



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 23, 2020 – 03:48 pm BST

PDB ID : 5H8L  
Title : Crystal structure of Medicago truncatula N-carbamoylputrescine amidohydrolase (MtCPA) C158S mutant in complex with putrescine  
Authors : Sekula, B.; Ruszkowski, M.; Malinska, M.; Dauter, Z.  
Deposited on : 2015-12-23  
Resolution : 2.29 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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 : 2.11  
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) : 2.11

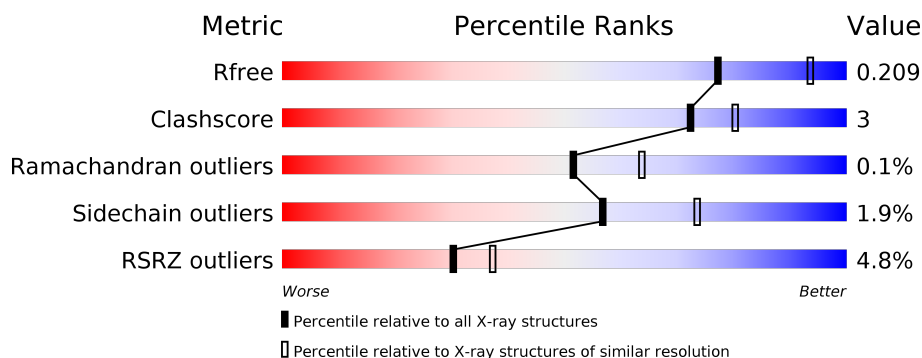
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.29 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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  $\geq 3$ , 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	304	<div> <div>7%</div> <div> <div></div> <div>81%</div> <div>13%</div> <div>••</div> </div> </div>
1	B	304	<div> <div>2%</div> <div> <div></div> <div>90%</div> <div>8%</div> <div>•</div> </div> </div>
1	C	304	<div> <div>%</div> <div> <div></div> <div>92%</div> <div>6%</div> <div>•</div> </div> </div>
1	D	304	<div> <div></div> <div> <div></div> <div>91%</div> <div>7%</div> <div>•</div> </div> </div>
1	E	304	<div> <div>%</div> <div> <div></div> <div>93%</div> <div>•••</div> </div> </div>
1	F	304	<div> <div>%</div> <div> <div></div> <div>89%</div> <div>8%</div> <div>••</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	304	
1	H	304	
1	I	304	
1	J	304	
1	K	304	
1	L	304	
1	M	304	
1	N	304	
1	O	304	
1	P	304	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	L	402	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 41665 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 N-carbamoylputrescine amidohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	295	Total	C	N	O	S	0	2	0
			2347	1501	404	433	9			
1	B	298	Total	C	N	O	S	0	1	0
			2372	1514	411	440	7			
1	C	301	Total	C	N	O	S	0	1	0
			2390	1524	412	445	9			
1	D	298	Total	C	N	O	S	0	2	0
			2372	1515	409	440	8			
1	E	297	Total	C	N	O	S	0	0	0
			2357	1505	408	437	7			
1	F	298	Total	C	N	O	S	0	1	0
			2368	1511	409	440	8			
1	G	295	Total	C	N	O	S	0	0	0
			2344	1497	405	435	7			
1	H	282	Total	C	N	O	S	0	1	0
			2248	1440	387	413	8			
1	I	292	Total	C	N	O	S	0	1	0
			2318	1481	400	429	8			
1	J	298	Total	C	N	O	S	0	1	0
			2368	1511	409	440	8			
1	K	301	Total	C	N	O	S	0	1	0
			2390	1524	412	445	9			
1	L	298	Total	C	N	O	S	0	2	0
			2372	1515	409	440	8			
1	M	297	Total	C	N	O	S	0	2	0
			2364	1511	408	437	8			
1	N	298	Total	C	N	O	S	0	0	0
			2365	1509	409	440	7			
1	O	297	Total	C	N	O	S	0	1	0
			2360	1507	408	437	8			
1	P	287	Total	C	N	O	S	0	1	0
			2284	1463	393	420	8			

There are 64 discrepancies between the modelled and reference sequences:

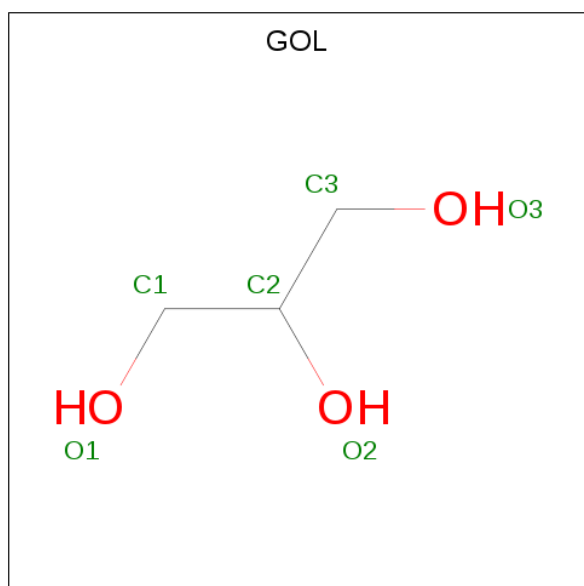
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP G7ITU5
A	-1	ASN	-	expression tag	UNP G7ITU5
A	0	ALA	-	expression tag	UNP G7ITU5
A	158	SER	CYS	engineered mutation	UNP G7ITU5
B	-2	SER	-	expression tag	UNP G7ITU5
B	-1	ASN	-	expression tag	UNP G7ITU5
B	0	ALA	-	expression tag	UNP G7ITU5
B	158	SER	CYS	engineered mutation	UNP G7ITU5
C	-2	SER	-	expression tag	UNP G7ITU5
C	-1	ASN	-	expression tag	UNP G7ITU5
C	0	ALA	-	expression tag	UNP G7ITU5
C	158	SER	CYS	engineered mutation	UNP G7ITU5
D	-2	SER	-	expression tag	UNP G7ITU5
D	-1	ASN	-	expression tag	UNP G7ITU5
D	0	ALA	-	expression tag	UNP G7ITU5
D	158	SER	CYS	engineered mutation	UNP G7ITU5
E	-2	SER	-	expression tag	UNP G7ITU5
E	-1	ASN	-	expression tag	UNP G7ITU5
E	0	ALA	-	expression tag	UNP G7ITU5
E	158	SER	CYS	engineered mutation	UNP G7ITU5
F	-2	SER	-	expression tag	UNP G7ITU5
F	-1	ASN	-	expression tag	UNP G7ITU5
F	0	ALA	-	expression tag	UNP G7ITU5
F	158	SER	CYS	engineered mutation	UNP G7ITU5
G	-2	SER	-	expression tag	UNP G7ITU5
G	-1	ASN	-	expression tag	UNP G7ITU5
G	0	ALA	-	expression tag	UNP G7ITU5
G	158	SER	CYS	engineered mutation	UNP G7ITU5
H	-2	SER	-	expression tag	UNP G7ITU5
H	-1	ASN	-	expression tag	UNP G7ITU5
H	0	ALA	-	expression tag	UNP G7ITU5
H	158	SER	CYS	engineered mutation	UNP G7ITU5
I	-2	SER	-	expression tag	UNP G7ITU5
I	-1	ASN	-	expression tag	UNP G7ITU5
I	0	ALA	-	expression tag	UNP G7ITU5
I	158	SER	CYS	engineered mutation	UNP G7ITU5
J	-2	SER	-	expression tag	UNP G7ITU5
J	-1	ASN	-	expression tag	UNP G7ITU5
J	0	ALA	-	expression tag	UNP G7ITU5
J	158	SER	CYS	engineered mutation	UNP G7ITU5
K	-2	SER	-	expression tag	UNP G7ITU5
K	-1	ASN	-	expression tag	UNP G7ITU5

