



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 06:37 PM UTC

PDB ID : 5H8J / pdb_00005h8j
Title : Crystal structure of *Medicago truncatula* N-carbamoylputrescine amidohydrolase (MtCPA) in complex with cadaverine
Authors : Sekula, B.; Ruszkowski, M.; Malinska, M.; Dauter, Z.
Deposited on : 2015-12-23
Resolution : 2.19 Å(reported)

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

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

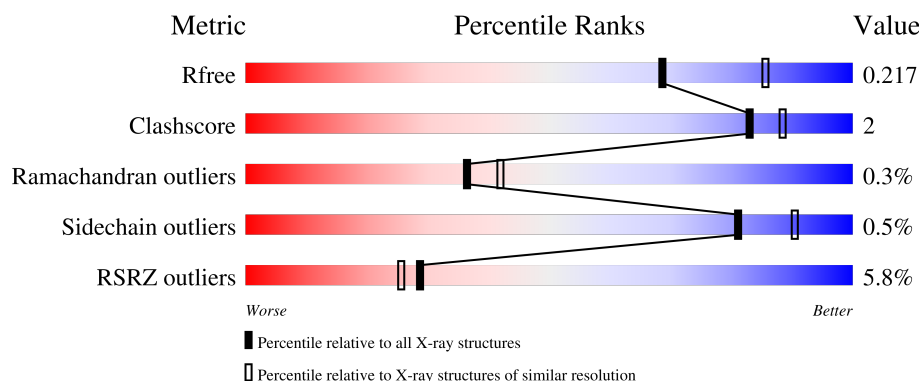
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	304	
1	B	304	
1	C	304	
1	D	304	
1	E	304	

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Mol	Chain	Length	Quality of chain
1	F	304	% 91% 6% ..
1	G	304	18% 86% 11% ..
1	H	304	43% 77% 11% 12%
1	I	304	4% 85% 11% .
1	J	304	% 91% 6% ..
1	K	304	% 94% ..
1	L	304	93% 5% .
1	M	304	% 91% 6% .
1	N	304	% 94% ..
1	O	304	2% 90% 8% .
1	P	304	9% 87% 5% 8%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 41003 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	296	Total	C	N	O	S	0	3	0
			2360	1508	407	435	10			
1	B	297	Total	C	N	O	S	0	0	0
			2357	1505	408	436	8			
1	C	301	Total	C	N	O	S	0	1	0
			2390	1524	412	444	10			
1	D	298	Total	C	N	O	S	0	3	0
			2382	1522	409	443	8			
1	E	297	Total	C	N	O	S	0	0	0
			2357	1505	408	436	8			
1	F	298	Total	C	N	O	S	0	0	0
			2365	1509	409	439	8			
1	G	295	Total	C	N	O	S	0	0	0
			2344	1497	405	434	8			
1	H	268	Total	C	N	O	S	0	0	0
			2131	1368	371	384	8			
1	I	292	Total	C	N	O	S	0	1	0
			2322	1484	400	429	9			
1	J	298	Total	C	N	O	S	0	1	0
			2369	1513	409	439	8			
1	K	301	Total	C	N	O	S	0	0	0
			2387	1522	412	444	9			
1	L	298	Total	C	N	O	S	0	2	0
			2374	1515	409	441	9			
1	M	298	Total	C	N	O	S	0	0	0
			2365	1509	409	439	8			
1	N	298	Total	C	N	O	S	0	1	0
			2369	1513	409	439	8			
1	O	298	Total	C	N	O	S	0	1	0
			2371	1513	409	441	8			
1	P	279	Total	C	N	O	S	0	1	0
			2217	1421	384	403	9			

There are 48 discrepancies between the modelled and reference sequences:

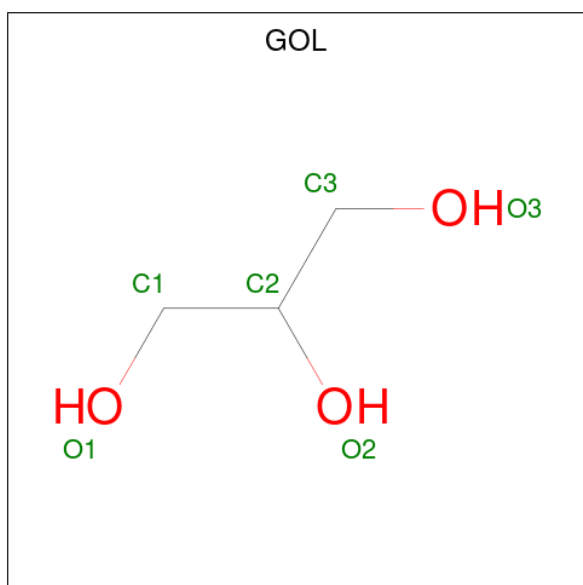
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP G7ITU5
A	-1	ASN	-	expression tag	UNP G7ITU5
A	0	ALA	-	expression tag	UNP G7ITU5
B	-2	SER	-	expression tag	UNP G7ITU5
B	-1	ASN	-	expression tag	UNP G7ITU5
B	0	ALA	-	expression tag	UNP G7ITU5
C	-2	SER	-	expression tag	UNP G7ITU5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

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Chain	Residue	Modelled	Actual	Comment	Reference
O	-2	SER	-	expression tag	UNP G7ITU5
O	-1	ASN	-	expression tag	UNP G7ITU5
O	0	ALA	-	expression tag	UNP G7ITU5
P	-2	SER	-	expression tag	UNP G7ITU5
P	-1	ASN	-	expression tag	UNP G7ITU5
P	0	ALA	-	expression tag	UNP G7ITU5

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0

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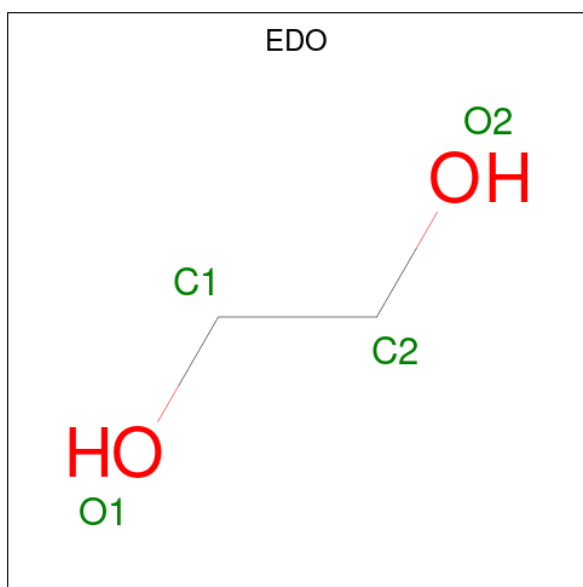
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	E	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	G	1	Total	C	O	0	0
			6	3	3		
2	H	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	K	1	Total	C	O	0	0
			6	3	3		
2	K	1	Total	C	O	0	0
			6	3	3		
2	M	1	Total	C	O	0	0
			6	3	3		
2	M	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	O	1	Total	C	O	0	0
			6	3	3		
2	O	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	G	1	Total	C	O	0	0
			7	4	3		
3	L	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



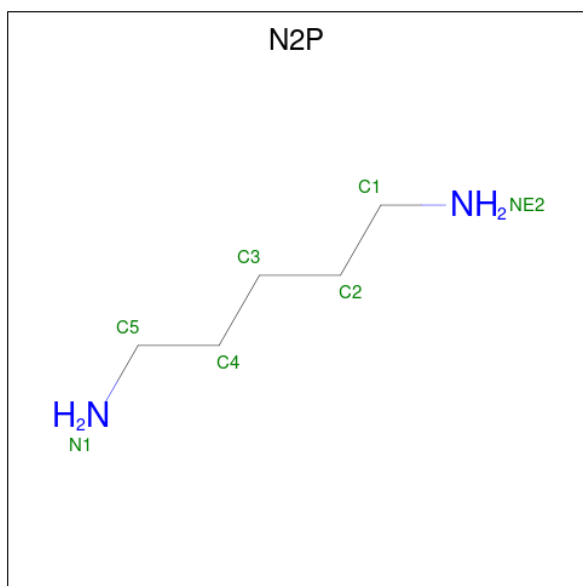
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 5 is PENTANE-1,5-DIAMINE (CCD ID: N2P) (formula: C₅H₁₄N₂).



