



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:30 PM UTC

PDB ID : 3M81 / pdb_00003m81
Title : Crystal structure of Acetyl xylan esterase (TM0077) from THERMOTOGA MARITIMA at 2.50 Å resolution (native apo structure)
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2010-03-17
Resolution : 2.50 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:

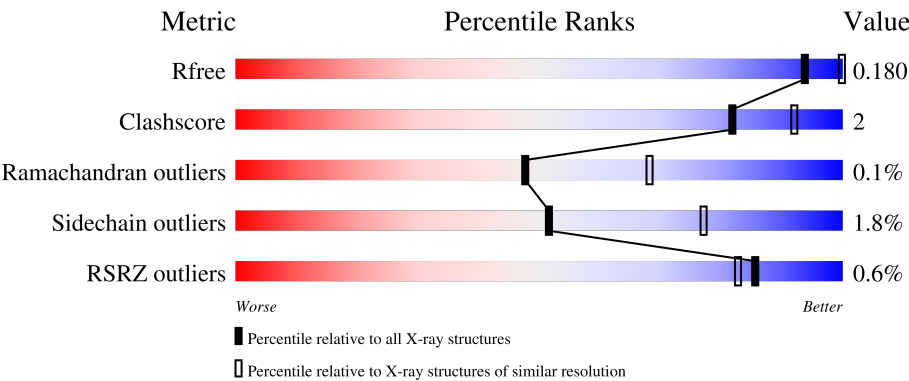
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	337	91%5%.
1	B	337	% 90%6%.
1	C	337	% 89%6%.
1	D	337	% 87%9%.
1	E	337	86%9%5%

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Mol	Chain	Length	Quality of chain
1	F	337	 87% 8%