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Chain	Residue	Modelled	Actual	Comment	Reference
K	0	ALA	-	expression tag	UNP G7ITU5
K	158	SER	CYS	engineered mutation	UNP G7ITU5
L	-2	SER	-	expression tag	UNP G7ITU5
L	-1	ASN	-	expression tag	UNP G7ITU5
L	0	ALA	-	expression tag	UNP G7ITU5
L	158	SER	CYS	engineered mutation	UNP G7ITU5
M	-2	SER	-	expression tag	UNP G7ITU5
M	-1	ASN	-	expression tag	UNP G7ITU5
M	0	ALA	-	expression tag	UNP G7ITU5
M	158	SER	CYS	engineered mutation	UNP G7ITU5
N	-2	SER	-	expression tag	UNP G7ITU5
N	-1	ASN	-	expression tag	UNP G7ITU5
N	0	ALA	-	expression tag	UNP G7ITU5
N	158	SER	CYS	engineered mutation	UNP G7ITU5
O	-2	SER	-	expression tag	UNP G7ITU5
O	-1	ASN	-	expression tag	UNP G7ITU5
O	0	ALA	-	expression tag	UNP G7ITU5
O	158	SER	CYS	engineered mutation	UNP G7ITU5
P	-2	SER	-	expression tag	UNP G7ITU5
P	-1	ASN	-	expression tag	UNP G7ITU5
P	0	ALA	-	expression tag	UNP G7ITU5
P	158	SER	CYS	engineered mutation	UNP G7ITU5

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 6	C 3	O 3	0	0
2	A	1	Total 6	C 3	O 3	0	0
2	B	1	Total 6	C 3	O 3	0	0
2	B	1	Total 6	C 3	O 3	0	0
2	C	1	Total 6	C 3	O 3	0	0
2	C	1	Total 6	C 3	O 3	0	0
2	D	1	Total 6	C 3	O 3	0	0
2	D	1	Total 6	C 3	O 3	0	0
2	E	1	Total 6	C 3	O 3	0	0
2	E	1	Total 6	C 3	O 3	0	0
2	E	1	Total 6	C 3	O 3	0	0
2	E	1	Total 6	C 3	O 3	0	0
2	F	1	Total 6	C 3	O 3	0	0
2	G	1	Total 6	C 3	O 3	0	0
2	G	1	Total 6	C 3	O 3	0	0
2	I	1	Total 6	C 3	O 3	0	0
2	J	1	Total 6	C 3	O 3	0	0
2	J	1	Total 6	C 3	O 3	0	0
2	J	1	Total 6	C 3	O 3	0	0
2	K	1	Total 6	C 3	O 3	0	0
2	K	1	Total 6	C 3	O 3	0	0
2	K	1	Total 6	C 3	O 3	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	L	1	Total	C	O	0	0
			6	3	3		
2	M	1	Total	C	O	0	0
			6	3	3		
2	N	1	Total	C	O	0	0
			6	3	3		
2	N	1	Total	C	O	0	0
			6	3	3		
2	N	1	Total	C	O	0	0
			6	3	3		
2	O	1	Total	C	O	0	0
			6	3	3		
2	P	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

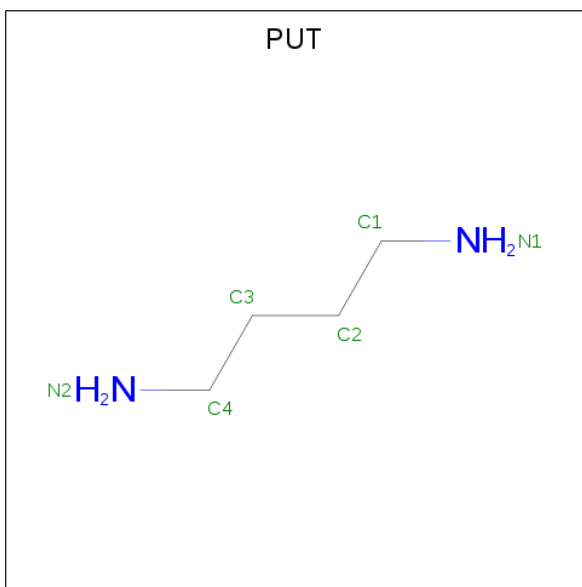
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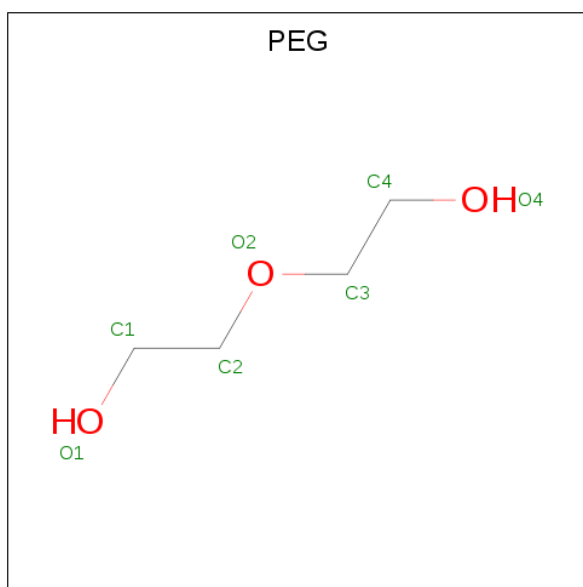
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total 4	C 2	O 2	0	0
3	D	1	Total 4	C 2	O 2	0	0
3	D	1	Total 4	C 2	O 2	0	0
3	F	1	Total 4	C 2	O 2	0	0
3	F	1	Total 4	C 2	O 2	0	0
3	I	1	Total 4	C 2	O 2	0	0
3	I	1	Total 4	C 2	O 2	0	0
3	K	1	Total 4	C 2	O 2	0	0
3	L	1	Total 4	C 2	O 2	0	0
3	L	1	Total 4	C 2	O 2	0	0
3	M	1	Total 4	C 2	O 2	0	0
3	N	1	Total 4	C 2	O 2	0	0
3	P	1	Total 4	C 2	O 2	0	0
3	P	1	Total 4	C 2	O 2	0	0