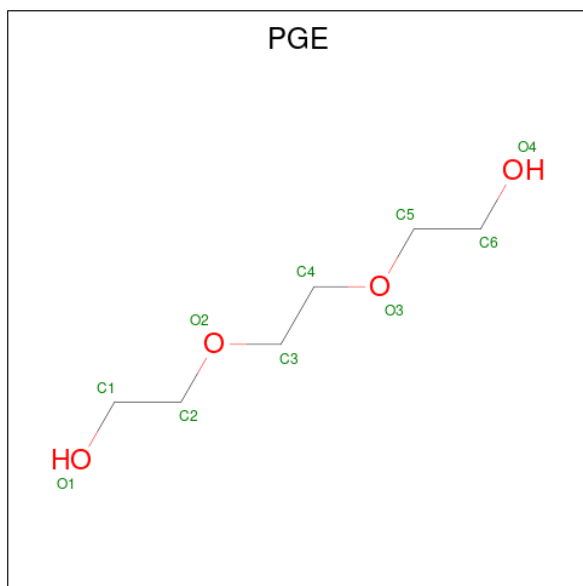
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C N 7 5 2	0	0
5	C	1	Total C N 7 5 2	0	0
5	D	1	Total C N 7 5 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	E	1	Total	C	N	0	0
			7	5	2		
5	F	1	Total	C	N	0	0
			7	5	2		
5	G	1	Total	C	N	0	0
			7	5	2		
5	J	1	Total	C	N	0	0
			7	5	2		
5	K	1	Total	C	N	0	0
			7	5	2		
5	L	1	Total	C	N	0	0
			7	5	2		
5	M	1	Total	C	N	0	0
			7	5	2		
5	N	1	Total	C	N	0	0
			7	5	2		
5	O	1	Total	C	N	0	0
			7	5	2		

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



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

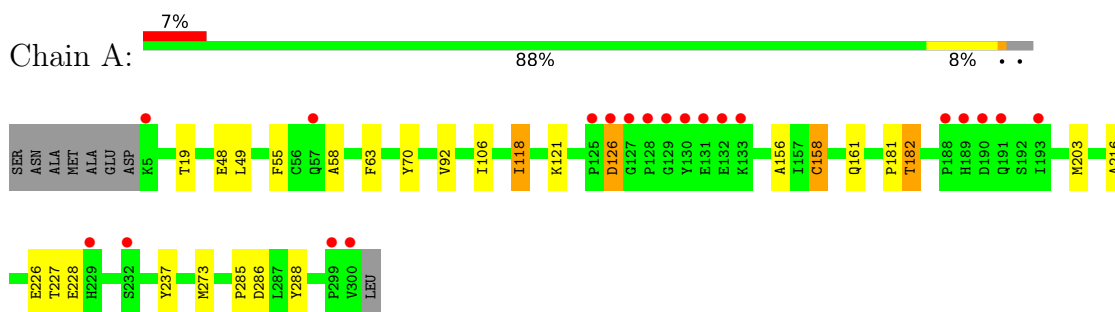
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	157	Total 157	O 157	0	0
7	B	233	Total 233	O 233	0	0
7	C	258	Total 258	O 258	0	0
7	D	288	Total 288	O 288	0	0
7	E	236	Total 236	O 236	0	0
7	F	177	Total 177	O 177	0	0
7	G	120	Total 120	O 120	0	0
7	H	38	Total 38	O 38	0	0
7	I	171	Total 171	O 171	0	0
7	J	232	Total 232	O 232	0	0
7	K	241	Total 241	O 241	0	0
7	L	277	Total 277	O 277	0	0
7	M	263	Total 263	O 263	0	0
7	N	253	Total 253	O 253	0	0
7	O	206	Total 206	O 206	0	0
7	P	92	Total 92	O 92	0	0

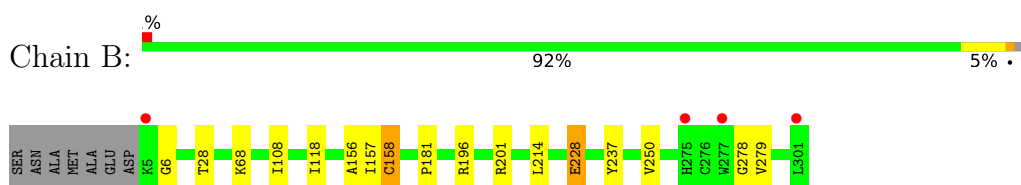
3 Residue-property plots [i](#)

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.