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
3	CL	D	412	-	-	X	-
4	ACT	D	416	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 16183 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 xylan esterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	1	0
			2605	1694	433	468	10			
1	B	322	Total	C	N	O	S	0	1	0
			2604	1694	432	468	10			
1	C	322	Total	C	N	O	S	0	1	0
			2603	1693	432	468	10			
1	D	322	Total	C	N	O	S	0	1	0
			2604	1694	431	469	10			
1	E	321	Total	C	N	O	S	0	2	0
			2607	1695	433	469	10			
1	F	322	Total	C	N	O	S	0	0	0
			2596	1689	431	466	10			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP Q9WXT2
A	-10	GLY	-	expression tag	UNP Q9WXT2
A	-9	SER	-	expression tag	UNP Q9WXT2
A	-8	ASP	-	expression tag	UNP Q9WXT2
A	-7	LYS	-	expression tag	UNP Q9WXT2
A	-6	ILE	-	expression tag	UNP Q9WXT2
A	-5	HIS	-	expression tag	UNP Q9WXT2
A	-4	HIS	-	expression tag	UNP Q9WXT2
A	-3	HIS	-	expression tag	UNP Q9WXT2
A	-2	HIS	-	expression tag	UNP Q9WXT2
A	-1	HIS	-	expression tag	UNP Q9WXT2
A	0	HIS	-	expression tag	UNP Q9WXT2
B	-11	MET	-	expression tag	UNP Q9WXT2
B	-10	GLY	-	expression tag	UNP Q9WXT2
B	-9	SER	-	expression tag	UNP Q9WXT2
B	-8	ASP	-	expression tag	UNP Q9WXT2
B	-7	LYS	-	expression tag	UNP Q9WXT2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	ILE	-	expression tag	UNP Q9WXT2
B	-5	HIS	-	expression tag	UNP Q9WXT2
B	-4	HIS	-	expression tag	UNP Q9WXT2
B	-3	HIS	-	expression tag	UNP Q9WXT2
B	-2	HIS	-	expression tag	UNP Q9WXT2
B	-1	HIS	-	expression tag	UNP Q9WXT2
B	0	HIS	-	expression tag	UNP Q9WXT2
C	-11	MET	-	expression tag	UNP Q9WXT2
C	-10	GLY	-	expression tag	UNP Q9WXT2
C	-9	SER	-	expression tag	UNP Q9WXT2
C	-8	ASP	-	expression tag	UNP Q9WXT2
C	-7	LYS	-	expression tag	UNP Q9WXT2
C	-6	ILE	-	expression tag	UNP Q9WXT2
C	-5	HIS	-	expression tag	UNP Q9WXT2
C	-4	HIS	-	expression tag	UNP Q9WXT2
C	-3	HIS	-	expression tag	UNP Q9WXT2
C	-2	HIS	-	expression tag	UNP Q9WXT2
C	-1	HIS	-	expression tag	UNP Q9WXT2
C	0	HIS	-	expression tag	UNP Q9WXT2
D	-11	MET	-	expression tag	UNP Q9WXT2
D	-10	GLY	-	expression tag	UNP Q9WXT2
D	-9	SER	-	expression tag	UNP Q9WXT2
D	-8	ASP	-	expression tag	UNP Q9WXT2
D	-7	LYS	-	expression tag	UNP Q9WXT2
D	-6	ILE	-	expression tag	UNP Q9WXT2
D	-5	HIS	-	expression tag	UNP Q9WXT2
D	-4	HIS	-	expression tag	UNP Q9WXT2
D	-3	HIS	-	expression tag	UNP Q9WXT2
D	-2	HIS	-	expression tag	UNP Q9WXT2
D	-1	HIS	-	expression tag	UNP Q9WXT2
D	0	HIS	-	expression tag	UNP Q9WXT2
E	-11	MET	-	expression tag	UNP Q9WXT2
E	-10	GLY	-	expression tag	UNP Q9WXT2
E	-9	SER	-	expression tag	UNP Q9WXT2
E	-8	ASP	-	expression tag	UNP Q9WXT2
E	-7	LYS	-	expression tag	UNP Q9WXT2
E	-6	ILE	-	expression tag	UNP Q9WXT2
E	-5	HIS	-	expression tag	UNP Q9WXT2
E	-4	HIS	-	expression tag	UNP Q9WXT2
E	-3	HIS	-	expression tag	UNP Q9WXT2
E	-2	HIS	-	expression tag	UNP Q9WXT2
E	-1	HIS	-	expression tag	UNP Q9WXT2

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Chain	Residue	Modelled	Actual	Comment	Reference
E	0	HIS	-	expression tag	UNP Q9WXT2
F	-11	MET	-	expression tag	UNP Q9WXT2
F	-10	GLY	-	expression tag	UNP Q9WXT2
F	-9	SER	-	expression tag	UNP Q9WXT2
F	-8	ASP	-	expression tag	UNP Q9WXT2
F	-7	LYS	-	expression tag	UNP Q9WXT2
F	-6	ILE	-	expression tag	UNP Q9WXT2
F	-5	HIS	-	expression tag	UNP Q9WXT2
F	-4	HIS	-	expression tag	UNP Q9WXT2
F	-3	HIS	-	expression tag	UNP Q9WXT2
F	-2	HIS	-	expression tag	UNP Q9WXT2
F	-1	HIS	-	expression tag	UNP Q9WXT2
F	0	HIS	-	expression tag	UNP Q9WXT2

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Ca 2 2	0	0
2	B	1	Total Ca 1 1	0	0
2	C	1	Total Ca 1 1	0	0
2	D	2	Total Ca 2 2	0	0
2	E	1	Total Ca 1 1	0	0
2	F	1	Total Ca 1 1	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