- Molecule 4 is 1,4-DIAMINOBUTANE (three-letter code: PUT) (formula: C<sub>4</sub>H<sub>12</sub>N<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	N	0	0
			6	4	2		
4	C	1	Total	C	N	0	0
			6	4	2		
4	D	1	Total	C	N	0	0
			6	4	2		
4	E	1	Total	C	N	0	0
			6	4	2		
4	F	1	Total	C	N	0	0
			6	4	2		
4	G	1	Total	C	N	0	0
			6	4	2		
4	J	1	Total	C	N	0	0
			6	4	2		
4	K	1	Total	C	N	0	0
			6	4	2		
4	L	1	Total	C	N	0	0
			6	4	2		
4	M	1	Total	C	N	0	0
			6	4	2		
4	N	1	Total	C	N	0	0
			6	4	2		
4	O	1	Total	C	N	0	0
			6	4	2		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			7	4	3		
5	K	1	Total	C	O	0	0
			7	4	3		
5	L	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	194	Total	O	0	0
			194	194		
6	B	264	Total	O	0	0
			264	264		
6	C	291	Total	O	0	0
			291	291		
6	D	312	Total	O	0	0
			312	312		
6	E	274	Total	O	0	0
			274	274		
6	F	205	Total	O	0	0
			205	205		
6	G	169	Total	O	0	0
			169	169		
6	H	61	Total	O	0	0
			61	61		
6	I	214	Total	O	0	0
			214	214		

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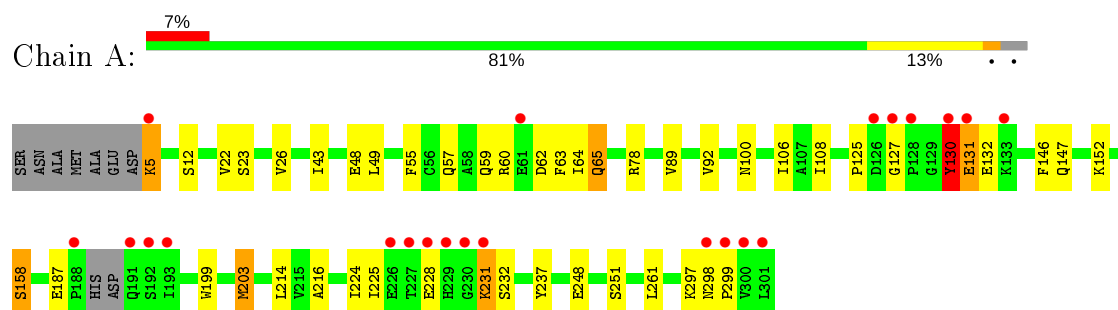
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	J	257	Total 257	O 257	0	0
6	K	270	Total 270	O 270	0	0
6	L	292	Total 292	O 292	0	0
6	M	279	Total 279	O 279	0	0
6	N	291	Total 291	O 291	0	0
6	O	226	Total 226	O 226	0	0
6	P	106	Total 106	O 106	0	0

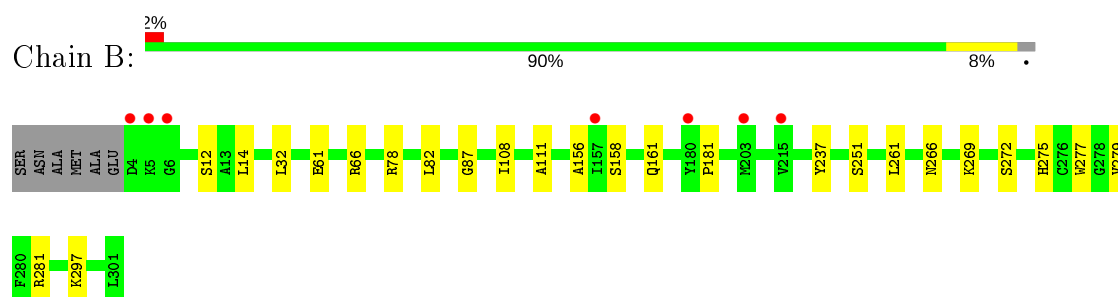
### 3 Residue-property plots [i](#)

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

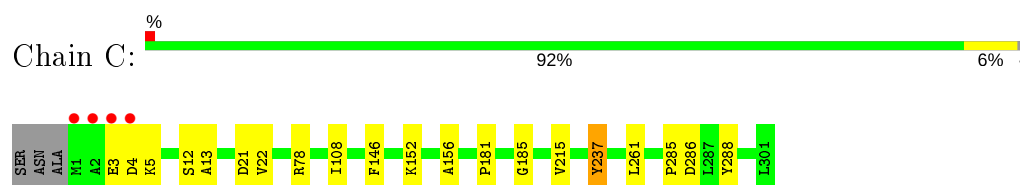
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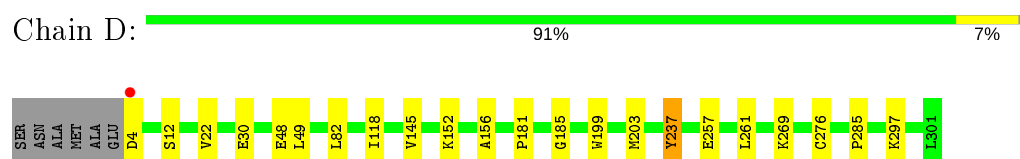
- Molecule 1: N-carbamoylputrescine amidohydrolase



