



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2022 – 05:28 PM EDT

PDB ID : 7KLI
Title : Crystal Structure of Enoyl-[acyl-carrier-protein] reductase [NADH] (InhA) from Mycobacterium abscessus
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2020-10-30
Resolution : 1.75 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

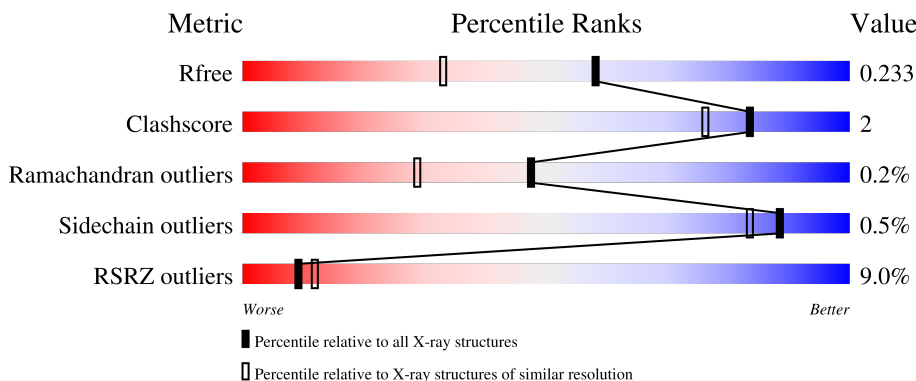
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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}	130704	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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	277	<div> <div>6%</div> <div> <div></div> <div>87%</div> <div>7%</div> <div>6%</div> </div> </div>
1	B	277	<div> <div>7%</div> <div> <div></div> <div>88%</div> <div>5%</div> <div>7%</div> </div> </div>
1	C	277	<div> <div>10%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div>.</div> </div> </div>
1	D	277	<div> <div>10%</div> <div> <div></div> <div>92%</div> <div>5%</div> <div>.</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8985 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 Enoyl-[acyl-carrier-protein] reductase [NADH].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	11	0
			1984	1267	338	370	9			
1	B	257	Total	C	N	O	S	0	8	0
			1962	1255	336	362	9			
1	C	267	Total	C	N	O	S	0	8	0
			1988	1275	335	367	11			
1	D	267	Total	C	N	O	S	0	7	0
			1985	1272	334	369	10			

There are 32 discrepancies between the modelled and reference sequences:

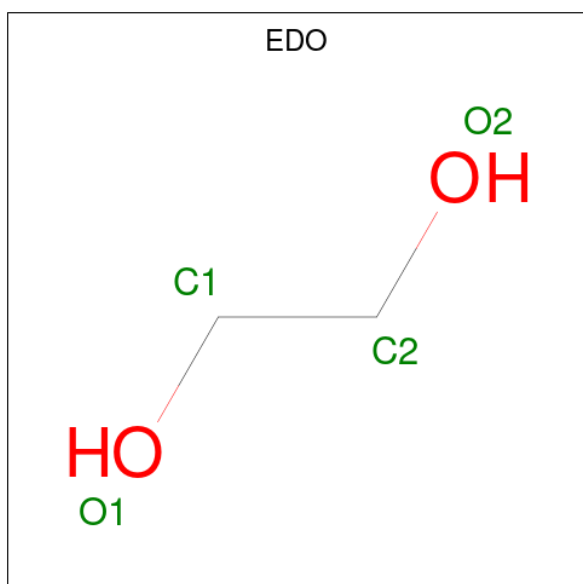
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP B1MC30
A	-6	ALA	-	expression tag	UNP B1MC30
A	-5	HIS	-	expression tag	UNP B1MC30
A	-4	HIS	-	expression tag	UNP B1MC30
A	-3	HIS	-	expression tag	UNP B1MC30
A	-2	HIS	-	expression tag	UNP B1MC30
A	-1	HIS	-	expression tag	UNP B1MC30
A	0	HIS	-	expression tag	UNP B1MC30
B	-7	MET	-	initiating methionine	UNP B1MC30
B	-6	ALA	-	expression tag	UNP B1MC30
B	-5	HIS	-	expression tag	UNP B1MC30
B	-4	HIS	-	expression tag	UNP B1MC30
B	-3	HIS	-	expression tag	UNP B1MC30
B	-2	HIS	-	expression tag	UNP B1MC30
B	-1	HIS	-	expression tag	UNP B1MC30
B	0	HIS	-	expression tag	UNP B1MC30
C	-7	MET	-	initiating methionine	UNP B1MC30
C	-6	ALA	-	expression tag	UNP B1MC30
C	-5	HIS	-	expression tag	UNP B1MC30
C	-4	HIS	-	expression tag	UNP B1MC30
C	-3	HIS	-	expression tag	UNP B1MC30