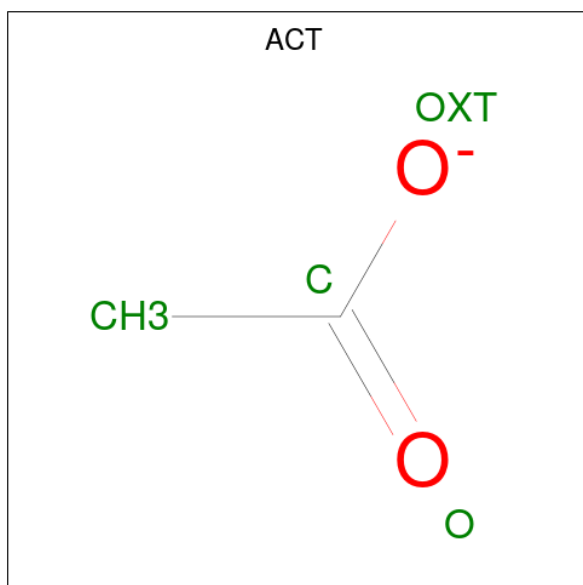
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0
3	B	1	Total Cl 1 1	0	0
3	C	1	Total Cl 1 1	0	0
3	D	1	Total Cl 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	1	Total	Cl	0	0
			1	1		
3	F	1	Total	Cl	0	0
			1	1		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $\text{C}_2\text{H}_3\text{O}_2^-$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		
5	E	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	98	Total	O	0	1
			99	99		
6	B	99	Total	O	0	0
			99	99		
6	C	88	Total	O	0	2
			88	88		
6	D	86	Total	O	0	2
			88	88		
6	E	64	Total	O	0	0
			64	64		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	F	72	Total	O	0	0
			72	72		

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 xylan esterase

Chain A: 



- Molecule 1: Acetyl xylan esterase

Chain B: 



- Molecule 1: Acetyl xylan esterase

Chain C: 



- Molecule 1: Acetyl xylan esterase

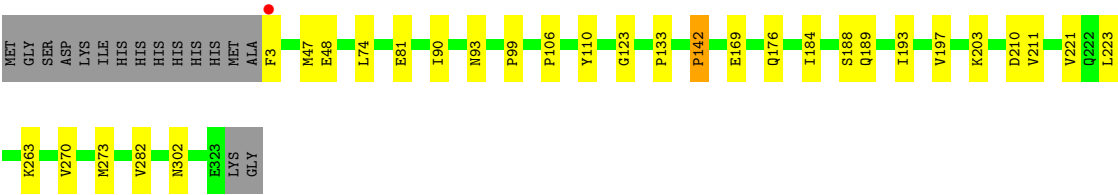
Chain D: 



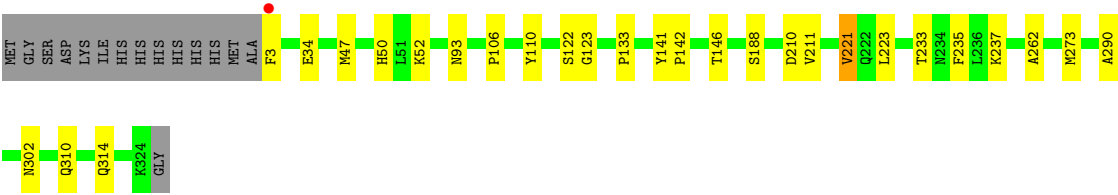
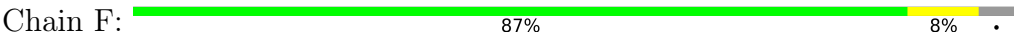
- Molecule 1: Acetyl xylan esterase

Chain E: 





● Molecule 1: Acetyl xylan esterase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	103.46Å 103.79Å 221.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.80 – 2.50 48.80 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.80-2.50) 100.0 (48.80-2.50)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019, PHENIX	Depositor
R, R_{free}	0.167 , 0.212 (Not available) , 0.180	Depositor DCC
R_{free} test set	4200 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	32.9	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 34.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.020 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16183	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.45% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, CL, EDO, CA