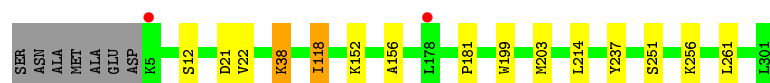
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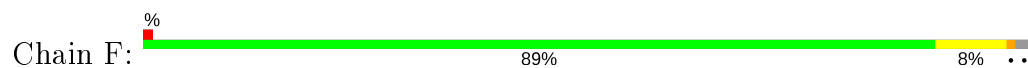
- Molecule 1: N-carbamoylputrescine amidohydrolase



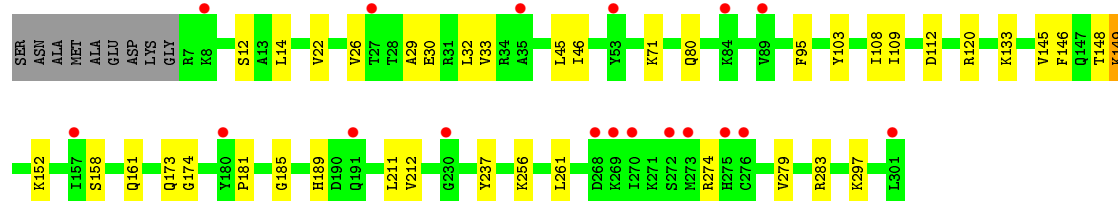
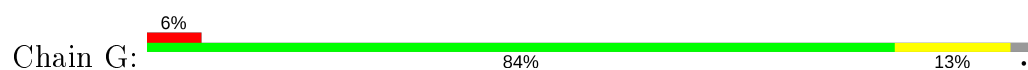
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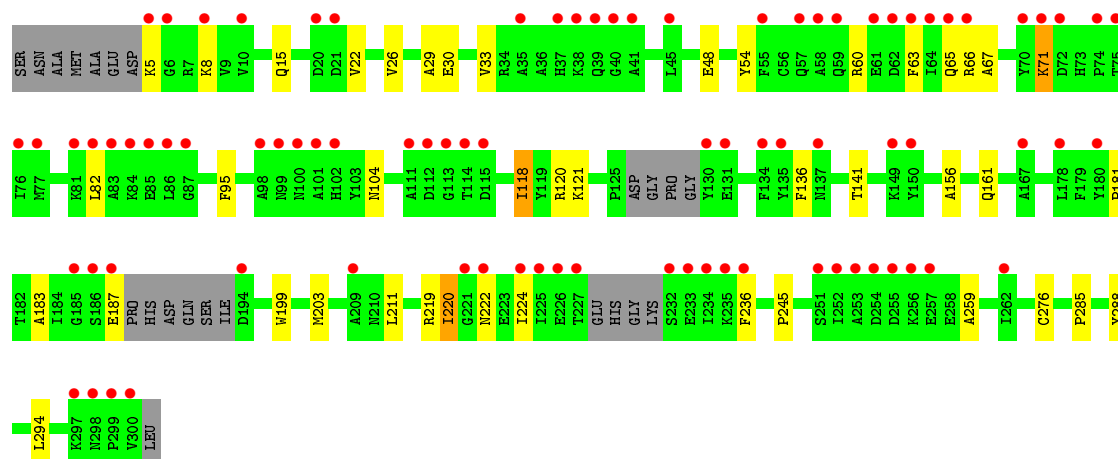
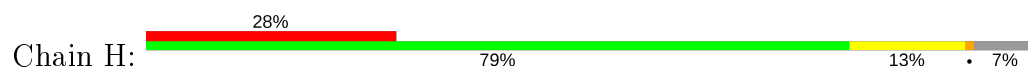
- Molecule 1: N-carbamoylputrescine amidohydrolase



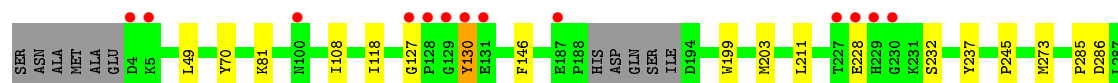
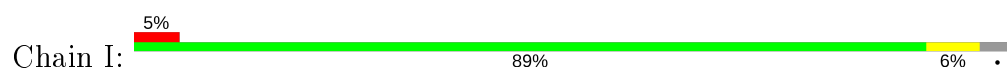
- Molecule 1: N-carbamoylputrescine amidohydrolase

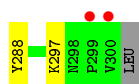


- Molecule 1: N-carbamoylputrescine amidohydrolase

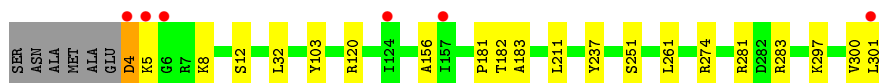
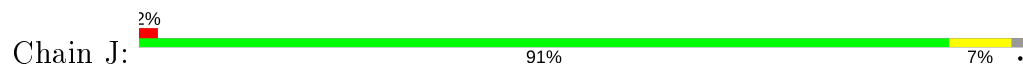


- Molecule 1: N-carbamoylputrescine amidohydrolase





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- Molecule 1: N-carbamoylputrescine amidohydrolase



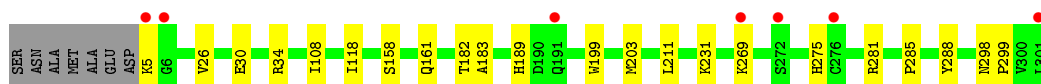
- Molecule 1: N-carbamoylputrescine amidohydrolase



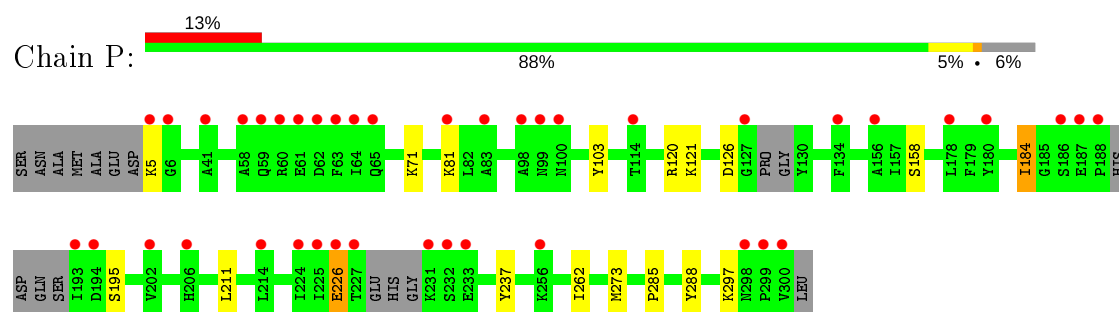
- Molecule 1: N-carbamoylputrescine amidohydrolase