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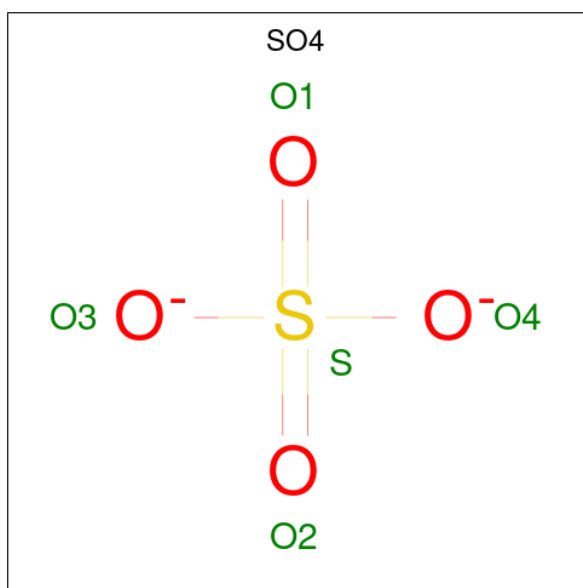
Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP B1MC30
C	-1	HIS	-	expression tag	UNP B1MC30
C	0	HIS	-	expression tag	UNP B1MC30
D	-7	MET	-	initiating methionine	UNP B1MC30
D	-6	ALA	-	expression tag	UNP B1MC30
D	-5	HIS	-	expression tag	UNP B1MC30
D	-4	HIS	-	expression tag	UNP B1MC30
D	-3	HIS	-	expression tag	UNP B1MC30
D	-2	HIS	-	expression tag	UNP B1MC30
D	-1	HIS	-	expression tag	UNP B1MC30
D	0	HIS	-	expression tag	UNP B1MC30

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	245	Total	O	0	15
			260	260		
4	B	241	Total	O	0	13
			254	254		
4	C	226	Total	O	0	13
			239	239		

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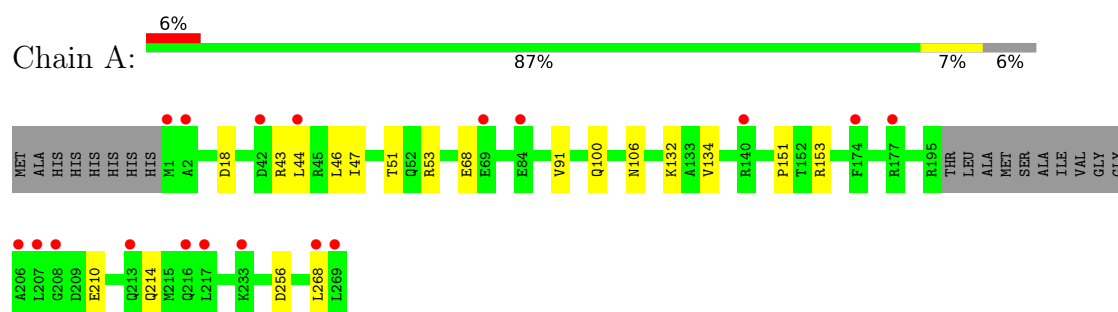
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	239	Total	O	0	10
			249	249		

