	110.80
1	E	90	SCY	CB-CA-C	-3.61	101.00	110.80
1	D	90	SCY	CB-CA-C	-3.57	101.10	110.80
1	G	90	SCY	CB-CA-C	-3.52	101.26	110.80
1	H	90	SCY	CB-CA-C	-3.48	101.34	110.80
1	F	90	SCY	CB-CA-C	-3.32	101.77	110.80
1	A	90	SCY	CB-CA-C	-3.11	102.36	110.80

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	90	SCY	OCD-CD-SG-CB
1	C	90	SCY	CE-CD-SG-CB
1	D	90	SCY	OCD-CD-SG-CB
1	D	90	SCY	CE-CD-SG-CB
1	E	90	SCY	OCD-CD-SG-CB
1	G	90	SCY	OCD-CD-SG-CB
1	G	90	SCY	CE-CD-SG-CB
1	H	90	SCY	OCD-CD-SG-CB
1	H	90	SCY	CE-CD-SG-CB
1	E	90	SCY	CE-CD-SG-CB
1	B	90	SCY	OCD-CD-SG-CB
1	B	90	SCY	CE-CD-SG-CB

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	E	90	SCY	1	0
1	A	90	SCY	1	0
1	H	90	SCY	1	0
1	C	90	SCY	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	418/428 (97%)	0.49	10 (2%) 59 63	35, 51, 80, 98	0
1	B	419/428 (97%)	0.29	3 (0%) 84 86	36, 47, 67, 83	0
1	C	420/428 (98%)	0.26	5 (1%) 76 79	33, 44, 68, 83	0
1	D	422/428 (98%)	0.18	4 (0%) 81 83	33, 43, 60, 90	0
1	E	420/428 (98%)	0.12	4 (0%) 79 82	31, 41, 59, 103	0
1	F	422/428 (98%)	0.20	6 (1%) 73 76	30, 44, 70, 87	0
1	G	419/428 (97%)	0.27	3 (0%) 84 86	34, 45, 60, 84	0
1	H	416/428 (97%)	0.95	37 (8%) 15 16	36, 68, 98, 113	0
All	All	3356/3424 (98%)	0.34	72 (2%) 63 67	30, 47, 79, 113	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	421	LEU	5.3
1	H	262	ASP	4.8
1	E	3	MET	4.6
1	D	425	HIS	4.2
1	A	224	GLY	3.7
1	H	373	LEU	3.2
1	G	421	LEU	3.1
1	G	2	ALA	3.1
1	H	261	PRO	3.0
1	H	228	ASP	3.0
1	D	421	LEU	3.0
1	F	421	LEU	3.0
1	A	225	VAL	2.9
1	C	422	GLU	2.9
1	B	421	LEU	2.9
1	B	224	GLY	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	423	HIS	2.7
1	F	3	MET	2.7
1	A	208	VAL	2.7
1	H	194	ALA	2.6
1	H	180	PHE	2.6
1	E	421	LEU	2.6
1	H	299	MET	2.6
1	C	2	ALA	2.6
1	H	75	ALA	2.6
1	D	208	VAL	2.5
1	H	214	PRO	2.5
1	D	3	MET	2.4
1	H	337	ILE	2.4
1	H	227	ALA	2.4
1	H	175	ARG	2.4
1	C	263	LYS	2.4
1	H	342	ALA	2.4
1	F	208	VAL	2.4
1	A	249	VAL	2.3
1	G	247	VAL	2.3
1	H	388	LEU	2.3
1	A	3	MET	2.3
1	H	273	ALA	2.3
1	H	341	GLU	2.3
1	H	225	VAL	2.3
1	H	366	VAL	2.3
1	H	420	CYS	2.3
1	F	241	PRO	2.2
1	F	352	ALA	2.2
1	H	351	ILE	2.2
1	H	4	LYS	2.2
1	A	170	TRP	2.2
1	H	197	PHE	2.2
1	F	232	LYS	2.2
1	A	223	ALA	2.2
1	H	352	ALA	2.2
1	C	334	ILE	2.2
1	H	169	ILE	2.2
1	H	153	ALA	2.2
1	H	290	ALA	2.2
1	H	173	THR	2.1
1	H	274	SER	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	223	ALA	2.1
1	H	340	ILE	2.1
1	A	197	PHE	2.1
1	C	199	ASP	2.1
1	H	170	TRP	2.1
1	H	284	VAL	2.1
1	H	393	GLN	2.1
1	H	339	LEU	2.1
1	H	347	ALA	2.1
1	H	208	VAL	2.0
1	H	105	LEU	2.0
1	H	260	ALA	2.0
1	A	4	LYS	2.0
1	E	4	LYS	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	E	90	9/10	0.87	0.13	33,38,46,71	0
1	SCY	G	90	9/10	0.87	0.12	35,41,51,58	0
1	SCY	H	90	9/10	0.87	0.14	49,57,65,74	0
1	SCY	C	90	9/10	0.88	0.13	36,42,48,67	0
1	SCY	A	90	9/10	0.89	0.13	41,46,53,69	0
1	SCY	D	90	9/10	0.90	0.12	35,37,47,61	0
1	SCY	B	90	9/10	0.90	0.11	39,42,48,64	0
1	SCY	F	90	9/10	0.92	0.11	36,41,51,59	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.