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	0.94	0/2685	0.89	2/3641 (0.1%)
1	B	0.97	0/2684	0.89	0/3640
1	C	0.94	0/2683	0.91	1/3639 (0.0%)
1	D	0.94	1/2684 (0.0%)	0.91	0/3640
1	E	0.95	0/2687	0.91	0/3644
1	F	0.93	0/2676	0.91	0/3629
All	All	0.95	1/16099 (0.0%)	0.90	3/21833 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	41	ASP	C-O	-5.66	1.21	1.23

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	132	TYR	CA-C-N	-5.51	114.91	120.31
1	A	132	TYR	C-N-CA	-5.51	114.91	120.31
1	C	188	SER	CB-CA-C	-5.46	110.30	116.63

There are no chirality outliers.

There are no planarity outliers.

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	2605	0	2514	8	0
1	B	2604	0	2509	9	0
1	C	2603	0	2507	11	0
1	D	2604	0	2507	22	0
1	E	2607	0	2512	14	0
1	F	2596	0	2507	17	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	1	0	0	1	0
3	B	1	0	0	1	0
3	C	1	0	0	1	0
3	D	1	0	0	2	0
3	E	1	0	0	1	0
3	F	1	0	0	1	0
4	A	4	0	3	0	0
4	D	4	0	3	3	0
4	E	4	0	3	1	0
5	A	8	0	12	1	0
5	B	4	0	6	0	0
5	C	4	0	6	0	0
5	D	4	0	6	3	0
5	E	4	0	6	0	0
5	F	4	0	6	0	0
6	A	99	0	0	0	0
6	B	99	0	0	0	0
6	C	88	0	0	0	0
6	D	88	0	0	3	0
6	E	64	0	0	0	0
6	F	72	0	0	0	0
All	All	16183	0	15107	77	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 2.