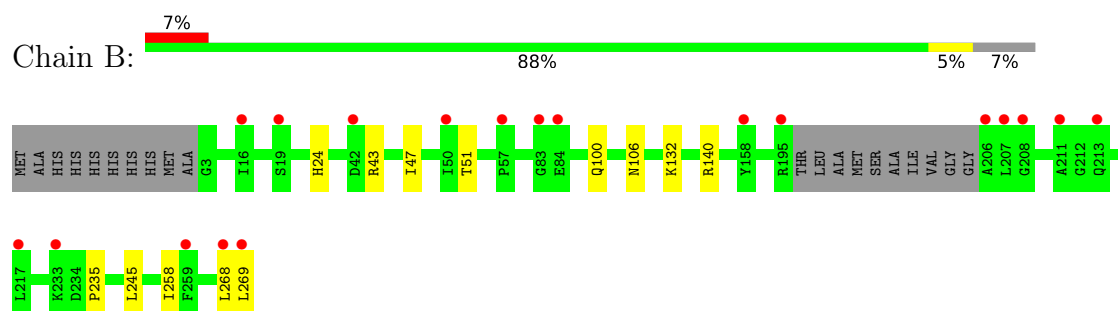
3 Residue-property plots [i](#)

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.

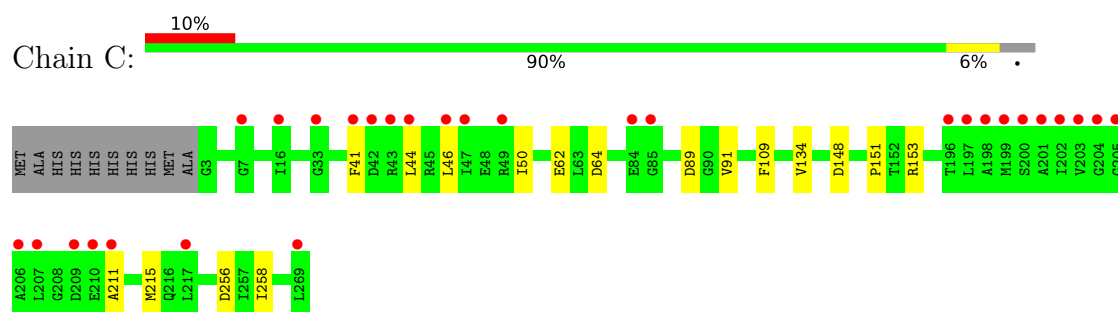
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



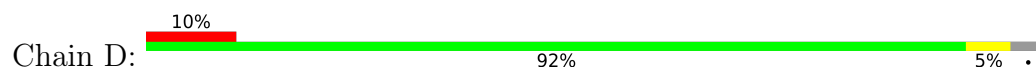
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

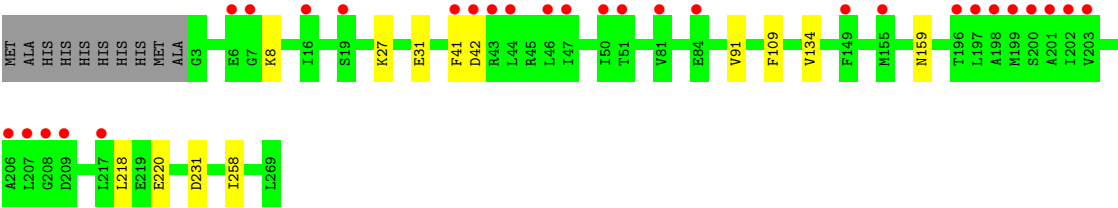


- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.84Å 98.03Å 147.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.14 – 1.75 47.14 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.14-1.75) 99.7 (47.14-1.75)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.72 (at 1.75Å)	Xtriage
Refinement program	PHENIX 1.19.2	Depositor
R, R_{free}	0.206 , 0.233 0.206 , 0.233	Depositor DCC
R_{free} test set	1983 reflections (1.73%)	wwPDB-VP
Wilson B-factor (Å ²)	12.9	Xtriage
Anisotropy	0.591	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 53.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.56$, $\langle L^2 \rangle = 0.41$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8985	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 56.12 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.8640e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, EDO

