



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:02 PM GMT

PDB ID : 4UOV
Title : The structure of a tetrameric alpha-carbonic anhydrase from *Thermovibrio ammonificans* reveals a core formed around intermolecular disulfides, which contribute to its thermostability.
Authors : James, P.; Isupov, M.; Sayer, C.; Berg, S.; Lioliou, M.; Kotlar, H.; Littlechild, J.
Deposited on : 2014-06-10
Resolution : 1.85 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
<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

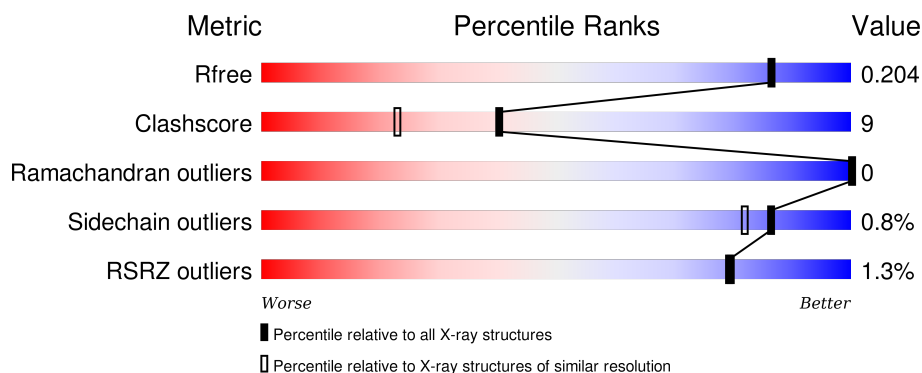
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

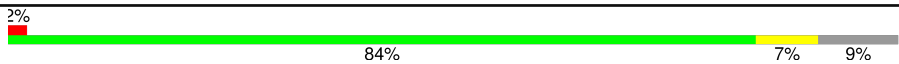
The reported resolution of this entry is 1.85 Å.

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	1745 (1.86-1.86)
Clashscore	102246	1898 (1.86-1.86)
Ramachandran outliers	100387	1875 (1.86-1.86)
Sidechain outliers	100360	1875 (1.86-1.86)
RSRZ outliers	91569	1747 (1.86-1.86)

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	247	
1	B	247	
1	C	247	
1	D	247	
1	E	247	

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Mol	Chain	Length	Quality of chain
1	F	247	 % 81% 10% 9%

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
10	PGE	B	305	-	-	-	X
10	PGE	D	305	-	-	-	X
10	PGE	E	305	-	-	-	X
10	PGE	F	309	-	-	-	X
11	PG5	B	312	-	-	X	X
4	PE3	A	300	-	-	-	X
4	PE3	B	300	-	-	-	X
4	PE3	F	300	-	-	-	X
5	B3P	A	301[A]	-	-	-	X
5	B3P	A	301[B]	-	-	-	X
5	B3P	C	301	-	-	-	X
5	B3P	D	301	-	-	-	X
5	B3P	E	301[A]	-	-	-	X
5	B3P	E	301[B]	-	-	-	X
7	CL	A	304	-	-	-	X
8	PEG	A	305	-	-	-	X
8	PEG	B	314	-	-	-	X
8	PEG	C	314	-	-	-	X
8	PEG	E	314	-	-	-	X
8	PEG	F	314	-	-	-	X
9	TLA	A	306	-	-	-	X
9	TLA	B	306	-	-	-	X
9	TLA	F	306	-	-	-	X

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 14247 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.

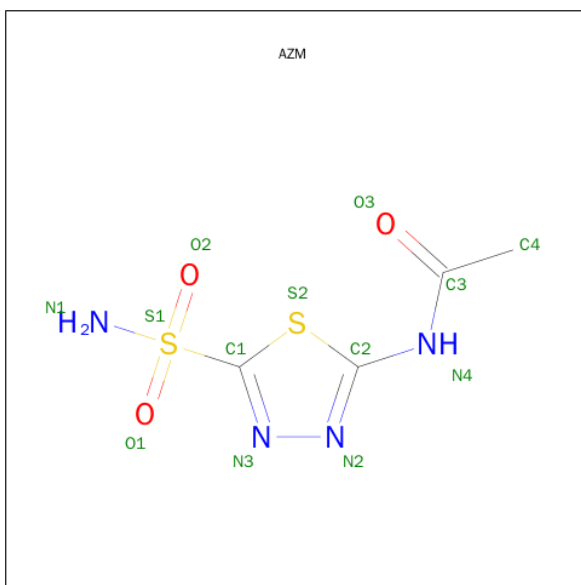
- Molecule 1 is a protein called CARBONATE DEHYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	225	Total	C	N	O	S	0	34	0
			1994	1296	342	343	13			
1	B	225	Total	C	N	O	S	0	33	0
			1997	1299	343	344	11			
1	C	225	Total	C	N	O	S	0	30	0
			1979	1286	342	339	12			
1	D	225	Total	C	N	O	S	0	31	0
			1991	1293	346	341	11			
1	E	225	Total	C	N	O	S	0	27	0
			1961	1273	337	339	12			
1	F	225	Total	C	N	O	S	0	26	0
			1962	1272	342	337	11			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

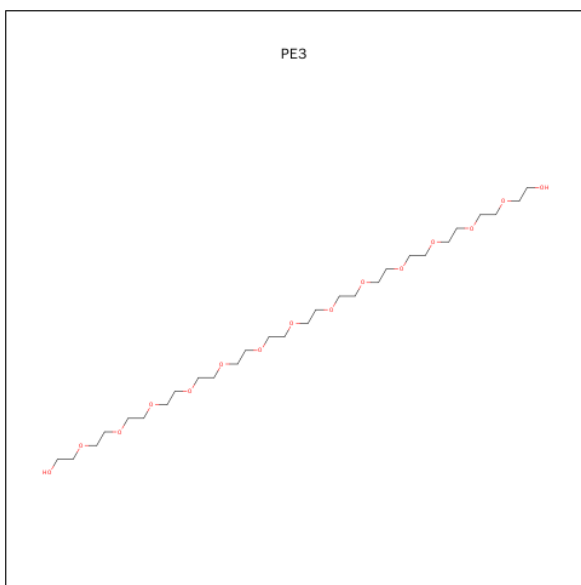
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Zn	0	0
			1	1		
2	E	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	F	1	Total	Zn	0	0
			1	1		

- Molecule 3 is 5-ACETAMIDO-1,3,4-THIADIAZOLE-2-SULFONAMIDE (three-letter code: AZM) (formula: C₄H₆N₄O₃S₂).



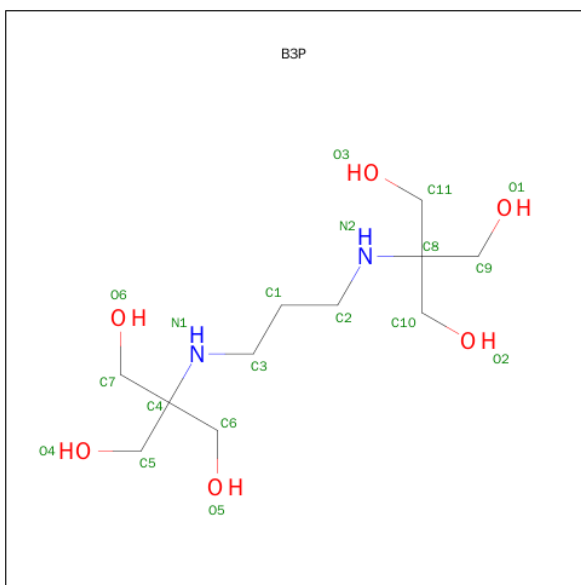
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			13	4	4	3	2		
3	B	1	Total	C	N	O	S	0	0
			13	4	4	3	2		
3	C	1	Total	C	N	O	S	0	0
			13	4	4	3	2		
3	D	1	Total	C	N	O	S	0	0
			13	4	4	3	2		
3	E	1	Total	C	N	O	S	0	0
			13	4	4	3	2		
3	F	1	Total	C	N	O	S	0	0
			13	4	4	3	2		

- Molecule 4 is 3,6,9,12,15,18,21,24,27,30,33,36,39-TRIDECAOXAHENTETRACONTANE-1,41-DIOL (three-letter code: PE3) (formula: C₂₈H₅₈O₁₅).