All (77) 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:D:185:ALA:HB1	5:D:420:EDO:H22	1.63	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:221:VAL:HG22	1:F:233:THR:OG1	1.88	0.73
1:D:185:ALA:HB1	5:D:420:EDO:C2	2.19	0.72
1:E:81:GLU:O	1:E:176[A]:GLN:HG3	1.96	0.65
1:C:188:SER:HB3	3:C:411:CL:CL	2.34	0.64
1:E:184:ILE:HD13	1:E:197:VAL:HG12	1.82	0.62
1:B:142:PRO:HA	4:D:416:ACT:H3	1.82	0.61
1:C:47:MET:HA	1:C:47:MET:HE2	1.83	0.60
1:E:93:ASN:HB3	1:E:123:GLY:HA3	1.84	0.59
1:D:120:GLN:HG2	1:D:144:PHE:CE2	2.39	0.56
1:A:64:TYR:OH	1:A:134:GLU:OE2	2.21	0.55
1:B:188:SER:HB3	3:B:410:CL:CL	2.45	0.54
1:C:53:THR:OG1	1:C:54:VAL:HG23	2.08	0.54
1:F:93:ASN:HB3	1:F:123:GLY:HA3	1.89	0.54
1:C:302[B]:ASN:CG	1:F:237:LYS:HZ2	2.16	0.53
1:D:188:SER:HB3	3:D:412:CL:CL	2.45	0.53
1:E:169:GLU:OE1	1:E:203:LYS:NZ	2.41	0.53
1:C:184:ILE:HD13	1:C:197:VAL:HG12	1.90	0.53
1:E:210:ASP:O	1:E:211:VAL:C	2.52	0.52
1:F:146:THR:HG21	1:F:235:PHE:HB2	1.91	0.52
1:A:93:ASN:HB3	1:A:123:GLY:HA3	1.92	0.52
1:E:142:PRO:HA	4:E:417:ACT:H1	1.92	0.52
1:F:188:SER:HB3	3:F:413:CL:CL	2.48	0.51
1:A:210:ASP:O	1:A:211:VAL:C	2.54	0.50
1:C:93:ASN:HB3	1:C:123:GLY:HA3	1.94	0.50
1:B:210:ASP:O	1:B:211:VAL:C	2.55	0.49
1:D:93:ASN:HB3	1:D:123:GLY:HA3	1.93	0.49
1:D:302:ASN:OD1	1:D:303:HIS:N	2.45	0.49
1:B:47:MET:HA	1:B:47:MET:HE2	1.95	0.49
1:F:310:GLN:O	1:F:314:GLN:HG3	2.13	0.48
1:B:93:ASN:HB3	1:B:123:GLY:HA3	1.94	0.48
1:D:120:GLN:HG2	1:D:144:PHE:CZ	2.49	0.48
1:E:3:PHE:CE1	1:E:273:MET:HE3	2.48	0.48
1:A:237:LYS:HZ2	1:E:302[B]:ASN:CG	2.22	0.47
1:D:3:PHE:CD1	1:D:273:MET:HE3	2.50	0.47
1:D:73:LEU:C	1:D:74:LEU:HD23	2.41	0.46
1:C:3:PHE:CE1	1:C:273:MET:HE3	2.50	0.46
1:F:3:PHE:HD1	1:F:273:MET:HE2	1.81	0.46
1:B:262:ALA:HB3	1:B:289:TYR:CE1	2.51	0.45
1:C:120:GLN:HG2	1:C:144:PHE:CE2	2.51	0.45
1:D:143:GLY:CA	4:D:416:ACT:H2	2.46	0.45
1:D:106:PRO:HA	1:D:110:TYR:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:VAL:HG12	6:D:507:HOH:O	2.16	0.45
1:E:106:PRO:HA	1:E:110:TYR:O	2.17	0.45
1:F:210:ASP:O	1:F:211:VAL:C	2.59	0.45
1:E:189:GLN:O	1:E:193:ILE:HG13	2.17	0.45
1:F:262:ALA:O	1:F:290:ALA:HB3	2.16	0.44
1:F:3:PHE:CE1	1:F:273:MET:HE3	2.53	0.44
1:E:188:SER:HB3	3:E:414:CL:CL	2.55	0.44
1:E:270:VAL:HB	1:E:282:VAL:HG11	2.00	0.43
1:A:218:ARG:HD2	1:A:222[B]:GLN:HE22	1.84	0.43
1:D:315:VAL:HG11	1:F:50:HIS:CB	2.48	0.43
1:D:15:ARG:NE	6:D:1589:HOH:O	2.47	0.43
1:A:16:PRO:O	1:A:288:TYR:OH	2.36	0.42
1:F:106:PRO:HA	1:F:110:TYR:O	2.20	0.42
1:A:185:ALA:HB1	5:A:423:EDO:H22	2.01	0.42
1:F:211:VAL:HG23	1:F:211:VAL:O	2.20	0.42
1:C:120:GLN:HG2	1:C:144:PHE:CZ	2.54	0.42
1:D:210:ASP:O	1:D:211:VAL:C	2.61	0.42
1:B:183:VAL:HG21	1:B:318:LEU:HD13	2.00	0.41
1:B:113:PHE:C	1:B:113:PHE:CD1	2.98	0.41
1:F:3:PHE:CD1	1:F:273:MET:CE	3.03	0.41
1:D:143:GLY:C	4:D:416:ACT:H2	2.45	0.41
1:E:133:PRO:HG3	1:F:133:PRO:HG3	2.02	0.41
1:C:210:ASP:O	1:C:211:VAL:C	2.63	0.41
1:D:16:PRO:O	1:D:288:TYR:OH	2.37	0.41
1:E:74:LEU:HD21	1:E:99:PRO:HB3	2.02	0.41
1:A:188:SER:HB3	3:A:409:CL:CL	2.58	0.41
1:D:151:ASP:OD1	1:D:152:PRO:HD2	2.20	0.41
1:F:3:PHE:HD1	1:F:273:MET:CE	2.33	0.41
1:D:166:ARG:NH2	6:D:509:HOH:O	2.54	0.41
1:D:185:ALA:HB1	5:D:420:EDO:H21	2.00	0.41
1:F:141:TYR:CG	1:F:142:PRO:HD2	2.56	0.41
1:C:106:PRO:HA	1:C:110:TYR:O	2.21	0.41
1:B:278:PRO:HA	1:B:279:PRO:HD3	1.96	0.40
1:D:92:TYR:HB2	3:D:412:CL:CL	2.58	0.40
1:D:74:LEU:HD11	1:D:99:PRO:O	2.21	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	321/337 (95%)	313 (98%)	8 (2%)	0	100	100
1	B	321/337 (95%)	310 (97%)	9 (3%)	2 (1%)	21	38
1	C	321/337 (95%)	312 (97%)	9 (3%)	0	100	100
1	D	321/337 (95%)	314 (98%)	7 (2%)	0	100	100
1	E	321/337 (95%)	313 (98%)	8 (2%)	0	100	100
1	F	320/337 (95%)	311 (97%)	9 (3%)	0	100	100
All	All	1925/2022 (95%)	1873 (97%)	50 (3%)	2 (0%)	48	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	302[A]	ASN
1	B	302[B]	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	269/285 (94%)	266 (99%)	3 (1%)	65	84
1	B	268/285 (94%)	264 (98%)	4 (2%)	57	80
1	C	268/285 (94%)	264 (98%)	4 (2%)	57	80
1	D	268/285 (94%)	265 (99%)	3 (1%)	65	84
1	E	269/285 (94%)	262 (97%)	7 (3%)	40	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	F	232/285 (81%)	225 (97%)	7 (3%)	36 64
All	All	1574/1710 (92%)	1546 (98%)	28 (2%)	51 77