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	0.41	0/2048	0.63	0/2774
1	B	0.43	0/2017	0.61	0/2732
1	C	0.41	0/2052	0.60	2/2789 (0.1%)
1	D	0.41	0/2046	0.61	0/2779
All	All	0.41	0/8163	0.61	2/11074 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	148	ASP	CB-CG-OD1	5.06	122.85	118.30
1	C	148	ASP	CB-CG-OD2	-5.04	113.76	118.30

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	1984	0	2009	13	0
1	B	1962	0	1986	10	0
1	C	1988	0	1995	10	0
1	D	1985	0	1983	10	0
2	A	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	4	0	6	0	0
2	C	8	0	12	0	0
2	D	8	0	12	0	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	1	0
4	A	260	0	0	2	1
4	B	254	0	0	2	2
4	C	239	0	0	1	1
4	D	249	0	0	5	0
All	All	8985	0	8009	38	2

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 (38) 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:153[B]:ARG:NH2	1:C:153[B]:ARG:HH12	1.93	0.67
1:B:269[B]:LEU:HD11	1:D:218:LEU:HA	1.84	0.59
1:D:220:GLU:HG3	4:D:466:HOH:O	2.03	0.59
1:A:18:ASP:OD2	1:A:53:ARG:NH1	2.35	0.59
1:C:258:ILE:HD12	1:D:258:ILE:HD12	1.85	0.58
1:D:27:LYS:NZ	1:D:31:GLU:OE2	2.36	0.58
1:B:268[B]:LEU:O	1:B:269[B]:LEU:HD23	2.05	0.56
1:A:68:GLU:OE2	4:A:401:HOH:O	2.18	0.56
1:B:100:GLN:NE2	4:B:406:HOH:O	2.35	0.55
1:C:41[B]:PHE:CD1	1:C:64:ASP:HB2	2.43	0.54
3:D:304:SO4:O4	4:D:401:HOH:O	2.18	0.53
1:A:210:GLU:HG3	1:A:214:GLN:HE21	1.73	0.53
1:C:91:VAL:HG21	1:C:134:VAL:HG21	1.90	0.53
1:C:89:ASP:OD1	4:C:401:HOH:O	2.19	0.52
1:A:91:VAL:HG21	1:A:134[B]:VAL:HG21	1.91	0.52
1:D:91:VAL:HG21	1:D:134:VAL:HG21	1.92	0.52
1:A:43:ARG:O	1:A:47[B]:ILE:HG12	2.10	0.51
1:D:8:LYS:NZ	4:D:405:HOH:O	2.33	0.51
1:B:245:LEU:HD11	1:B:258[A]:ILE:HG13	1.93	0.51
1:B:245:LEU:HD11	1:B:258[B]:ILE:HD13	1.95	0.48
1:A:100:GLN:HG3	4:A:521[A]:HOH:O	2.14	0.48
1:A:268[B]:LEU:HD13	1:C:153[B]:ARG:HH21	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:LYS:HD2	1:D:109:PHE:HB3	1.95	0.47
1:C:44:LEU:HD11	1:C:62:GLU:HB2	1.97	0.47
1:B:43:ARG:O	1:B:47[B]:ILE:HG12	2.15	0.46
1:B:140:ARG:NH1	4:B:407[A]:HOH:O	2.36	0.45
1:B:132:LYS:HD2	1:C:109:PHE:HB3	1.99	0.44
1:D:159:ASN:ND2	4:D:449[B]:HOH:O	2.14	0.44
1:C:211:ALA:HB1	1:C:215:MET:HB3	1.99	0.44
1:B:24:HIS:CE1	1:B:235:PRO:HG3	2.53	0.44
1:A:43:ARG:HB3	1:A:46:LEU:HB3	1.98	0.43
1:C:46:LEU:O	1:C:50:ILE:HD12	2.20	0.42
1:D:41[B]:PHE:CG	1:D:42:ASP:N	2.87	0.42
1:A:91:VAL:HG21	1:A:134[B]:VAL:HG11	2.03	0.41
1:A:44:LEU:HD23	1:A:44:LEU:HA	1.93	0.41
1:D:231:ASP:OD2	4:D:402:HOH:O	2.22	0.41