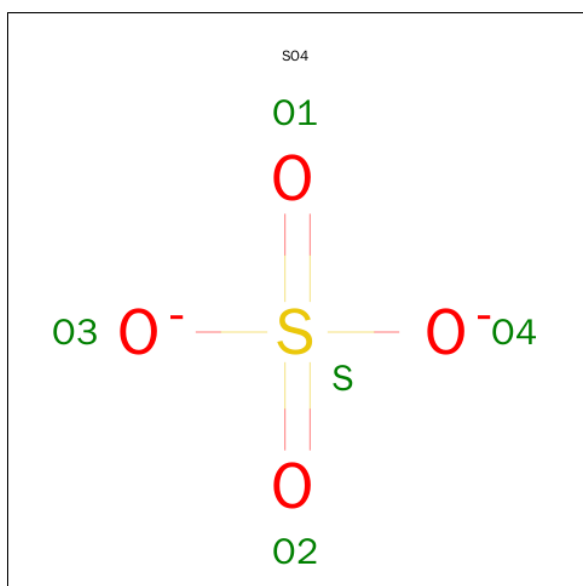
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			21	14	7		
4	B	1	Total	C	O	0	0
			21	14	7		
4	F	1	Total	C	O	0	0
			21	14	7		

- Molecule 5 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYL AMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: B3P) (formula: $C_{11}H_{26}N_2O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	1
			24	14	2	8		
5	C	1	Total	C	N	O	0	0
			19	11	2	6		
5	D	1	Total	C	N	O	0	0
			19	11	2	6		
5	E	1	Total	C	N	O	0	1
			22	11	2	9		
5	F	1	Total	C	N	O	0	1
			21	11	2	8		

- Molecule 6 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).

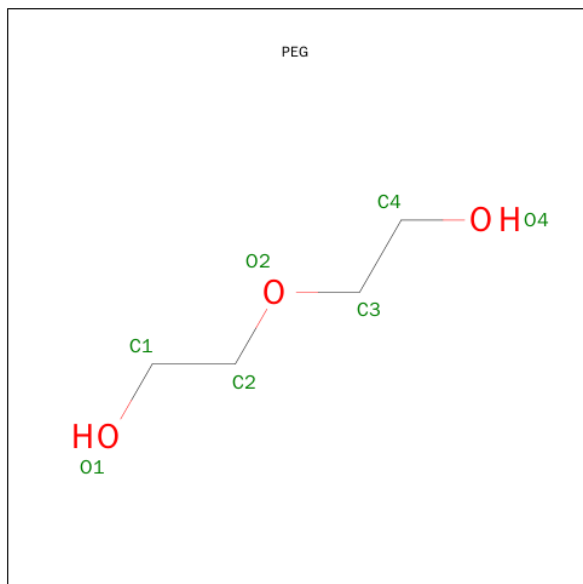


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	C	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		
6	E	1	Total	O	S	0	0
			5	4	1		
6	F	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	2	Total	Cl	0	0
			2	2		
7	E	2	Total	Cl	0	0
			2	2		
7	B	3	Total	Cl	0	0
			3	3		
7	C	3	Total	Cl	0	0
			3	3		
7	A	3	Total	Cl	0	0
			3	3		
7	F	2	Total	Cl	0	0
			2	2		

- Molecule 8 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



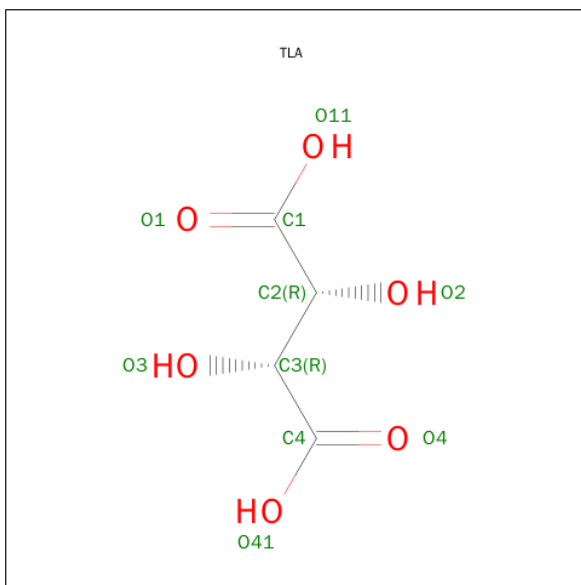
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			7	4	3		
8	B	1	Total	C	O	0	0
			7	4	3		
8	B	1	Total	C	O	0	0
			7	4	3		
8	C	1	Total	C	O	0	0
			7	4	3		
8	E	1	Total	C	O	0	0
			7	4	3		
8	E	1	Total	C	O	0	0
			7	4	3		

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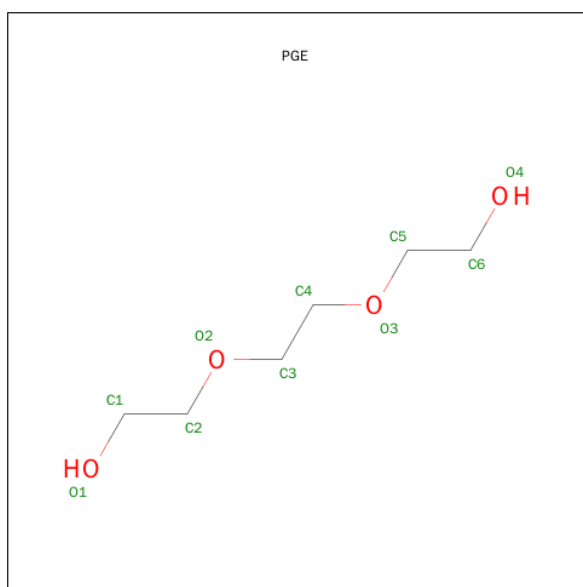
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	F	1	Total	C	O	0	0
			7	4	3		
8	F	1	Total	C	O	0	0
			7	4	3		

- Molecule 9 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula: C₄H₆O₆).



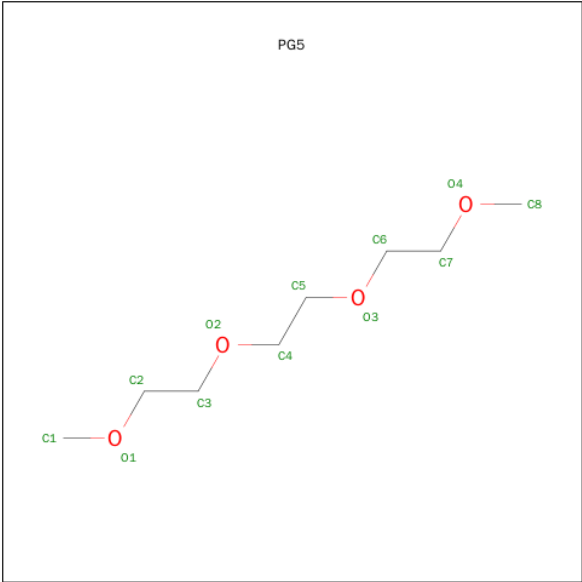
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			10	4	6		
9	B	1	Total	C	O	0	0
			10	4	6		
9	F	1	Total	C	O	0	0
			10	4	6		

- Molecule 10 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	C	O	0	0
			10	6	4		
10	C	1	Total	C	O	0	0
			10	6	4		
10	D	1	Total	C	O	0	0
			10	6	4		
10	E	1	Total	C	O	0	0
			10	6	4		
10	F	1	Total	C	O	0	0
			10	6	4		

- Molecule 11 is 1-METHOXY-2-[2-(2-METHOXY-ETHOXY)]-ETHANE (three-letter code: PG5) (formula: C₈H₁₈O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	B	1	Total	C	O	0	0
			12	8	4		