- Molecule 1: N-carbamoylputrescine amidohydrolase



- Molecule 1: N-carbamoylputrescine amidohydrolase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	152.21Å 211.08Å 208.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.53 – 2.29 39.53 – 2.29	Depositor EDS
% Data completeness (in resolution range)	97.5 (39.53-2.29) 97.5 (39.53-2.29)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.61 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, $R_{free}$	0.154 , 0.206 0.163 , 0.209	Depositor DCC
$R_{free}$ test set	2944 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.1	Xtriage
Anisotropy	0.728	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 37.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	41665	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	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 41.50 % 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 2.3407e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: PUT, GOL, PEG, 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	0.83	1/2407 (0.0%)	0.81	1/3253 (0.0%)
1	B	0.85	0/2432	0.79	0/3290
1	C	0.88	0/2449	0.81	0/3312
1	D	0.86	1/2434 (0.0%)	0.80	0/3293
1	E	0.81	0/2413	0.78	1/3264 (0.0%)
1	F	0.82	0/2427	0.79	1/3283 (0.0%)
1	G	0.85	0/2400	0.78	0/3248
1	H	0.77	0/2300	0.76	0/3107
1	I	0.84	0/2375	0.79	0/3212
1	J	0.84	0/2427	0.78	0/3283
1	K	0.84	0/2449	0.79	0/3312
1	L	0.86	0/2434	0.79	0/3293
1	M	0.84	0/2426	0.78	0/3282
1	N	0.88	0/2421	0.80	0/3275
1	O	0.79	0/2419	0.77	0/3272
1	P	0.74	0/2337	0.74	0/3157
All	All	0.83	2/38550 (0.0%)	0.79	3/52136 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	130	TYR	CE1-CZ	5.68	1.46	1.38
1	D	257	GLU	CB-CG	5.26	1.62	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	118	ILE	CB-CA-C	-5.24	101.11	111.60
1	F	225	ILE	CB-CA-C	-5.01	101.57	111.60
1	A	225	ILE	CB-CA-C	-5.01	101.58	111.60

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	2347	0	2321	39	0
1	B	2372	0	2330	15	0
1	C	2390	0	2351	12	0
1	D	2372	0	2337	14	0
1	E	2357	0	2319	11	0
1	F	2368	0	2328	17	0
1	G	2344	0	2303	25	0
1	H	2248	0	2222	27	0
1	I	2318	0	2281	14	0
1	J	2368	0	2328	10	0
1	K	2390	0	2351	12	0
1	L	2372	0	2337	12	0
1	M	2364	0	2333	16	0
1	N	2365	0	2323	12	0
1	O	2360	0	2324	13	0
1	P	2284	0	2260	7	0
2	A	12	0	16	2	0
2	B	12	0	16	1	0
2	C	12	0	16	1	0
2	D	12	0	16	2	0
2	E	24	0	32	0	0
2	F	6	0	8	0	0
2	G	12	0	16	0	0
2	I	6	0	8	0	0
2	J	18	0	24	0	0
2	K	18	0	24	1	0
2	L	6	0	8	4	0
2	M	6	0	8	0	0
2	N	24	0	32	0	0
2	O	6	0	8	0	0
2	P	6	0	8	0	0
3	A	4	0	6	2	0
3	B	8	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	4	0	6	2	0
3	D	8	0	12	1	0
3	F	8	0	12	1	0
3	I	8	0	12	1	0
3	K	4	0	6	1	0
3	L	8	0	12	2	0
3	M	4	0	6	0	0
3	N	4	0	6	0	0
3	P	8	0	12	0	0
4	B	6	0	12	0	0
4	C	6	0	12	0	0
4	D	6	0	12	0	0
4	E	6	0	12	0	0
4	F	6	0	12	0	0
4	G	6	0	12	0	0
4	J	6	0	12	0	0
4	K	6	0	12	1	0
4	L	6	0	12	0	0
4	M	6	0	12	0	0
4	N	6	0	12	0	0
4	O	6	0	12	0	0
5	D	7	0	10	0	0
5	K	7	0	10	0	0
5	L	7	0	10	0	0
6	A	194	0	0	13	0
6	B	264	0	0	3	0
6	C	291	0	0	5	0
6	D	312	0	0	3	0
6	E	274	0	0	5	0
6	F	205	0	0	2	0
6	G	169	0	0	5	0
6	H	61	0	0	2	0
6	I	214	0	0	3	0
6	J	257	0	0	4	0
6	K	270	0	0	4	0
6	L	292	0	0	1	0
6	M	279	0	0	3	0
6	N	291	0	0	5	0
6	O	226	0	0	2	0
6	P	106	0	0	1	0
All	All	41665	0	37564	257	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 3.

The worst 5 of 257 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:65:GLN:HB2	6:A:653:HOH:O	1.70	0.89
1:C:21:ASP:HB2	6:C:740:HOH:O	1.73	0.88
1:N:21:ASP:HB2	6:N:739:HOH:O	1.73	0.86
1:N:269:LYS:HE2	6:N:685:HOH:O	1.79	0.83
1:N:84:LYS:HE2	6:N:660:HOH:O	1.81	0.80

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	293/304 (96%)	278 (95%)	14 (5%)	1 (0%)	41	50
1	B	297/304 (98%)	285 (96%)	11 (4%)	1 (0%)	41	50
1	C	300/304 (99%)	291 (97%)	9 (3%)	0	100	100
1	D	298/304 (98%)	290 (97%)	8 (3%)	0	100	100
1	E	295/304 (97%)	285 (97%)	10 (3%)	0	100	100
1	F	297/304 (98%)	284 (96%)	12 (4%)	1 (0%)	41	50
1	G	293/304 (96%)	281 (96%)	11 (4%)	1 (0%)	41	50
1	H	275/304 (90%)	254 (92%)	21 (8%)	0	100	100
1	I	289/304 (95%)	277 (96%)	12 (4%)	0	100	100
1	J	297/304 (98%)	286 (96%)	11 (4%)	0	100	100
1	K	300/304 (99%)	291 (97%)	9 (3%)	0	100	100
1	L	298/304 (98%)	290 (97%)	8 (3%)	0	100	100
1	M	297/304 (98%)	286 (96%)	10 (3%)	1 (0%)	41	50