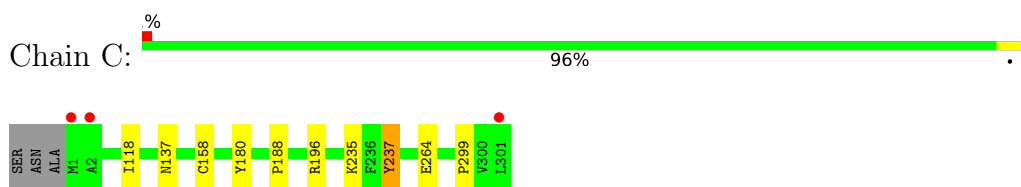
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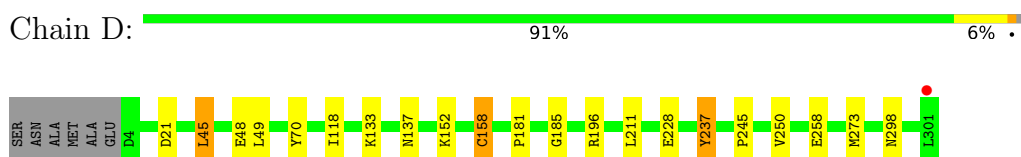
- Molecule 1: N-carbamoylputrescine amidohydrolase



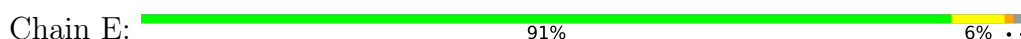
- Molecule 1: N-carbamoylputrescine amidohydrolase



- Molecule 1: N-carbamoylputrescine amidohydrolase



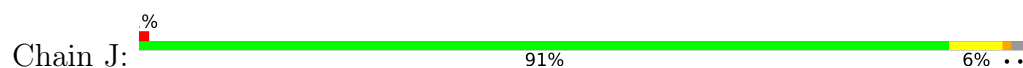
- Molecule 1: N-carbamoylputrescine amidohydrolase







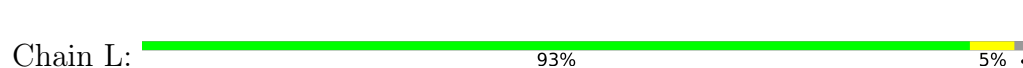
- Molecule 1: N-carbamoylputrescine amidohydrolase



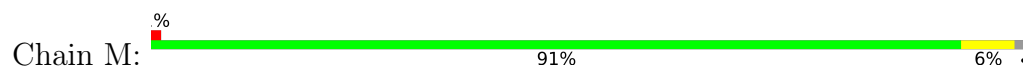
- Molecule 1: N-carbamoylputrescine amidohydrolase



- Molecule 1: N-carbamoylputrescine amidohydrolase



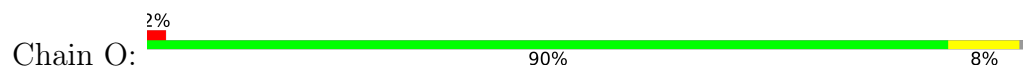
- Molecule 1: N-carbamoylputrescine amidohydrolase



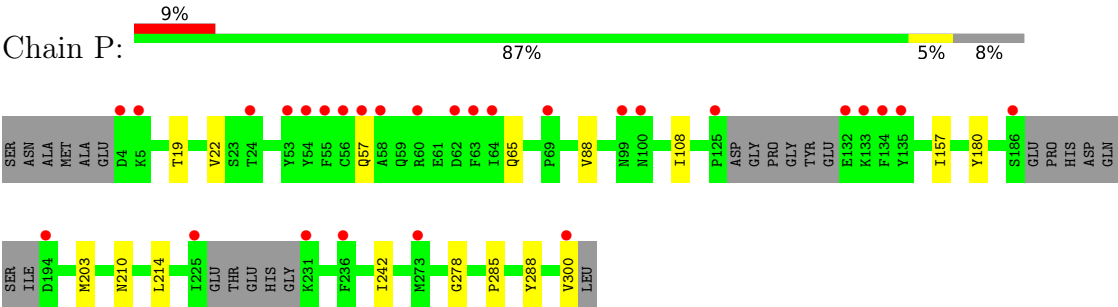
- Molecule 1: N-carbamoylputrescine amidohydrolase