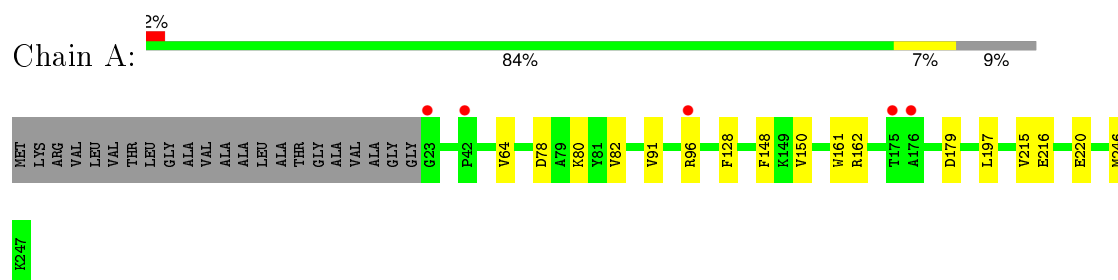
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	360	Total	O	0	0
			360	360		
12	B	363	Total	O	0	0
			363	363		
12	C	291	Total	O	0	0
			291	291		
12	D	242	Total	O	0	0
			242	242		
12	E	339	Total	O	0	0
			339	339		
12	F	323	Total	O	0	0
			323	323		

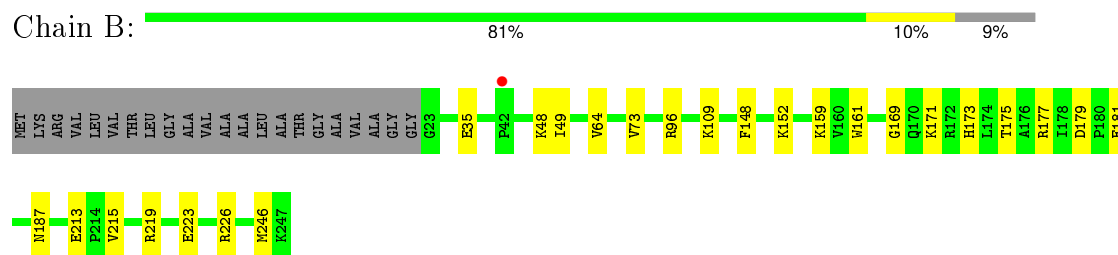
3 Residue-property plots [i](#)

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.

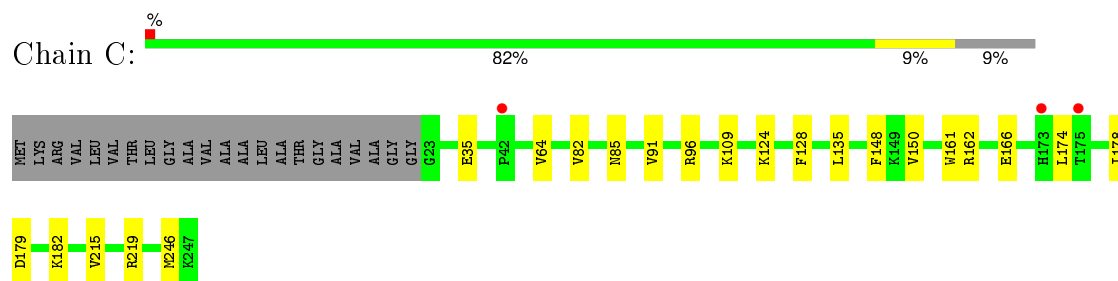
• Molecule 1: CARBONATE DEHYDRATASE



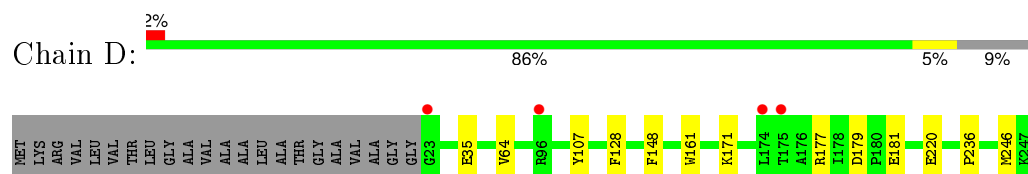
• Molecule 1: CARBONATE DEHYDRATASE



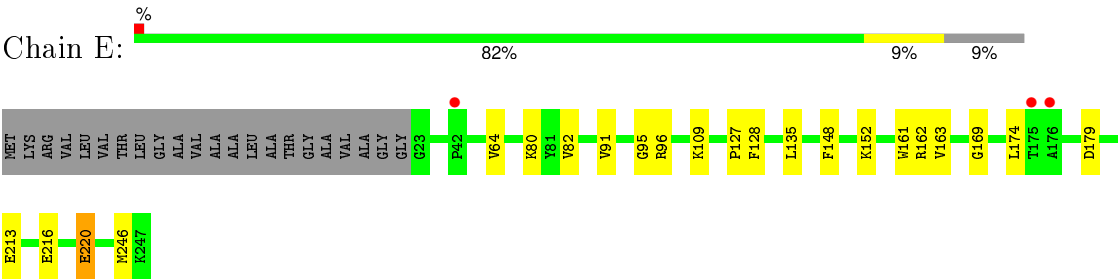
• Molecule 1: CARBONATE DEHYDRATASE



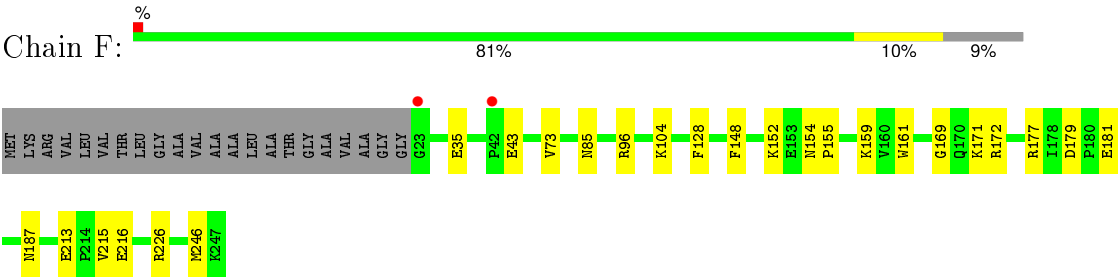
• Molecule 1: CARBONATE DEHYDRATASE



• Molecule 1: CARBONATE DEHYDRATASE



● Molecule 1: CARBONATE DEHYDRATASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	163.75Å 283.02Å 52.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.01 – 1.85 42.01 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.5 (42.01-1.85) 99.5 (42.01-1.85)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 1.86Å)	Xtriage
Refinement program	REFMAC 5.8.0071	Depositor
R, R_{free}	0.172 , 0.204 0.172 , 0.204	Depositor DCC
R_{free} test set	10402 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.365	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 60.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 206860 reflections (0.000%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14247	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.78% 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: ZN, PGE, AZM, CL, B3P, PE3, TLA, PG5, SO4, PEG

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	0.42	0/2152	0.54	0/2904
1	B	0.44	0/2153	0.56	0/2904
1	C	0.41	0/2126	0.53	0/2867
1	D	0.41	0/2141	0.54	0/2886
1	E	0.44	0/2097	0.55	0/2829
1	F	0.46	0/2096	0.56	0/2825
All	All	0.43	0/12765	0.55	0/17215

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1994	0	2071	26	0
1	B	1997	0	2070	76	1
1	C	1979	0	2048	30	0
1	D	1991	0	2064	28	0
1	E	1961	0	2024	38	0
1	F	1962	0	2022	39	1
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	13	0	6	0	0
3	B	13	0	5	0	0
3	C	13	0	6	1	0
3	D	13	0	6	0	0
3	E	13	0	6	0	0
3	F	13	0	6	0	0
4	A	21	0	24	0	0
4	B	21	0	24	0	0
4	F	21	0	24	0	0
5	A	24	0	20	0	0
5	C	19	0	26	3	0
5	D	19	0	26	0	0
5	E	22	0	18	0	0
5	F	21	0	12	0	0
6	A	5	0	0	1	0
6	B	5	0	0	0	0
6	C	5	0	0	1	0
6	D	5	0	0	0	0
6	E	5	0	0	0	0
6	F	5	0	0	0	0
7	A	3	0	0	1	0
7	B	3	0	0	1	0
7	C	3	0	0	1	0
7	D	2	0	0	1	0
7	E	2	0	0	1	0
7	F	2	0	0	1	0
8	A	7	0	10	1	0
8	B	14	0	20	2	0
8	C	7	0	10	1	0
8	E	14	0	20	3	0
8	F	14	0	20	4	0
9	A	10	0	4	0	0
9	B	10	0	4	0	0
9	F	10	0	4	0	0
10	B	10	0	14	5	0
10	C	10	0	14	0	0
10	D	10	0	14	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	E	10	0	14	1	0
10	F	10	0	14	0	0
11	B	12	0	18	31	0
12	A	360	0	0	13	0
12	B	363	0	0	16	0
12	C	291	0	0	8	0
12	D	242	0	0	15	0
12	E	339	0	0	18	0
12	F	323	0	0	11	0
All	All	14247	0	12688	218	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 9.