All (2) 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)
4:A:562:HOH:O	4:B:450:HOH:O[3_454]	2.18	0.02
4:B:526:HOH:O	4:C:436:HOH:O[4_445]	2.19	0.01

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	265/277 (96%)	256 (97%)	8 (3%)	1 (0%)	34 17
1	B	260/277 (94%)	251 (96%)	9 (4%)	0	100 100
1	C	273/277 (99%)	263 (96%)	9 (3%)	1 (0%)	34 17
1	D	272/277 (98%)	262 (96%)	10 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1070/1108 (97%)	1032 (96%)	36 (3%)	2 (0%)	47 29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	151	PRO
1	C	151	PRO

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	209/217 (96%)	207 (99%)	2 (1%)	76 63
1	B	206/217 (95%)	205 (100%)	1 (0%)	88 83
1	C	205/217 (94%)	204 (100%)	1 (0%)	88 83
1	D	204/217 (94%)	204 (100%)	0	100 100
All	All	824/868 (95%)	820 (100%)	4 (0%)	88 83

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	ASN
1	A	256	ASP
1	B	106	ASN
1	C	256	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	214	GLN
1	B	24	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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

14 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	EDO	D	301	-	3,3,3	0.50	0	2,2,2	1.17	0
3	SO4	D	303	-	4,4,4	0.16	0	6,6,6	0.18	0
3	SO4	B	302	-	4,4,4	0.19	0	6,6,6	0.21	0
2	EDO	C	302	-	3,3,3	0.52	0	2,2,2	0.24	0
3	SO4	B	303	-	4,4,4	0.15	0	6,6,6	0.09	0
3	SO4	D	304	-	4,4,4	0.15	0	6,6,6	0.18	0
2	EDO	C	301	-	3,3,3	0.41	0	2,2,2	0.87	0
2	EDO	A	301	-	3,3,3	0.34	0	2,2,2	0.86	0
3	SO4	C	303	-	4,4,4	0.15	0	6,6,6	0.09	0
3	SO4	A	302	-	4,4,4	0.13	0	6,6,6	0.09	0
3	SO4	A	303	-	4,4,4	0.14	0	6,6,6	0.06	0
2	EDO	B	301	-	3,3,3	0.45	0	2,2,2	1.11	0
3	SO4	C	304	-	4,4,4	0.18	0	6,6,6	0.05	0
2	EDO	D	302	-	3,3,3	0.52	0	2,2,2	0.31	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
2	EDO	D	301	-	-	0/1/1/1	-
2	EDO	C	302	-	-	0/1/1/1	-
2	EDO	C	301	-	-	0/1/1/1	-
2	EDO	A	301	-	-	0/1/1/1	-
2	EDO	B	301	-	-	0/1/1/1	-
2	EDO	D	302	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	304	SO4	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	259/277 (93%)	0.64	18 (6%) 16 22	5, 12, 27, 46	0
1	B	257/277 (92%)	0.72	19 (7%) 14 19	5, 12, 28, 57	0
1	C	267/277 (96%)	0.80	29 (10%) 5 7	6, 14, 37, 52	0
1	D	267/277 (96%)	0.78	29 (10%) 5 7	5, 14, 37, 50	0
All	All	1050/1108 (94%)	0.74	95 (9%) 9 12	5, 13, 35, 57	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	207	LEU	10.5
1	A	207	LEU	8.6
1	C	198	ALA	7.5
1	A	1	MET	7.3
1	A	2	ALA	7.3
1	C	203[A]	VAL	6.9
1	C	202	ILE	6.6
1	C	41[A]	PHE	6.4
1	D	42	ASP	5.7
1	D	44	LEU	5.3
1	B	269[A]	LEU	5.0
1	C	44	LEU	4.9
1	C	201	ALA	4.9
1	A	206	ALA	4.7
1	D	197	LEU	4.7
1	D	203	VAL	4.6
1	B	84	GLU	4.6
1	C	42	ASP	4.6
1	A	269[A]	LEU	4.4
1	C	199	MET	4.4
1	C	84	GLU	4.3