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLU
1	A	47	MET
1	A	221	VAL
1	B	48	GLU
1	B	52	LYS
1	B	221	VAL
1	B	223	LEU
1	C	46	ARG
1	C	74	LEU
1	C	221	VAL
1	C	223	LEU
1	D	74	LEU
1	D	221	VAL
1	D	223	LEU
1	E	47	MET
1	E	48	GLU
1	E	90	ILE
1	E	142	PRO
1	E	221	VAL
1	E	223	LEU
1	E	263	LYS
1	F	34	GLU
1	F	47	MET
1	F	52	LYS
1	F	122	SER
1	F	221	VAL
1	F	223	LEU
1	F	302	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	GLN
1	B	222	GLN
1	C	88	GLN

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Mol	Chain	Res	Type
1	C	140	GLN
1	C	222	GLN
1	D	50	HIS
1	E	50	HIS
1	E	140	GLN
1	F	50	HIS
1	F	100	HIS
1	F	140	GLN
1	F	176	GLN
1	F	222	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 24 ligands modelled in this entry, 14 are monoatomic - leaving 10 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	ACT	E	417	-	3,3,3	0.95	0	3,3,3	1.80	1 (33%)
5	EDO	E	421	-	3,3,3	0.68	0	2,2,2	0.02	0
5	EDO	F	422	-	3,3,3	0.51	0	2,2,2	0.04	0
5	EDO	C	419	-	3,3,3	0.56	0	2,2,2	0.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	D	420	-	3,3,3	0.33	0	2,2,2	0.42	0
4	ACT	A	415	-	3,3,3	1.05	0	3,3,3	1.92	2 (66%)
5	EDO	A	424	-	3,3,3	0.49	0	2,2,2	0.43	0
5	EDO	A	423	-	3,3,3	0.52	0	2,2,2	0.20	0
4	ACT	D	416	-	3,3,3	0.90	0	3,3,3	0.99	0
5	EDO	B	418	-	3,3,3	0.69	0	2,2,2	0.07	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
5	EDO	E	421	-	-	0/1/1/1	-
5	EDO	F	422	-	-	1/1/1/1	-
5	EDO	C	419	-	-	1/1/1/1	-
5	EDO	D	420	-	-	1/1/1/1	-
5	EDO	A	424	-	-	1/1/1/1	-
5	EDO	A	423	-	-	0/1/1/1	-
5	EDO	B	418	-	-	0/1/1/1	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	415	ACT	OXT-C-O	-2.58	112.47	122.03
4	E	417	ACT	OXT-C-O	-2.48	112.84	122.03
4	A	415	ACT	OXT-C-CH3	2.08	123.77	115.05