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	N	296/304 (97%)	286 (97%)	9 (3%)	1 (0%)	41	50
1	O	296/304 (97%)	285 (96%)	10 (3%)	1 (0%)	41	50
1	P	280/304 (92%)	271 (97%)	9 (3%)	0	100	100
All	All	4701/4864 (97%)	4520 (96%)	174 (4%)	7 (0%)	51	64

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	158	SER
1	A	158	SER
1	N	158	SER
1	F	158	SER
1	M	158	SER

### 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	247/252 (98%)	234 (95%)	13 (5%)	22	31
1	B	249/252 (99%)	244 (98%)	5 (2%)	55	72
1	C	251/252 (100%)	247 (98%)	4 (2%)	62	78
1	D	250/252 (99%)	249 (100%)	1 (0%)	91	96
1	E	247/252 (98%)	244 (99%)	3 (1%)	71	84
1	F	249/252 (99%)	243 (98%)	6 (2%)	49	66
1	G	246/252 (98%)	239 (97%)	7 (3%)	43	60
1	H	236/252 (94%)	227 (96%)	9 (4%)	33	47
1	I	243/252 (96%)	239 (98%)	4 (2%)	62	78
1	J	249/252 (99%)	243 (98%)	6 (2%)	49	66
1	K	251/252 (100%)	248 (99%)	3 (1%)	71	84
1	L	250/252 (99%)	249 (100%)	1 (0%)	91	96
1	M	249/252 (99%)	246 (99%)	3 (1%)	71	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	N	248/252 (98%)	246 (99%)	2 (1%)	81	91
1	O	248/252 (98%)	246 (99%)	2 (1%)	81	91
1	P	240/252 (95%)	232 (97%)	8 (3%)	38	53
All	All	3953/4032 (98%)	3876 (98%)	77 (2%)	57	73

5 of 77 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	149	LYS
1	H	187	GLU
1	P	81	LYS
1	G	161	GLN
1	H	5	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 27 such sidechains are listed below:

Mol	Chain	Res	Type
1	H	161	GLN
1	J	65	GLN
1	P	99	ASN
1	H	222	ASN
1	C	65	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 carbohydrates in this entry.