- Molecule 1: N-carbamoylputrescine amidohydrolase



- 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, α , β , γ	152.27Å 211.04Å 208.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.56 – 2.19 39.56 – 2.19	Depositor EDS
% Data completeness (in resolution range)	99.1 (39.56-2.19) 99.1 (39.56-2.19)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.171 , 0.211 0.181 , 0.217	Depositor DCC
R_{free} test set	3402 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	0.469	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 51.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	41003	wwPDB-VP
Average B, all atoms (Å ²)	52.0	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 41.46 % 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 2.3540e-04. 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: N2P, GOL, PEG, EDO, PGE

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	1.17	3/2425 (0.1%)	1.02	4/3281 (0.1%)
1	B	1.22	3/2413 (0.1%)	1.02	6/3264 (0.2%)
1	C	1.21	3/2449 (0.1%)	1.01	4/3312 (0.1%)
1	D	1.28	4/2447 (0.2%)	1.06	6/3310 (0.2%)
1	E	1.18	2/2413 (0.1%)	1.01	3/3264 (0.1%)
1	F	1.12	2/2421 (0.1%)	1.00	5/3275 (0.2%)
1	G	1.29	9/2400 (0.4%)	1.05	6/3248 (0.2%)
1	H	1.18	2/2179 (0.1%)	1.03	7/2944 (0.2%)
1	I	1.22	5/2378 (0.2%)	1.05	9/3213 (0.3%)
1	J	1.25	3/2428 (0.1%)	1.04	5/3285 (0.2%)
1	K	1.17	2/2443 (0.1%)	1.00	4/3304 (0.1%)
1	L	1.24	4/2436 (0.2%)	1.03	5/3295 (0.2%)
1	M	1.19	3/2421 (0.1%)	1.02	4/3275 (0.1%)
1	N	1.22	0/2428	1.01	5/3285 (0.2%)
1	O	1.24	5/2430 (0.2%)	1.00	6/3287 (0.2%)
1	P	1.07	2/2268 (0.1%)	0.96	4/3063 (0.1%)
All	All	1.20	52/38379 (0.1%)	1.02	83/51905 (0.2%)

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	ASP	CA-C	8.98	1.61	1.53
1	L	249	ILE	C-O	-8.54	1.15	1.24
1	K	252	ILE	C-O	-7.34	1.15	1.23
1	O	215	VAL	C-O	7.18	1.31	1.24
1	J	180	TYR	CA-C	6.74	1.59	1.52
1	L	241	PHE	C-O	6.65	1.31	1.23
1	A	126	ASP	N-CA	6.35	1.53	1.46
1	G	138	PRO	C-O	6.31	1.30	1.23
1	K	250	VAL	C-O	-6.08	1.18	1.24
1	G	91	PRO	CA-C	6.07	1.58	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	180	TYR	C-O	-6.02	1.19	1.24
1	G	73	HIS	CA-C	6.01	1.59	1.52
1	D	21	ASP	CA-C	5.89	1.59	1.52
1	C	188	PRO	C-O	-5.83	1.16	1.24
1	M	279	VAL	N-CA	-5.81	1.39	1.46
1	F	167	ALA	C-O	-5.71	1.17	1.24
1	H	172	LEU	N-CA	5.71	1.53	1.46
1	B	201	ARG	C-O	-5.65	1.17	1.24
1	L	64	ILE	C-O	-5.65	1.17	1.24
1	B	250	VAL	C-O	-5.63	1.18	1.24
1	O	249	ILE	C-O	-5.60	1.18	1.24
1	O	31	ARG	N-CA	5.58	1.53	1.46
1	I	126	ASP	N-CA	5.56	1.53	1.46
1	G	260	VAL	CA-C	5.55	1.58	1.52
1	E	46	ILE	CA-C	5.54	1.59	1.52
1	D	45	LEU	CA-C	5.50	1.59	1.52
1	I	93	SER	C-O	-5.50	1.17	1.23
1	P	242	ILE	C-O	5.50	1.29	1.24
1	O	241	PHE	C-O	5.48	1.30	1.23
1	B	118	ILE	C-O	-5.37	1.18	1.23
1	I	281	ARG	N-CA	-5.36	1.39	1.46
1	I	284	ARG	CA-C	5.31	1.58	1.53
1	G	107	ALA	CA-C	5.29	1.59	1.52
1	M	46	ILE	CA-C	5.28	1.59	1.52
1	O	138	PRO	C-O	5.28	1.29	1.23
1	J	216	ALA	N-CA	5.26	1.52	1.46
1	F	21	ASP	CA-C	5.26	1.59	1.52
1	A	286	ASP	CA-C	5.21	1.59	1.52
1	C	264	GLU	C-O	-5.17	1.17	1.24
1	G	213	PRO	C-O	5.17	1.29	1.23
1	M	99	ASN	C-O	-5.15	1.19	1.24
1	G	158	CYS	N-CA	5.14	1.52	1.46
1	I	164	PRO	C-O	-5.14	1.17	1.24
1	D	250	VAL	C-O	-5.12	1.19	1.24
1	E	289	LYS	C-O	-5.12	1.18	1.24
1	G	138	PRO	CA-C	5.11	1.59	1.52
1	D	298	ASN	C-O	-5.05	1.17	1.24
1	J	101	ALA	N-CA	-5.05	1.39	1.45
1	P	210	ASN	CA-C	5.03	1.59	1.52
1	G	26	VAL	CA-C	5.01	1.59	1.52
1	H	206	HIS	N-CA	5.01	1.52	1.46
1	L	266	ASN	CA-C	5.01	1.59	1.52