All (218) 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:B:246[B]:MET:CE	1:D:64[B]:VAL:HG11	1.55	1.35
1:A:64[B]:VAL:HG11	1:C:246[B]:MET:CE	1.57	1.32
1:B:171[B]:LYS:NZ	11:B:312:PG5:H12	1.46	1.26
1:E:220[B]:GLU:CD	12:E:2279:HOH:O	1.73	1.24
1:B:173[B]:HIS:NE2	11:B:312:PG5:O1	1.71	1.23
1:B:173[B]:HIS:CE1	11:B:312:PG5:O1	1.94	1.20
1:A:64[B]:VAL:CG1	1:C:246[B]:MET:HE1	1.70	1.20
1:E:220[B]:GLU:OE1	12:E:2279:HOH:O	1.56	1.20
1:E:96[B]:ARG:CZ	12:E:2115:HOH:O	1.90	1.19
1:B:171[B]:LYS:CE	11:B:312:PG5:H12	1.72	1.18
1:B:173[B]:HIS:NE2	11:B:312:PG5:C1	2.06	1.17
1:B:171[B]:LYS:HE3	11:B:312:PG5:C2	1.73	1.17
1:B:109[B]:LYS:HE2	10:B:305:PGE:H42	1.25	1.13
1:E:64[B]:VAL:HG11	1:F:246[B]:MET:CE	1.78	1.13
1:B:173[B]:HIS:CD2	12:B:2134:HOH:O	1.99	1.13
1:C:162[B]:ARG:HG2	1:C:162[B]:ARG:HH11	1.09	1.13
1:B:171[B]:LYS:HE3	11:B:312:PG5:H22	1.12	1.11
1:A:246[B]:MET:HE1	1:C:64[B]:VAL:CG1	1.81	1.11
1:A:220[A]:GLU:OE2	12:A:2309:HOH:O	1.67	1.11
1:B:171[B]:LYS:CE	11:B:312:PG5:C1	2.30	1.10
1:B:171[B]:LYS:CE	11:B:312:PG5:H22	1.82	1.10
1:A:246[B]:MET:CE	1:C:64[B]:VAL:HG11	1.82	1.09
1:B:171[B]:LYS:NZ	11:B:312:PG5:C1	2.16	1.07
1:D:179[B]:ASP:OD2	12:D:2181:HOH:O	1.70	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177[B]:ARG:HH11	1:B:177[B]:ARG:HG2	0.94	1.06
8:A:305:PEG:O1	12:A:2186:HOH:O	1.69	1.06
1:F:226[B]:ARG:HH11	1:F:226[B]:ARG:HG2	0.90	1.05
1:B:96[B]:ARG:NH2	12:B:2130:HOH:O	1.84	1.05
1:D:220[B]:GLU:CG	12:D:2211:HOH:O	2.02	1.05
1:B:171[B]:LYS:HE3	11:B:312:PG5:C1	1.84	1.04
1:B:246[B]:MET:CE	1:D:64[B]:VAL:CG1	2.35	1.04
1:B:173[B]:HIS:NE2	11:B:312:PG5:H11	1.72	1.04
1:B:177[B]:ARG:HD2	12:B:2126:HOH:O	1.58	1.03
1:F:177[B]:ARG:HG2	1:F:177[B]:ARG:HH11	0.93	1.03
1:A:64[B]:VAL:CG1	1:C:246[B]:MET:CE	2.32	1.01
1:E:152[B]:LYS:HE3	12:E:2212:HOH:O	0.83	1.00
1:E:64[B]:VAL:CG1	1:F:246[B]:MET:HE3	1.89	1.00
1:C:179[B]:ASP:OD2	12:C:2223:HOH:O	1.82	0.98
1:B:177[B]:ARG:HH11	1:B:177[B]:ARG:CG	1.75	0.97
6:A:302:SO4:O4	12:A:2192:HOH:O	1.83	0.97
1:E:179[B]:ASP:OD2	12:E:2241:HOH:O	1.81	0.97
1:B:246[B]:MET:HE1	1:D:64[B]:VAL:HG11	0.99	0.97
1:C:179[B]:ASP:HB2	12:C:2222:HOH:O	1.63	0.97
1:E:64[B]:VAL:CG1	1:F:246[B]:MET:CE	2.44	0.95
1:F:226[B]:ARG:HG2	1:F:226[B]:ARG:NH1	1.70	0.95
1:B:177[B]:ARG:NH1	1:B:177[B]:ARG:HG2	1.75	0.95
1:D:181[A]:GLU:OE1	12:D:2072:HOH:O	1.83	0.94
1:B:246[B]:MET:HE1	1:D:64[B]:VAL:CG1	1.93	0.94
1:B:171[B]:LYS:HG3	11:B:312:PG5:H22	1.47	0.94
1:A:246[B]:MET:HE1	1:C:64[B]:VAL:HG11	1.41	0.94
1:D:35[A]:GLU:OE1	12:D:2012:HOH:O	1.84	0.93
1:F:177[B]:ARG:NH1	1:F:177[B]:ARG:HG2	1.66	0.93
1:C:162[B]:ARG:NH1	1:C:162[B]:ARG:HG2	1.81	0.92
1:F:177[B]:ARG:CG	1:F:177[B]:ARG:HH11	1.82	0.92
1:E:64[B]:VAL:HG11	1:F:246[B]:MET:HE1	1.51	0.92
1:B:246[B]:MET:HE3	1:D:64[B]:VAL:CG1	1.99	0.91
1:F:226[B]:ARG:HH11	1:F:226[B]:ARG:CG	1.83	0.91
1:B:96[B]:ARG:NE	12:B:2130:HOH:O	1.99	0.91
1:B:171[B]:LYS:HE3	11:B:312:PG5:H12	1.43	0.90
1:E:213[B]:GLU:OE1	12:E:2276:HOH:O	1.90	0.88
1:E:80[B]:LYS:NZ	12:E:2118:HOH:O	1.71	0.87
1:B:173[B]:HIS:CD2	11:B:312:PG5:O1	2.27	0.87
1:B:171[B]:LYS:HZ1	11:B:312:PG5:H12	1.05	0.87
1:E:64[B]:VAL:HG13	1:F:246[B]:MET:HE3	1.57	0.84
1:E:96[B]:ARG:NH2	12:E:2115:HOH:O	2.00	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:220[B]:GLU:HG2	12:D:2211:HOH:O	1.71	0.83
1:F:43[A]:GLU:CG	12:F:2033:HOH:O	2.27	0.83
1:D:171[B]:LYS:CE	12:D:2177:HOH:O	2.27	0.83
1:A:162[A]:ARG:NE	12:A:2246:HOH:O	1.96	0.83
1:B:213[B]:GLU:CD	12:B:2297:HOH:O	2.16	0.82
1:B:171[B]:LYS:CG	11:B:312:PG5:H22	2.07	0.82