## 5.6 Ligand geometry

62 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	$\# Z  > 2$
3	EDO	A	403	-	3,3,3	0.51	0	2,2,2	0.31	0
4	PUT	D	402	-	5,5,5	0.21	0	4,4,4	0.82	0
2	GOL	I	401	-	5,5,5	0.47	0	5,5,5	0.17	0
4	PUT	O	401	-	5,5,5	0.35	0	4,4,4	0.78	0
4	PUT	G	401	-	5,5,5	0.22	0	4,4,4	1.11	1 (25%)
5	PEG	D	406	-	6,6,6	0.68	0	5,5,5	0.19	0
2	GOL	P	401	-	5,5,5	0.44	0	5,5,5	0.15	0
4	PUT	C	401	-	5,5,5	0.26	0	4,4,4	0.92	0
2	GOL	G	402	-	5,5,5	0.44	0	5,5,5	0.58	0
5	PEG	L	405	-	6,6,6	0.59	0	5,5,5	0.34	0
2	GOL	J	403	-	5,5,5	0.57	0	5,5,5	0.70	0
2	GOL	O	402	-	5,5,5	0.25	0	5,5,5	0.67	0
3	EDO	B	404	-	3,3,3	0.61	0	2,2,2	0.15	0
2	GOL	F	402	-	5,5,5	0.51	0	5,5,5	0.24	0
4	PUT	K	401	-	5,5,5	0.38	0	4,4,4	0.63	0
4	PUT	L	401	-	5,5,5	0.43	0	4,4,4	0.77	0
2	GOL	N	405	-	5,5,5	0.65	0	5,5,5	0.52	0
2	GOL	L	402	-	5,5,5	0.79	0	5,5,5	0.51	0
3	EDO	M	403	-	3,3,3	0.78	0	2,2,2	0.13	0
2	GOL	C	403	-	5,5,5	0.51	0	5,5,5	0.41	0
3	EDO	K	405	-	3,3,3	0.60	0	2,2,2	0.11	0
2	GOL	B	403	-	5,5,5	0.31	0	5,5,5	0.36	0
3	EDO	L	403	-	3,3,3	0.54	0	2,2,2	0.24	0
2	GOL	E	402	-	5,5,5	0.37	0	5,5,5	0.36	0
4	PUT	J	401	-	5,5,5	0.40	0	4,4,4	0.98	0
4	PUT	N	402	-	5,5,5	0.28	0	4,4,4	0.61	0
3	EDO	I	403	-	3,3,3	0.51	0	2,2,2	0.43	0
3	EDO	P	402	-	3,3,3	0.60	0	2,2,2	0.21	0
3	EDO	P	403	-	3,3,3	0.53	0	2,2,2	0.29	0
2	GOL	K	402	-	5,5,5	0.61	0	5,5,5	0.41	0
4	PUT	B	401	-	5,5,5	0.34	0	4,4,4	0.76	0
4	PUT	M	401	-	5,5,5	0.47	0	4,4,4	0.66	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	PUT	E	401	-	5,5,5	0.18	0	4,4,4	0.84	0
3	EDO	D	405	-	3,3,3	0.50	0	2,2,2	0.35	0
2	GOL	E	404	-	5,5,5	0.36	0	5,5,5	0.32	0
2	GOL	G	403	-	5,5,5	0.46	0	5,5,5	0.24	0
3	EDO	I	402	-	3,3,3	0.65	0	2,2,2	0.05	0
4	PUT	F	401	-	5,5,5	0.36	0	4,4,4	0.46	0
2	GOL	K	404	-	5,5,5	0.53	0	5,5,5	0.27	0
3	EDO	L	404	-	3,3,3	0.55	0	2,2,2	0.12	0
2	GOL	J	404	-	5,5,5	0.34	0	5,5,5	0.30	0
2	GOL	C	402	-	5,5,5	0.43	0	5,5,5	0.47	0
2	GOL	A	402	-	5,5,5	0.47	0	5,5,5	0.23	0
2	GOL	D	401	-	5,5,5	0.42	0	5,5,5	0.31	0
2	GOL	E	405	-	5,5,5	0.58	0	5,5,5	0.43	0
2	GOL	N	404	-	5,5,5	0.46	0	5,5,5	0.24	0
5	PEG	K	406	-	6,6,6	0.59	0	5,5,5	0.14	0
3	EDO	B	405	-	3,3,3	0.69	0	2,2,2	0.02	0
2	GOL	B	402	-	5,5,5	0.69	0	5,5,5	0.59	0
2	GOL	N	401	-	5,5,5	0.23	0	5,5,5	0.23	0
2	GOL	N	403	-	5,5,5	0.78	0	5,5,5	0.64	0
2	GOL	E	403	-	5,5,5	0.48	0	5,5,5	0.44	0
2	GOL	J	402	-	5,5,5	0.38	0	5,5,5	0.39	0
2	GOL	D	403	-	5,5,5	0.79	0	5,5,5	0.51	0
2	GOL	K	403	-	5,5,5	0.65	0	5,5,5	0.47	0
3	EDO	D	404	-	3,3,3	0.63	0	2,2,2	0.24	0
3	EDO	C	404	-	3,3,3	0.71	0	2,2,2	0.42	0
3	EDO	F	404	-	3,3,3	0.59	0	2,2,2	0.30	0
2	GOL	A	401	-	5,5,5	0.59	0	5,5,5	0.64	0
2	GOL	M	402	-	5,5,5	0.40	0	5,5,5	0.21	0
3	EDO	N	406	-	3,3,3	0.61	0	2,2,2	0.02	0
3	EDO	F	403	-	3,3,3	0.61	0	2,2,2	0.21	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	EDO	A	403	-	-	1/1/1/1	-
4	PUT	D	402	-	-	0/3/3/3	-
2	GOL	I	401	-	-	2/4/4/4	-
4	PUT	O	401	-	-	0/3/3/3	-
4	PUT	G	401	-	-	0/3/3/3	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PEG	D	406	-	-	1/4/4/4	-
2	GOL	P	401	-	-	4/4/4/4	-
4	PUT	C	401	-	-	0/3/3/3	-
2	GOL	G	402	-	-	1/4/4/4	-
5	PEG	L	405	-	-	2/4/4/4	-
2	GOL	J	403	-	-	2/4/4/4	-
2	GOL	O	402	-	-	3/4/4/4	-
3	EDO	B	404	-	-	1/1/1/1	-
2	GOL	F	402	-	-	4/4/4/4	-
4	PUT	K	401	-	-	0/3/3/3	-
4	PUT	L	401	-	-	0/3/3/3	-
2	GOL	N	405	-	-	4/4/4/4	-
2	GOL	L	402	-	-	4/4/4/4	-
3	EDO	M	403	-	-	1/1/1/1	-
2	GOL	C	403	-	-	4/4/4/4	-
3	EDO	K	405	-	-	1/1/1/1	-
2	GOL	B	403	-	-	0/4/4/4	-
3	EDO	L	403	-	-	0/1/1/1	-
2	GOL	E	402	-	-	2/4/4/4	-
4	PUT	J	401	-	-	0/3/3/3	-
4	PUT	N	402	-	-	0/3/3/3	-
3	EDO	I	403	-	-	0/1/1/1	-
3	EDO	P	402	-	-	1/1/1/1	-
3	EDO	P	403	-	-	1/1/1/1	-
2	GOL	K	402	-	-	4/4/4/4	-
4	PUT	B	401	-	-	0/3/3/3	-
4	PUT	M	401	-	-	0/3/3/3	-
4	PUT	E	401	-	-	0/3/3/3	-
3	EDO	D	405	-	-	0/1/1/1	-
2	GOL	E	404	-	-	2/4/4/4	-
2	GOL	G	403	-	-	4/4/4/4	-
3	EDO	I	402	-	-	1/1/1/1	-
4	PUT	F	401	-	-	0/3/3/3	-
2	GOL	K	404	-	-	4/4/4/4	-
3	EDO	L	404	-	-	1/1/1/1	-
2	GOL	J	404	-	-	3/4/4/4	-
2	GOL	C	402	-	-	2/4/4/4	-
2	GOL	A	402	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	D	401	-	-	2/4/4/4	-
2	GOL	E	405	-	-	2/4/4/4	-
2	GOL	N	404	-	-	2/4/4/4	-
5	PEG	K	406	-	-	3/4/4/4	-
3	EDO	B	405	-	-	0/1/1/1	-
2	GOL	B	402	-	-	4/4/4/4	-
2	GOL	N	401	-	-	2/4/4/4	-
2	GOL	N	403	-	-	1/4/4/4	-
2	GOL	E	403	-	-	2/4/4/4	-
2	GOL	J	402	-	-	3/4/4/4	-
2	GOL	D	403	-	-	2/4/4/4	-
2	GOL	K	403	-	-	4/4/4/4	-
3	EDO	D	404	-	-	1/1/1/1	-
3	EDO	C	404	-	-	0/1/1/1	-
3	EDO	F	404	-	-	1/1/1/1	-
2	GOL	A	401	-	-	2/4/4/4	-
2	GOL	M	402	-	-	4/4/4/4	-
3	EDO	N	406	-	-	0/1/1/1	-
3	EDO	F	403	-	-	1/1/1/1	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	401	PUT	C2-C3-C4	-2.03	99.39	113.77

There are no chirality outliers.

5 of 100 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	I	401	GOL	O1-C1-C2-C3
2	F	402	GOL	C1-C2-C3-O3
2	N	405	GOL	O1-C1-C2-O2
2	N	405	GOL	C1-C2-C3-O3
2	L	402	GOL	O1-C1-C2-C3

There are no ring outliers.