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	F	422	EDO	O1-C1-C2-O2
5	A	424	EDO	O1-C1-C2-O2
5	C	419	EDO	O1-C1-C2-O2
5	D	420	EDO	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	417	ACT	1	0
5	D	420	EDO	3	0
5	A	423	EDO	1	0
4	D	416	ACT	3	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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(Å²)	Q<0.9	
1	A	322/337 (95%)	-0.65	1 (0%)	90	87	15, 26, 44, 59	1 (0%)
1	B	322/337 (95%)	-0.62	2 (0%)	85	83	11, 26, 43, 60	1 (0%)
1	C	322/337 (95%)	-0.63	4 (1%)	76	73	12, 26, 44, 60	1 (0%)
1	D	322/337 (95%)	-0.61	2 (0%)	85	83	14, 26, 44, 60	1 (0%)
1	E	321/337 (95%)	-0.54	1 (0%)	90	87	12, 27, 44, 60	2 (0%)
1	F	322/337 (95%)	-0.56	1 (0%)	90	87	20, 27, 44, 60	0
All	All	1931/2022 (95%)	-0.60	11 (0%)	85	83	11, 26, 44, 60	6 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	3	PHE	3.1
1	A	3	PHE	3.0
1	C	3	PHE	2.9
1	D	3	PHE	2.8
1	E	3	PHE	2.5
1	B	79	GLU	2.5
1	B	3	PHE	2.4
1	C	135	GLY	2.3
1	C	324	LYS	2.2
1	D	324	LYS	2.1
1	C	134	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides 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. 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
2	CA	F	406	1/1	0.83	0.13	72,72,72,72	0
2	CA	A	404	1/1	0.84	0.10	69,69,69,69	0
2	CA	D	405	1/1	0.85	0.09	66,66,66,66	0
2	CA	E	407	1/1	0.87	0.12	65,65,65,65	0
3	CL	E	414	1/1	0.87	0.11	46,46,46,46	0
4	ACT	E	417	4/4	0.88	0.15	46,49,49,49	0
2	CA	C	403	1/1	0.90	0.06	74,74,74,74	0
2	CA	B	408	1/1	0.92	0.06	64,64,64,64	0
3	CL	D	412	1/1	0.92	0.09	44,44,44,44	0
4	ACT	A	415	4/4	0.93	0.10	32,37,37,37	0
3	CL	F	413	1/1	0.93	0.12	47,47,47,47	0
3	CL	C	411	1/1	0.94	0.07	40,40,40,40	0
4	ACT	D	416	4/4	0.94	0.09	32,32,34,35	0
3	CL	A	409	1/1	0.94	0.08	47,47,47,47	0
5	EDO	A	424	4/4	0.94	0.10	34,36,40,43	0
5	EDO	C	419	4/4	0.94	0.09	35,38,39,41	0
3	CL	B	410	1/1	0.95	0.07	42,42,42,42	0
2	CA	D	402	1/1	0.95	0.04	30,30,30,30	0
5	EDO	D	420	4/4	0.95	0.10	38,42,48,50	0
5	EDO	A	423	4/4	0.96	0.09	30,34,34,35	0
5	EDO	B	418	4/4	0.96	0.08	25,29,34,37	0
5	EDO	F	422	4/4	0.96	0.10	29,44,48,54	0
5	EDO	E	421	4/4	0.97	0.09	27,41,41,44	0
2	CA	A	401	1/1	0.97	0.03	25,25,25,25	0

6.5 Other polymers ⓘ

There are no such residues in this entry.