7:D:304:CL:CL	12:D:2090:HOH:O	2.34	0.82
1:A:179[B]:ASP:OD2	12:A:2262:HOH:O	1.97	0.82
1:E:220[A]:GLU:OE1	12:E:2222:HOH:O	1.98	0.82
1:E:220[B]:GLU:OE2	12:E:2279:HOH:O	1.85	0.82
1:F:171[A]:LYS:NZ	12:F:2121:HOH:O	2.13	0.81
1:B:171[B]:LYS:HG3	11:B:312:PG5:C2	2.09	0.81
1:D:177[B]:ARG:HH11	1:D:177[B]:ARG:HA	1.46	0.79
1:A:246[B]:MET:HE1	1:C:64[B]:VAL:HG13	1.63	0.79
1:A:64[B]:VAL:HG11	1:C:246[B]:MET:HE1	0.82	0.78
1:B:171[B]:LYS:HZ1	11:B:312:PG5:C1	1.85	0.78
1:F:35[B]:GLU:HG2	12:F:2019:HOH:O	1.83	0.78
1:A:246[B]:MET:HE2	1:C:64[B]:VAL:HG11	1.67	0.77
7:C:304:CL:CL	12:C:2123:HOH:O	2.39	0.77
1:B:177[B]:ARG:NH2	1:B:179[B]:ASP:OD2	2.17	0.77
1:B:96[B]:ARG:CZ	12:B:2130:HOH:O	2.19	0.75
1:B:173[B]:HIS:HD2	12:B:2134:HOH:O	1.49	0.75
7:A:304:CL:CL	12:A:2143:HOH:O	2.41	0.75
1:B:171[B]:LYS:CD	11:B:312:PG5:H22	2.16	0.75
1:B:109[B]:LYS:CE	10:B:305:PGE:H42	2.13	0.72
1:A:64[B]:VAL:CG1	1:C:246[B]:MET:HE3	2.18	0.72
7:F:304:CL:CL	12:F:2125:HOH:O	2.45	0.72
1:B:213[B]:GLU:HG2	12:B:2276:HOH:O	1.90	0.71
1:D:35[B]:GLU:HG2	12:D:2012:HOH:O	1.91	0.71
1:B:246[B]:MET:HE3	1:D:64[B]:VAL:HG13	1.73	0.71
1:F:215[B]:VAL:HG11	12:F:2223:HOH:O	1.91	0.70
1:C:124[B]:LYS:HE3	6:C:302:SO4:O2	1.90	0.70
1:E:246[B]:MET:HE1	1:F:246[B]:MET:HG2	1.71	0.70
1:C:35[B]:GLU:OE2	12:C:2025:HOH:O	0.70	0.70
1:C:162[B]:ARG:CG	1:C:162[B]:ARG:HH11	1.95	0.70
1:B:177[B]:ARG:CD	12:B:2126:HOH:O	2.28	0.69
1:B:109[B]:LYS:HZ1	10:B:305:PGE:H2	1.56	0.69
7:B:304:CL:CL	12:B:2146:HOH:O	2.47	0.69
1:B:171[B]:LYS:CE	11:B:312:PG5:H13	2.22	0.68
1:A:246[B]:MET:CE	1:C:64[B]:VAL:CG1	2.51	0.67
1:D:35[B]:GLU:OE1	12:D:2016:HOH:O	2.12	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:171[B]:LYS:HE3	12:D:2177:HOH:O	1.93	0.66
1:B:215[B]:VAL:HG11	12:B:2241:HOH:O	1.95	0.66
1:B:177[B]:ARG:HH22	1:B:179[B]:ASP:CG	2.00	0.65
1:A:197:LEU:HD12	12:A:2294:HOH:O	1.96	0.65
1:D:35[B]:GLU:OE1	1:D:35[B]:GLU:HA	1.96	0.65
1:D:177[B]:ARG:NH1	1:D:177[B]:ARG:HA	2.12	0.65
7:E:304:CL:CL	12:E:2125:HOH:O	2.50	0.65
1:E:152[B]:LYS:CE	12:E:2212:HOH:O	1.65	0.64
1:B:173[B]:HIS:CD2	11:B:312:PG5:C1	2.81	0.64
1:A:64[B]:VAL:HG13	1:C:246[B]:MET:HE3	1.81	0.62
1:F:159:LYS:HA	8:F:310:PEG:H31	1.83	0.60
1:B:173[B]:HIS:CE1	11:B:312:PG5:H32	2.36	0.60
8:F:310:PEG:H22	12:F:2120:HOH:O	2.03	0.59
1:A:215[B]:VAL:HG11	12:A:2236:HOH:O	2.00	0.59
1:E:246[B]:MET:CE	1:F:246[B]:MET:HG2	2.32	0.59
1:D:171[B]:LYS:CD	12:D:2177:HOH:O	2.50	0.59
1:B:48[A]:LYS:HG3	1:B:49:ILE:HG23	1.84	0.59
1:D:107:TYR:CD2	10:D:305:PGE:H52	2.38	0.58
1:E:64[B]:VAL:HG11	1:F:246[B]:MET:HE3	1.56	0.58
5:C:301:B3P:H31	12:C:2269:HOH:O	2.03	0.58
1:D:171[B]:LYS:HD2	12:D:2177:HOH:O	2.02	0.57
1:E:96[B]:ARG:NH1	12:E:2115:HOH:O	2.18	0.57
1:D:220[B]:GLU:HG3	12:D:2211:HOH:O	1.84	0.57
1:F:177[B]:ARG:NH1	1:F:177[B]:ARG:CG	2.49	0.57
1:E:82[B]:VAL:HG22	1:E:91:VAL:HG22	1.87	0.56
1:F:226[B]:ARG:NH1	1:F:226[B]:ARG:CG	2.53	0.56
1:C:96[B]:ARG:NH2	12:C:2116:HOH:O	2.39	0.56
1:E:64[B]:VAL:CG1	1:F:246[B]:MET:HE1	2.23	0.56
1:B:223[B]:GLU:HA	1:B:223[B]:GLU:OE1	2.06	0.56
1:B:109[B]:LYS:NZ	10:B:305:PGE:H52	2.20	0.55
1:A:78:ASP:HB2	1:A:96[A]:ARG:HH21	1.72	0.55
1:B:177[B]:ARG:NH1	1:B:177[B]:ARG:CG	2.47	0.54
1:B:213[B]:GLU:CG	12:B:2297:HOH:O	2.55	0.54
1:B:169:GLY:HA2	8:B:314:PEG:H32	1.89	0.53
1:A:82[A]:VAL:HG22	1:A:91:VAL:HG22	1.90	0.53
1:B:171[B]:LYS:HZ2	11:B:312:PG5:C1	2.17	0.53
1:B:177[B]:ARG:HG3	12:B:2126:HOH:O	2.08	0.53
1:B:219:ARG:O	1:B:223[B]:GLU:HG2	2.08	0.53
1:E:152[A]:LYS:NZ	12:E:2210:HOH:O	2.40	0.52