16 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	EDO	2	0
4	K	401	PUT	1	0
2	L	402	GOL	4	0
2	C	403	GOL	1	0
3	K	405	EDO	1	0
2	B	403	GOL	1	0
3	L	403	EDO	2	0
2	K	402	GOL	1	0
3	I	402	EDO	1	0
2	D	401	GOL	1	0
3	B	405	EDO	1	0
2	D	403	GOL	1	0
3	D	404	EDO	1	0
3	C	404	EDO	2	0
2	A	401	GOL	2	0
3	F	403	EDO	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	295/304 (97%)	-0.02	22 (7%) 14 19	28, 45, 95, 132	0
1	B	298/304 (98%)	-0.20	7 (2%) 60 67	25, 36, 59, 115	0
1	C	301/304 (99%)	-0.34	4 (1%) 77 81	25, 35, 54, 112	0
1	D	298/304 (98%)	-0.44	1 (0%) 94 96	24, 33, 49, 92	0
1	E	297/304 (97%)	-0.34	2 (0%) 87 91	26, 37, 57, 92	0
1	F	298/304 (98%)	-0.41	3 (1%) 82 86	28, 44, 64, 93	0
1	G	295/304 (97%)	0.44	18 (6%) 21 27	47, 65, 84, 101	0
1	H	282/304 (92%)	1.45	85 (30%) 0 0	50, 85, 118, 136	0
1	I	292/304 (96%)	-0.16	15 (5%) 28 35	28, 41, 84, 138	0
1	J	298/304 (98%)	-0.27	6 (2%) 65 71	26, 36, 58, 118	0
1	K	301/304 (99%)	-0.28	8 (2%) 54 62	26, 37, 59, 111	0
1	L	298/304 (98%)	-0.48	1 (0%) 94 96	24, 34, 50, 91	0
1	M	297/304 (97%)	-0.34	2 (0%) 87 91	25, 35, 52, 88	0
1	N	298/304 (98%)	-0.47	3 (1%) 82 86	25, 35, 52, 92	0
1	O	297/304 (97%)	-0.21	7 (2%) 59 66	32, 46, 66, 108	0
1	P	287/304 (94%)	0.69	41 (14%) 2 3	34, 61, 104, 119	0
All	All	4732/4864 (97%)	-0.09	225 (4%) 30 37	24, 40, 87, 138	0

The worst 5 of 225 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	58	ALA	7.4
1	H	300	VAL	7.3
1	J	4	ASP	6.9
1	B	4	ASP	6.3
1	A	230	GLY	6.3

## 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 carbohydrates in this entry.

## 6.4 Ligands [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	EDO	B	404	4/4	0.65	0.20	71,72,89,91	0
3	EDO	N	406	4/4	0.72	0.29	57,58,64,65	0
3	EDO	M	403	4/4	0.78	0.21	62,65,71,71	0
2	GOL	N	405	6/6	0.78	0.29	62,69,85,89	0
2	GOL	E	405	6/6	0.81	0.35	71,83,91,92	0
3	EDO	P	402	4/4	0.81	0.22	69,80,85,86	0
2	GOL	A	402	6/6	0.82	0.23	59,66,75,98	0
5	PEG	L	405	7/7	0.83	0.37	62,68,83,83	0
3	EDO	D	404	4/4	0.83	0.27	62,65,75,75	0
2	GOL	A	401	6/6	0.83	0.19	55,74,75,75	0
3	EDO	L	403	4/4	0.83	0.33	59,59,63,63	0
2	GOL	P	401	6/6	0.84	0.26	75,79,82,83	0
5	PEG	D	406	7/7	0.84	0.26	57,72,84,89	0
3	EDO	L	404	4/4	0.84	0.20	70,78,84,85	0
2	GOL	J	403	6/6	0.84	0.23	62,64,74,80	0
2	GOL	J	404	6/6	0.85	0.24	61,73,77,80	0
3	EDO	D	405	4/4	0.85	0.19	69,75,75,80	0
2	GOL	G	402	6/6	0.86	0.16	79,88,91,94	0
2	GOL	D	403	6/6	0.87	0.39	48,60,69,72	0
3	EDO	I	403	4/4	0.87	0.19	55,82,83,85	0
3	EDO	C	404	4/4	0.88	0.20	52,57,63,63	0
3	EDO	K	405	4/4	0.88	0.24	51,59,62,63	0
3	EDO	B	405	4/4	0.88	0.21	49,62,74,83	0
4	PUT	G	401	6/6	0.89	0.39	66,68,72,76	0
2	GOL	F	402	6/6	0.89	0.24	57,66,70,71	0
3	EDO	F	403	4/4	0.89	0.24	54,63,64,73	0
3	EDO	F	404	4/4	0.90	0.42	71,72,75,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	PEG	K	406	7/7	0.90	0.28	52,81,94,96	0
2	GOL	N	403	6/6	0.91	0.15	44,58,71,75	0
2	GOL	L	402	6/6	0.91	0.13	45,55,58,59	0
2	GOL	B	403	6/6	0.91	0.14	53,67,70,73	0
2	GOL	O	402	6/6	0.91	0.15	59,64,68,73	0
3	EDO	P	403	4/4	0.91	0.32	75,77,79,80	0
2	GOL	E	402	6/6	0.91	0.15	51,64,78,84	0
2	GOL	I	401	6/6	0.92	0.14	54,60,69,70	0
2	GOL	K	404	6/6	0.92	0.28	53,58,64,71	0
2	GOL	B	402	6/6	0.93	0.11	49,55,62,67	0
2	GOL	E	404	6/6	0.94	0.22	49,54,60,61	0
2	GOL	G	403	6/6	0.94	0.18	46,60,62,65	0
4	PUT	F	401	6/6	0.94	0.32	59,62,67,67	0
2	GOL	N	404	6/6	0.94	0.20	44,51,60,61	0
2	GOL	C	403	6/6	0.94	0.22	43,57,66,67	0
3	EDO	I	402	4/4	0.94	0.19	53,58,60,69	0
4	PUT	K	401	6/6	0.94	0.28	41,42,45,45	0
2	GOL	K	403	6/6	0.95	0.13	45,52,63,67	0
4	PUT	E	401	6/6	0.95	0.32	36,37,43,45	0
3	EDO	A	403	4/4	0.95	0.16	52,61,61,65	0
4	PUT	O	401	6/6	0.95	0.28	47,48,50,50	0
2	GOL	K	402	6/6	0.95	0.17	48,59,59,69	0
4	PUT	B	401	6/6	0.95	0.28	39,40,40,40	0
2	GOL	D	401	6/6	0.95	0.15	49,57,75,78	0
4	PUT	J	401	6/6	0.96	0.33	38,40,42,43	0
4	PUT	M	401	6/6	0.96	0.32	34,35,37,39	0
2	GOL	E	403	6/6	0.96	0.17	62,71,87,99	0
2	GOL	J	402	6/6	0.96	0.10	45,52,58,66	0
2	GOL	M	402	6/6	0.96	0.16	44,54,70,70	0
2	GOL	C	402	6/6	0.96	0.14	45,52,69,77	0
4	PUT	N	402	6/6	0.96	0.36	47,47,49,51	0
2	GOL	N	401	6/6	0.97	0.11	46,58,69,73	0
4	PUT	D	402	6/6	0.97	0.27	37,41,44,45	0
4	PUT	L	401	6/6	0.97	0.32	41,45,50,50	0
4	PUT	C	401	6/6	0.97	0.30	36,38,41,44	0

## 6.5 Other polymers

There are no such residues in this entry.