All (83) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	5	LYS	N-CA-C	7.53	121.30	110.24
1	H	6	GLY	N-CA-C	7.32	120.48	112.29
1	G	92	VAL	N-CA-C	6.81	117.22	106.88
1	K	237	TYR	N-CA-C	6.71	121.08	113.02
1	D	258	GLU	N-CA-C	-6.64	100.46	110.28
1	P	157	ILE	N-CA-C	6.52	117.51	108.12
1	L	258	GLU	N-CA-C	-6.40	101.00	110.48
1	C	196	ARG	N-CA-C	6.39	118.25	111.28
1	B	237	TYR	N-CA-C	6.37	120.75	112.92
1	I	152	LYS	N-CA-C	-6.27	99.00	108.96
1	E	118	ILE	N-CA-C	6.21	119.20	108.95
1	J	196	ARG	N-CA-C	6.19	118.82	111.33
1	M	5	LYS	N-CA-C	6.12	118.45	110.43
1	H	23	SER	N-CA-C	6.11	117.81	111.03
1	B	118	ILE	N-CA-C	6.09	119.00	108.95
1	J	237	TYR	N-CA-C	6.05	120.37	112.92
1	G	137	ASN	N-CA-C	-6.00	102.21	109.83
1	D	196	ARG	N-CA-C	5.96	117.78	111.28
1	D	237	TYR	N-CA-C	5.91	120.19	112.92
1	E	19	THR	N-CA-C	-5.91	100.67	109.94
1	O	244	GLY	CA-C-N	5.87	125.55	119.56
1	O	244	GLY	C-N-CA	5.87	125.55	119.56
1	G	104	ASN	N-CA-C	-5.84	98.21	108.20
1	D	298	ASN	N-CA-C	5.81	119.15	108.69
1	B	28	THR	N-CA-C	-5.80	105.04	111.36
1	B	6	GLY	N-CA-C	5.80	118.00	111.85
1	I	157	ILE	N-CA-C	5.70	116.96	108.46
1	D	137	ASN	N-CA-C	-5.67	102.63	109.83
1	P	278	GLY	N-CA-C	5.64	123.11	115.32
1	H	135	TYR	N-CA-C	5.64	118.84	108.58
1	B	196	ARG	N-CA-C	5.60	117.38	111.28
1	N	118	ILE	CB-CA-C	-5.58	102.86	110.90
1	L	136	PHE	N-CA-C	5.56	118.02	109.07
1	O	258	GLU	N-CA-C	-5.54	101.44	110.20
1	A	19	THR	N-CA-C	-5.54	101.80	110.17
1	H	298	ASN	N-CA-C	5.52	116.91	109.24
1	H	118	ILE	N-CA-C	5.50	115.81	108.11
1	M	136	PHE	N-CA-C	5.50	117.86	108.90
1	M	157	ILE	N-CA-C	5.50	116.03	108.12
1	G	76	ILE	N-CA-C	-5.47	105.17	110.42
1	N	298	ASN	N-CA-C	5.44	116.81	109.24