1:B:64[B]:VAL:HG11	1:D:246[B]:MET:HE1	1.91	0.52
1:E:127:PRO:HA	8:E:313:PEG:H22	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82[A]:VAL:HG22	1:C:91:VAL:HG22	1.92	0.52
1:C:109[B]:LYS:HE3	1:C:135[B]:LEU:HD12	1.92	0.51
1:E:95:GLY:HA3	10:E:305:PGE:H5	1.92	0.51
1:B:177[B]:ARG:CG	12:B:2126:HOH:O	2.59	0.51
1:E:246[B]:MET:HE3	1:F:246[B]:MET:CE	2.41	0.50
1:F:43[A]:GLU:HG3	12:F:2033:HOH:O	2.02	0.50
1:A:179[B]:ASP:CG	12:A:2262:HOH:O	2.47	0.50
1:D:220[B]:GLU:OE2	12:D:2211:HOH:O	2.20	0.50
1:F:73:VAL:HG11	1:F:181[B]:GLU:HG3	1.94	0.50
1:C:128:PHE:HB3	1:C:148:PHE:HB2	1.93	0.50
1:F:169:GLY:HA2	8:F:314:PEG:H21	1.93	0.49
1:B:213[B]:GLU:HG2	12:B:2297:HOH:O	2.11	0.49
1:B:173[B]:HIS:CE1	11:B:312:PG5:C2	2.94	0.49
1:E:152[A]:LYS:HG2	1:E:216:GLU:OE1	2.13	0.49
1:E:169:GLY:HA2	8:E:314:PEG:H21	1.95	0.48
1:B:109[B]:LYS:HE2	10:B:305:PGE:C4	2.19	0.48
1:F:128:PHE:HB3	1:F:148:PHE:HB2	1.95	0.47
1:B:187:ASN:HB3	1:B:213[B]:GLU:HG3	1.96	0.46
1:B:177[B]:ARG:NH2	1:B:179[B]:ASP:CG	2.62	0.46
1:B:171[B]:LYS:HD2	11:B:312:PG5:H13	1.97	0.46
1:A:215[B]:VAL:HG12	1:A:216:GLU:N	2.31	0.46
1:D:35[B]:GLU:OE1	1:D:35[B]:GLU:CA	2.61	0.46
1:C:219:ARG:NH1	12:C:2250:HOH:O	2.41	0.46
1:F:96[B]:ARG:HG3	1:F:96[B]:ARG:HH11	1.80	0.46
1:B:64[A]:VAL:HG11	1:D:246[A]:MET:SD	2.56	0.45
12:A:2322:HOH:O	8:E:313:PEG:H41	2.16	0.45
1:B:171[B]:LYS:CD	11:B:312:PG5:H13	2.47	0.45
1:F:187:ASN:HB3	1:F:213[B]:GLU:HG3	1.99	0.45
1:C:166:GLU:O	5:C:301:B3P:O3	2.25	0.45
1:B:35[B]:GLU:HA	1:B:35[B]:GLU:OE1	2.16	0.44
1:F:177[B]:ARG:HH22	1:F:179:ASP:HB2	1.83	0.44
1:A:162[A]:ARG:CZ	12:A:2246:HOH:O	2.50	0.44
1:E:246[B]:MET:HE3	1:F:246[B]:MET:HE2	1.98	0.44
1:E:80[A]:LYS:NZ	12:E:2117:HOH:O	2.50	0.44
1:A:162[A]:ARG:NH2	12:A:2246:HOH:O	2.48	0.44
1:E:152[B]:LYS:CG	12:E:2212:HOH:O	2.61	0.43
1:F:85:ASN:HA	8:F:314:PEG:H22	2.01	0.43
1:C:148:PHE:CD2	1:C:215[A]:VAL:HG23	2.54	0.43
1:B:226[B]:ARG:NH2	12:B:2302:HOH:O	2.51	0.43
1:F:43[A]:GLU:HG2	12:F:2033:HOH:O	2.05	0.43
1:F:215[B]:VAL:HG12	1:F:216:GLU:N	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:171[B]:LYS:NZ	12:D:2177:HOH:O	2.40	0.43
1:F:154:ASN:HA	1:F:155:PRO:HD3	1.91	0.42
3:C:299:AZM:S2	3:C:299:AZM:O3	2.77	0.42
1:A:150[B]:VAL:HG22	12:A:2223:HOH:O	2.18	0.42
1:F:43[A]:GLU:CD	12:F:2033:HOH:O	2.53	0.42
1:B:173[B]:HIS:CG	11:B:312:PG5:O1	2.72	0.42
1:B:171[B]:LYS:HE3	11:B:312:PG5:O1	2.07	0.42
1:B:152[B]:LYS:HB3	1:B:152[B]:LYS:HE2	1.82	0.42
1:B:159:LYS:HA	8:B:310:PEG:H22	2.01	0.42
1:B:171[B]:LYS:NZ	11:B:312:PG5:H13	2.20	0.42
1:E:64[B]:VAL:HG13	1:F:246[B]:MET:CE	2.32	0.42
1:C:174:LEU:HD22	1:C:178:ILE:HD11	2.01	0.41
1:B:173[B]:HIS:CD2	11:B:312:PG5:H11	2.48	0.41
1:C:179[A]:ASP:HB2	1:C:182:LYS:HD2	2.02	0.41
1:A:78:ASP:O	1:A:80[B]:LYS:HE3	2.20	0.41
1:E:82[B]:VAL:CG2	1:E:174:LEU:HD21	2.50	0.41
1:E:162[B]:ARG:HG3	1:E:163:VAL:HG13	2.03	0.41
1:E:109:LYS:HE3	1:E:135[B]:LEU:HD12	2.02	0.41
1:D:128:PHE:HB3	1:D:148:PHE:HB2	2.02	0.41
1:B:73:VAL:HG11	1:B:181[B]:GLU:HG3	2.03	0.41
1:E:152[B]:LYS:CD	12:E:2212:HOH:O	2.35	0.41
5:C:301:B3P:H112	5:C:301:B3P:H21	1.70	0.41
1:A:128:PHE:HB3	1:A:148:PHE:HB2	2.02	0.41
1:C:150[B]:VAL:HG22	12:C:2197:HOH:O	2.20	0.41
1:F:104[B]:LYS:NZ	12:F:2157:HOH:O	2.52	0.41
1:B:148:PHE:CD2	1:B:215[A]:VAL:HG23	2.56	0.40
1:F:104[A]:LYS:HE3	12:F:2157:HOH:O	2.21	0.40
1:E:128:PHE:HB3	1:E:148:PHE:HB2	2.03	0.40
1:E:162[A]:ARG:NH2	12:E:2220:HOH:O	2.54	0.40
1:C:85:ASN:HA	8:C:314:PEG:H22	2.03	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:B:175:THR:OG1	1:F:172[B]:ARG:NH2[3_444]	2.14	0.06