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Mol	Chain	Res	Type	RSRZ
1	D	41[A]	PHE	4.2
1	C	196	THR	4.1
1	D	199	MET	3.8
1	B	208	GLY	3.8
1	D	196	THR	3.7
1	C	85	GLY	3.7
1	B	206	ALA	3.5
1	B	268[A]	LEU	3.5
1	D	46	LEU	3.4
1	C	197	LEU	3.4
1	A	268[A]	LEU	3.4
1	D	198	ALA	3.3
1	B	217[A]	LEU	3.3
1	C	206	ALA	3.2
1	D	202	ILE	3.2
1	C	200	SER	3.1
1	D	200	SER	3.1
1	A	217	LEU	3.0
1	B	158	TYR	3.0
1	A	208	GLY	2.9
1	B	16	ILE	2.9
1	C	16	ILE	2.9
1	D	81	VAL	2.9
1	C	210	GLU	2.8
1	D	16	ILE	2.8
1	C	204	GLY	2.8
1	B	213	GLN	2.8
1	A	44	LEU	2.8
1	B	259	PHE	2.8
1	C	217[A]	LEU	2.7
1	D	51	THR	2.7
1	D	47	ILE	2.6
1	D	149	PHE	2.5
1	B	57	PRO	2.5
1	C	46	LEU	2.5
1	A	213	GLN	2.5
1	D	208	GLY	2.5
1	D	207	LEU	2.5
1	D	84	GLU	2.4
1	D	43	ARG	2.4
1	B	211	ALA	2.4
1	A	140	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	42	ASP	2.4
1	D	209	ASP	2.4
1	A	84[A]	GLU	2.4
1	C	49	ARG	2.3
1	C	43	ARG	2.3
1	B	83	GLY	2.3
1	C	269	LEU	2.3
1	B	195	ARG	2.3
1	A	233	LYS	2.3
1	A	177[A]	ARG	2.3
1	B	233	LYS	2.3
1	B	19	SER	2.2
1	C	209	ASP	2.2
1	C	33	GLY	2.2
1	D	201	ALA	2.2
1	D	50	ILE	2.2
1	D	217[A]	LEU	2.2
1	C	47	ILE	2.2
1	C	205	GLY	2.2
1	D	155[A]	MET	2.1
1	B	42	ASP	2.1
1	B	50	ILE	2.1
1	D	7	GLY	2.1
1	A	69	GLU	2.1
1	C	207	LEU	2.1
1	D	6	GLU	2.1
1	C	211	ALA	2.1
1	A	216	GLN	2.0
1	A	174	PHE	2.0
1	D	19[A]	SER	2.0
1	D	206	ALA	2.0
1	C	7	GLY	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 [i](#)

There are no monosaccharides 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
3	SO4	A	302	5/5	0.72	0.33	20,25,41,45	5
2	EDO	D	302	4/4	0.73	0.17	24,28,30,31	0
3	SO4	B	302	5/5	0.76	0.27	20,25,34,39	5
3	SO4	C	304	5/5	0.76	0.27	24,29,33,34	5
3	SO4	C	303	5/5	0.80	0.24	24,29,42,42	5
2	EDO	C	302	4/4	0.81	0.20	26,29,30,30	0
3	SO4	A	303	5/5	0.82	0.32	33,40,46,51	5
3	SO4	D	303	5/5	0.82	0.23	25,27,32,37	5
3	SO4	D	304	5/5	0.84	0.24	22,25,30,31	5
2	EDO	D	301	4/4	0.90	0.19	8,10,12,16	0
3	SO4	B	303	5/5	0.91	0.20	44,47,56,57	5
2	EDO	C	301	4/4	0.95	0.11	9,10,11,12	0
2	EDO	B	301	4/4	0.96	0.14	6,9,10,12	0
2	EDO	A	301	4/4	0.97	0.10	9,9,11,12	0

6.5 Other polymers

There are no such residues in this entry.