1	H	48	GLU	CB-CA-C	-5.44	101.18	109.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	136	PHE	N-CA-C	5.44	117.76	108.90
1	O	196	ARG	N-CA-C	5.42	117.18	111.28
1	C	118	ILE	N-CA-C	5.36	117.80	108.95
1	M	237	TYR	N-CA-C	5.36	118.83	112.72
1	I	19	THR	N-CA-C	-5.33	102.13	110.17
1	I	56	CYS	N-CA-C	5.31	119.38	113.01
1	K	252	ILE	N-CA-C	5.30	115.35	108.35
1	C	237	TYR	N-CA-C	5.29	119.42	112.92
1	F	187	GLU	CA-C-N	5.27	125.08	119.28
1	F	187	GLU	C-N-CA	5.27	125.08	119.28
1	K	22	VAL	N-CA-C	5.27	115.48	110.42
1	P	19	THR	N-CA-C	-5.26	102.67	110.46
1	A	118	ILE	CB-CA-C	-5.26	102.56	110.55
1	F	298	ASN	N-CA-C	5.26	116.55	109.24
1	N	136	PHE	N-CA-C	5.25	117.46	108.90
1	L	104	ASN	N-CA-C	-5.22	99.27	108.20
1	I	118	ILE	N-CA-C	5.21	117.36	108.86
1	H	278	GLY	N-CA-C	5.21	122.28	115.36
1	D	152	LYS	N-CA-C	-5.19	100.70	108.96
1	J	137	ASN	CA-C-N	-5.19	114.41	119.76
1	J	137	ASN	C-N-CA	-5.19	114.41	119.76
1	C	137	ASN	N-CA-C	-5.19	103.05	109.64
1	J	118	ILE	N-CA-C	5.17	117.48	108.95
1	L	259	ALA	N-CA-C	5.15	116.89	108.76
1	A	237	TYR	N-CA-C	5.15	119.25	112.92
1	G	137	ASN	CA-C-N	-5.14	114.62	119.76
1	G	137	ASN	C-N-CA	-5.14	114.62	119.76
1	B	157	ILE	N-CA-C	5.13	115.50	108.12
1	K	288	TYR	N-CA-C	5.13	119.22	112.92
1	L	196	ARG	N-CA-C	5.13	116.87	111.28
1	E	136	PHE	N-CA-C	5.12	117.24	108.90
1	F	237	TYR	N-CA-C	5.10	118.54	112.72
1	A	118	ILE	N-CA-C	5.09	117.35	108.95
1	I	232	SER	N-CA-C	-5.09	100.87	108.96
1	O	240	SER	N-CA-C	-5.08	103.44	110.35
1	F	157	ILE	N-CA-C	5.07	115.42	108.12
1	I	195	SER	N-CA-C	5.07	119.15	112.92
1	P	22	VAL	N-CA-C	5.05	115.27	110.42
1	O	104	ASN	N-CA-C	-5.02	100.33	108.41
1	N	237	TYR	N-CA-C	5.01	119.34	112.88
1	N	104	ASN	N-CA-C	-5.01	99.64	108.20