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	257/247 (104%)	253 (98%)	4 (2%)	0	100	100
1	B	256/247 (104%)	252 (98%)	4 (2%)	0	100	100
1	C	253/247 (102%)	249 (98%)	4 (2%)	0	100	100
1	D	254/247 (103%)	250 (98%)	4 (2%)	0	100	100
1	E	250/247 (101%)	245 (98%)	5 (2%)	0	100	100
1	F	249/247 (101%)	245 (98%)	4 (2%)	0	100	100
All	All	1519/1482 (102%)	1494 (98%)	25 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	230/208 (111%)	229 (100%)	1 (0%)	93	92
1	B	229/208 (110%)	228 (100%)	1 (0%)	93	92
1	C	226/208 (109%)	225 (100%)	1 (0%)	93	92
1	D	227/208 (109%)	225 (99%)	2 (1%)	84	79
1	E	223/208 (107%)	220 (99%)	3 (1%)	76	65
1	F	222/208 (107%)	219 (99%)	3 (1%)	74	63
All	All	1357/1248 (109%)	1346 (99%)	11 (1%)	86	82

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	161	TRP
1	B	161	TRP
1	C	161	TRP
1	D	161	TRP
1	D	236	PRO
1	E	161	TRP
1	E	220[A]	GLU
1	E	220[B]	GLU
1	F	152[A]	LYS
1	F	152[B]	LYS
1	F	161	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	140	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 61 ligands modelled in this entry, 21 are monoatomic - leaving 40 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	AZM	A	299	2	9,13,13	0.49	0	7,19,19	1.78	3 (42%)
4	PE3	A	300	-	20,20,42	0.54	0	19,19,41	0.31	0
5	B3P	A	301[A]	-	18,18,18	0.49	0	23,23,23	1.80	2 (8%)
5	B3P	A	301[B]	-	18,18,18	0.53	0	23,23,23	1.68	2 (8%)
6	SO4	A	302	-	4,4,4	0.29	0	6,6,6	0.10	0
8	PEG	A	305	-	6,6,6	0.40	0	5,5,5	0.38	0
9	TLA	A	306	-	3,9,9	0.57	0	6,12,12	0.91	0
3	AZM	B	299	2	9,13,13	0.68	0	7,19,19	2.23	3 (42%)
4	PE3	B	300	-	20,20,42	0.54	0	19,19,41	0.22	0
6	SO4	B	302	-	4,4,4	0.27	0	6,6,6	0.08	0
10	PGE	B	305	-	9,9,9	0.47	0	8,8,8	0.34	0
9	TLA	B	306	-	3,9,9	0.64	0	6,12,12	1.03	0
8	PEG	B	310	-	6,6,6	0.49	0	5,5,5	0.19	0
11	PG5	B	312	-	11,11,11	0.50	0	10,10,10	0.30	0
8	PEG	B	314	-	6,6,6	0.44	0	5,5,5	0.36	0
3	AZM	C	299	2	9,13,13	0.55	0	7,19,19	2.03	4 (57%)
5	B3P	C	301	-	18,18,18	0.63	0	23,23,23	5.37	13 (56%)
6	SO4	C	302	-	4,4,4	0.35	0	6,6,6	0.31	0
10	PGE	C	305	-	9,9,9	0.46	0	8,8,8	0.28	0
8	PEG	C	314	-	6,6,6	0.46	0	5,5,5	0.33	0
3	AZM	D	299	2	9,13,13	0.37	0	7,19,19	1.57	2 (28%)
5	B3P	D	301	-	18,18,18	0.42	0	23,23,23	1.96	2 (8%)
6	SO4	D	302	-	4,4,4	0.35	0	6,6,6	0.08	0
10	PGE	D	305	-	9,9,9	0.45	0	8,8,8	0.23	0
3	AZM	E	299	2	9,13,13	0.37	0	7,19,19	1.52	2 (28%)
5	B3P	E	301[A]	-	18,18,18	0.52	0	23,23,23	1.66	2 (8%)
5	B3P	E	301[B]	-	18,18,18	0.52	0	23,23,23	1.66	2 (8%)
6	SO4	E	302	-	4,4,4	0.26	0	6,6,6	0.06	0
10	PGE	E	305	-	9,9,9	0.44	0	8,8,8	0.29	0
8	PEG	E	313	-	6,6,6	0.46	0	5,5,5	0.49	0
8	PEG	E	314	-	6,6,6	0.41	0	5,5,5	0.39	0
3	AZM	F	299	2	9,13,13	0.79	0	7,19,19	2.05	2 (28%)
4	PE3	F	300	-	20,20,42	0.49	0	19,19,41	0.23	0
5	B3P	F	301[A]	-	18,18,18	0.43	0	23,23,23	1.68	3 (13%)
5	B3P	F	301[B]	-	18,18,18	0.43	0	23,23,23	1.68	3 (13%)
6	SO4	F	302	-	4,4,4	0.28	0	6,6,6	0.07	0
9	TLA	F	306	-	3,9,9	0.53	0	6,12,12	1.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	PGE	F	309	-	9,9,9	0.45	0	8,8,8	0.31	0
8	PEG	F	310	-	6,6,6	0.55	0	5,5,5	0.29	0
8	PEG	F	314	-	6,6,6	0.41	0	5,5,5	0.37	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
3	AZM	A	299	2	-	0/2/10/10	0/0/1/1
4	PE3	A	300	-	-	0/18/18/40	0/0/0/0
5	B3P	A	301[A]	-	-	0/28/28/28	0/0/0/0
5	B3P	A	301[B]	-	-	0/28/28/28	0/0/0/0
6	SO4	A	302	-	-	0/0/0/0	0/0/0/0
8	PEG	A	305	-	-	0/4/4/4	0/0/0/0
9	TLA	A	306	-	-	0/4/12/12	0/0/0/0
3	AZM	B	299	2	-	0/2/10/10	0/0/1/1
4	PE3	B	300	-	-	0/18/18/40	0/0/0/0
6	SO4	B	302	-	-	0/0/0/0	0/0/0/0
10	PGE	B	305	-	-	0/7/7/7	0/0/0/0
9	TLA	B	306	-	-	0/4/12/12	0/0/0/0
8	PEG	B	310	-	-	0/4/4/4	0/0/0/0
11	PG5	B	312	-	-	0/9/9/9	0/0/0/0
8	PEG	B	314	-	-	0/4/4/4	0/0/0/0
3	AZM	C	299	2	-	0/2/10/10	0/0/1/1
5	B3P	C	301	-	-	0/28/28/28	0/0/0/0
6	SO4	C	302	-	-	0/0/0/0	0/0/0/0
10	PGE	C	305	-	-	0/7/7/7	0/0/0/0
8	PEG	C	314	-	-	0/4/4/4	0/0/0/0
3	AZM	D	299	2	-	0/2/10/10	0/0/1/1
5	B3P	D	301	-	-	0/28/28/28	0/0/0/0
6	SO4	D	302	-	-	0/0/0/0	0/0/0/0
10	PGE	D	305	-	-	0/7/7/7	0/0/0/0
3	AZM	E	299	2	-	0/2/10/10	0/0/1/1
5	B3P	E	301[A]	-	-	0/28/28/28	0/0/0/0
5	B3P	E	301[B]	-	-	0/28/28/28	0/0/0/0
6	SO4	E	302	-	-	0/0/0/0	0/0/0/0
10	PGE	E	305	-	-	0/7/7/7	0/0/0/0
8	PEG	E	313	-	-	0/4/4/4	0/0/0/0
8	PEG	E	314	-	-	0/4/4/4	0/0/0/0
3	AZM	F	299	2	-	0/2/10/10	0/0/1/1
4	PE3	F	300	-	-	0/18/18/40	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	B3P	F	301[A]	-	-	0/28/28/28	0/0/0/0
5	B3P	F	301[B]	-	-	0/28/28/28	0/0/0/0
6	SO4	F	302	-	-	0/0/0/0	0/0/0/0
9	TLA	F	306	-	-	0/4/12/12	0/0/0/0
10	PGE	F	309	-	-	0/7/7/7	0/0/0/0
8	PEG	F	310	-	-	0/4/4/4	0/0/0/0
8	PEG	F	314	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	301	B3P	C10-C8-C9	-11.72	85.60	110.14
5	C	301	B3P	C6-C4-C5	-11.00	87.11	110.14
5	C	301	B3P	C7-C4-C5	-9.91	89.39	110.14
5	C	301	B3P	C9-C8-N2	-7.77	86.52	109.19
5	C	301	B3P	C5-C4-N1	-7.64	86.92	109.19
5	C	301	B3P	C11-C8-C9	-7.25	94.97	110.14
3	B	299	AZM	C1-S1-N1	-3.36	102.57	108.36
3	C	299	AZM	C1-S1-N1	-2.93	103.32	108.36
5	F	301[A]	B3P	C7-C4-C5	-2.49	104.92	110.14
5	F	301[B]	B3P	C7-C4-C5	-2.49	104.92	110.14
3	C	299	AZM	O3-C3-N4	-2.28	120.41	123.13
3	B	299	AZM	O3-C3-N4	-2.25	120.45	123.13
3	A	299	AZM	C1-S1-N1	-2.08	104.77	108.36
3	D	299	AZM	C1-S1-N1	-2.08	104.78	108.36
3	A	299	AZM	O3-C3-N4	-2.01	120.73	123.13
3	E	299	AZM	C1-S1-N1	-2.00	104.91	108.36
3	C	299	AZM	O2-S1-N1	2.08	109.97	107.28
5	C	301	B3P	C11-C8-N2	2.38	116.12	109.19
3	D	299	AZM	O2-S1-N1	2.43	110.42	107.28
3	C	299	AZM	O2-S1-O1	2.43	122.22	118.80
3	E	299	AZM	O1-S1-N1	2.45	110.45	107.28
5	C	301	B3P	C6-C4-N1	2.94	117.78	109.19
3	A	299	AZM	O2-S1-N1	2.97	111.12	107.28
3	F	299	AZM	O2-S1-N1	2.98	111.14	107.28
3	F	299	AZM	O1-S1-N1	3.63	111.98	107.28
3	B	299	AZM	O1-S1-N1	3.83	112.24	107.28
5	C	301	B3P	C7-C4-C6	4.25	119.04	110.14
5	A	301[B]	B3P	C3-N1-C4	4.27	122.44	116.07
5	F	301[A]	B3P	C2-N2-C8	4.37	122.58	116.07
5	F	301[B]	B3P	C2-N2-C8	4.37	122.58	116.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	301	B3P	C11-C8-C10	4.60	119.77	110.14
5	C	301	B3P	C7-C4-N1	4.65	122.75	109.19
5	C	301	B3P	C10-C8-N2	5.07	123.99	109.19
5	C	301	B3P	C3-N1-C4	5.15	123.74	116.07
5	E	301[B]	B3P	C2-N2-C8	5.30	123.96	116.07
5	E	301[A]	B3P	C2-N2-C8	5.30	123.96	116.07
5	E	301[B]	B3P	C3-N1-C4	5.34	124.02	116.07
5	E	301[A]	B3P	C3-N1-C4	5.34	124.02	116.07
5	A	301[A]	B3P	C3-N1-C4	5.39	124.10	116.07
5	F	301[A]	B3P	C3-N1-C4	5.80	124.71	116.07
5	F	301[B]	B3P	C3-N1-C4	5.80	124.71	116.07
5	D	301	B3P	C2-N2-C8	5.90	124.86	116.07
5	A	301[B]	B3P	C2-N2-C8	6.15	125.24	116.07
5	A	301[A]	B3P	C2-N2-C8	6.15	125.24	116.07
5	D	301	B3P	C3-N1-C4	6.28	125.42	116.07

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

16 monomers are involved in 55 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	302	SO4	1	0
8	A	305	PEG	1	0
10	B	305	PGE	5	0
8	B	310	PEG	1	0
11	B	312	PG5	31	0
8	B	314	PEG	1	0
3	C	299	AZM	1	0
5	C	301	B3P	3	0
6	C	302	SO4	1	0
8	C	314	PEG	1	0
10	D	305	PGE	1	0
10	E	305	PGE	1	0
8	E	313	PEG	2	0
8	E	314	PEG	1	0
8	F	310	PEG	2	0
8	F	314	PEG	2	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	225/247 (91%)	-0.14	5 (2%) 65 63	14, 24, 42, 65	0
1	B	225/247 (91%)	-0.24	1 (0%) 93 92	13, 22, 39, 61	0
1	C	225/247 (91%)	-0.17	3 (1%) 79 79	14, 26, 45, 71	0
1	D	225/247 (91%)	-0.11	4 (1%) 71 71	14, 26, 43, 67	0
1	E	225/247 (91%)	-0.20	3 (1%) 79 79	13, 24, 43, 58	0
1	F	225/247 (91%)	-0.33	2 (0%) 85 85	12, 21, 36, 61	0
All	All	1350/1482 (91%)	-0.20	18 (1%) 79 79	12, 24, 43, 71	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	175	THR	4.0
1	C	175	THR	3.2
1	C	173[A]	HIS	3.2
1	D	23	GLY	3.0
1	A	175	THR	2.6
1	D	174	LEU	2.5
1	E	175	THR	2.4
1	A	42	PRO	2.3
1	A	96[A]	ARG	2.2
1	E	176	ALA	2.2
1	F	42	PRO	2.2
1	D	96[A]	ARG	2.2
1	F	23	GLY	2.1
1	B	42	PRO	2.1
1	C	42	PRO	2.1
1	E	42	PRO	2.1
1	A	176	ALA	2.0
1	A	23	GLY	2.0

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
8	PEG	F	314	7/7	0.89	0.18	16.86	25,30,41,43	7
8	PEG	B	314	7/7	0.87	0.19	15.34	36,53,64,68	0
5	B3P	C	301	19/19	0.67	0.43	13.10	38,46,53,54	19
5	B3P	A	301[B]	19/19	0.88	0.23	10.16	25,33,48,48	19
5	B3P	A	301[A]	19/19	0.88	0.23	10.01	22,32,48,48	19
5	B3P	E	301[A]	19/19	0.86	0.22	6.55	26,30,46,49	19
5	B3P	D	301	19/19	0.70	0.42	5.63	32,40,52,54	19
4	PE3	B	300	21/43	0.90	0.20	5.30	38,43,64,68	0
4	PE3	A	300	21/43	0.89	0.17	4.56	35,41,63,66	0
9	TLA	B	306	10/10	0.85	0.23	4.45	20,25,27,30	10
8	PEG	E	314	7/7	0.86	0.20	4.02	26,37,47,50	7
4	PE3	F	300	21/43	0.92	0.15	3.95	33,40,62,67	0
8	PEG	A	305	7/7	0.87	0.17	3.68	33,35,43,52	7
5	B3P	E	301[B]	19/19	0.86	0.22	3.59	23,34,46,49	19
10	PGE	E	305	10/10	0.81	0.17	3.18	33,42,47,51	10
9	TLA	F	306	10/10	0.90	0.16	2.86	20,24,25,25	10
7	CL	A	304	1/1	0.99	0.16	2.79	24,24,24,24	1
10	PGE	B	305	10/10	0.78	0.17	2.77	34,43,54,59	10
11	PG5	B	312	12/12	0.85	0.21	2.57	33,40,44,47	12
8	PEG	C	314	7/7	0.85	0.18	2.35	46,61,79,85	0
10	PGE	D	305	10/10	0.75	0.24	2.26	36,49,55,64	10
9	TLA	A	306	10/10	0.90	0.15	2.15	20,25,27,28	10
10	PGE	F	309	10/10	0.76	0.18	2.02	34,38,42,42	10
5	B3P	F	301[B]	19/19	0.91	0.12	1.79	27,33,47,52	2

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	AZM	F	299	13/13	0.98	0.10	1.74	17,25,44,45	0
6	SO4	F	302	5/5	0.96	0.14	1.58	50,51,52,57	5
10	PGE	C	305	10/10	0.87	0.21	1.52	43,47,56,56	10
5	B3P	F	301[A]	19/19	0.91	0.12	1.51	26,33,47,52	2
3	AZM	B	299	13/13	0.98	0.10	1.29	17,29,54,56	0
3	AZM	A	299	13/13	0.99	0.10	0.88	21,28,42,43	0
6	SO4	D	302	5/5	0.95	0.13	0.62	71,73,74,79	0
6	SO4	C	302	5/5	0.96	0.11	0.43	44,50,52,53	5
6	SO4	B	302	5/5	0.97	0.12	0.22	43,46,49,52	5
6	SO4	E	302	5/5	0.97	0.11	0.20	37,41,42,47	5
7	CL	F	308	1/1	0.94	0.11	-0.20	25,25,25,25	1
7	CL	D	307	1/1	0.76	0.13	-0.44	48,48,48,48	1
7	CL	C	303	1/1	0.91	0.17	-0.50	37,37,37,37	1
7	CL	E	307	1/1	0.91	0.11	-0.51	37,37,37,37	1
7	CL	F	304	1/1	0.98	0.07	-0.64	32,32,32,32	0
3	AZM	C	299	13/13	0.99	0.08	-0.71	17,27,36,41	0
7	CL	A	307	1/1	0.90	0.12	-0.80	37,37,37,37	1
3	AZM	E	299	13/13	0.98	0.07	-0.82	18,27,37,41	0
7	CL	C	304	1/1	0.99	0.07	-0.93	34,34,34,34	0
3	AZM	D	299	13/13	0.98	0.07	-1.04	18,29,36,36	0
7	CL	B	308	1/1	0.98	0.08	-1.05	30,30,30,30	1
7	CL	C	307	1/1	0.91	0.10	-1.06	33,33,33,33	1
7	CL	D	304	1/1	0.98	0.05	-1.61	34,34,34,34	0
7	CL	E	304	1/1	0.99	0.05	-1.90	34,34,34,34	0
7	CL	B	303	1/1	0.96	0.09	-2.22	29,29,29,29	0
7	CL	A	303	1/1	0.92	0.10	-4.72	34,34,34,34	1
2	ZN	B	298	1/1	0.99	0.07	-	17,17,17,17	0
8	PEG	F	310	7/7	0.75	0.29	-	30,32,44,45	7
2	ZN	C	298	1/1	1.00	0.07	-	20,20,20,20	0
7	CL	B	304	1/1	0.97	0.11	-	31,31,31,31	1
2	ZN	D	298	1/1	1.00	0.07	-	20,20,20,20	0
8	PEG	B	310	7/7	0.79	0.26	-	24,29,32,38	7
2	ZN	A	298	1/1	1.00	0.07	-	19,19,19,19	0
6	SO4	A	302	5/5	0.95	0.17	-	37,41,46,51	5
2	ZN	E	298	1/1	1.00	0.08	-	18,18,18,18	0
2	ZN	F	298	1/1	1.00	0.07	-	17,17,17,17	0
8	PEG	E	313	7/7	0.88	0.17	-	39,41,45,50	7

6.5 Other polymers ⓘ

There are no such residues in this entry.