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	2360	0	2329	14	0
1	B	2357	0	2319	9	0
1	C	2390	0	2351	5	0
1	D	2382	0	2346	12	0
1	E	2357	0	2319	12	0
1	F	2365	0	2323	13	0
1	G	2344	0	2303	17	0
1	H	2131	0	2113	17	0
1	I	2322	0	2289	14	0
1	J	2369	0	2332	12	0
1	K	2387	0	2346	11	0
1	L	2374	0	2334	8	0
1	M	2365	0	2323	14	0
1	N	2369	0	2332	11	0
1	O	2371	0	2329	13	0
1	P	2217	0	2198	6	0
2	A	18	0	24	0	0
2	B	6	0	8	1	0
2	C	18	0	24	0	0
2	D	6	0	8	0	0
2	E	12	0	16	0	0
2	F	12	0	16	1	0
2	G	6	0	8	0	0
2	H	6	0	8	0	0
2	I	18	0	24	0	0
2	K	12	0	16	0	0
2	M	18	0	24	0	0
2	N	6	0	8	0	0
2	O	12	0	16	0	0
3	A	7	0	10	0	0
3	G	7	0	10	0	0
3	L	7	0	10	0	0
4	A	4	0	6	0	0
4	C	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	4	0	6	0	0
4	F	4	0	6	0	0
4	I	4	0	6	0	0
4	J	4	0	6	0	0
4	K	4	0	6	0	0
4	L	4	0	6	0	0
4	N	4	0	6	0	0
5	B	7	0	14	3	0
5	C	7	0	14	3	0
5	D	7	0	14	3	0
5	E	7	0	14	3	0
5	F	7	0	14	3	0
5	G	7	0	14	2	0
5	J	7	0	14	3	0
5	K	7	0	14	3	0
5	L	7	0	14	3	0
5	M	7	0	14	2	0
5	N	7	0	14	3	0
5	O	7	0	14	3	0
6	B	10	0	14	0	0
7	A	157	0	0	1	0
7	B	233	0	0	2	0
7	C	258	0	0	2	0
7	D	288	0	0	0	0
7	E	236	0	0	1	0
7	F	177	0	0	3	0
7	G	120	0	0	0	0
7	H	38	0	0	0	0
7	I	171	0	0	0	0
7	J	232	0	0	0	0
7	K	241	0	0	0	0
7	L	277	0	0	0	0
7	M	263	0	0	0	0
7	N	253	0	0	2	0
7	O	206	0	0	0	0
7	P	92	0	0	0	0
All	All	41003	0	37352	181	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All (181) 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:M:158:CYS:SG	5:M:401:N2P:H5C1	1.98	1.04
1:F:158:CYS:SG	5:F:402:N2P:H5C1	1.96	1.04
1:O:211:LEU:HD11	1:O:277:TRP:CE3	2.09	0.87
1:J:158:CYS:SG	5:J:401:N2P:H5C1	2.14	0.87
1:M:158:CYS:SG	5:M:401:N2P:C5	2.66	0.83
1:D:158:CYS:SG	5:D:401:N2P:H5C1	2.19	0.80
1:I:245:PRO:HB2	1:I:273:MET:HE1	1.62	0.80
1:J:158:CYS:SG	5:J:401:N2P:C5	2.70	0.79
1:O:211:LEU:HD11	1:O:277:TRP:HE3	1.48	0.79
1:A:58:ALA:HB2	1:A:227:THR:HG22	1.65	0.79
1:C:158:CYS:SG	5:C:401:N2P:C5	2.72	0.78
1:B:158:CYS:SG	5:B:401:N2P:C5	2.73	0.77
1:E:158:CYS:SG	5:E:401:N2P:H5C1	2.25	0.77
1:O:211:LEU:HD13	1:O:274:ARG:HA	1.65	0.77
1:L:245:PRO:HG2	1:L:273:MET:HE1	1.67	0.76
1:O:158:CYS:SG	5:O:401:N2P:H5C1	2.26	0.75
1:F:158:CYS:SG	5:F:402:N2P:C5	2.74	0.75
1:N:158:CYS:SG	5:N:401:N2P:H5C1	2.28	0.74
1:H:118:ILE:H	1:H:118:ILE:HD13	1.52	0.73
1:A:203[A]:MET:HE1	1:A:216:ALA:HB2	1.73	0.70
1:K:245:PRO:HB2	1:K:273:MET:HE1	1.74	0.69
1:I:211:LEU:HD23	1:I:273:MET:HE3	1.74	0.68
1:B:158:CYS:SG	5:B:401:N2P:H5C1	2.34	0.67
1:D:158:CYS:SG	5:D:401:N2P:C5	2.83	0.67
1:K:211:LEU:HD23	1:K:273:MET:HE3	1.75	0.67
1:C:158:CYS:SG	5:C:401:N2P:H5C1	2.34	0.67
1:I:203[B]:MET:HE3	1:I:214:LEU:HD21	1.77	0.65
1:M:245:PRO:CG	1:M:273:MET:HE1	2.27	0.64
1:L:158:CYS:SG	5:L:401:N2P:H5C1	2.37	0.64
1:A:203[A]:MET:CE	1:A:216:ALA:HB2	2.29	0.61
1:A:55:PHE:HB2	1:A:63:PHE:CD2	2.36	0.60
1:G:158:CYS:SG	5:G:401:N2P:H5C1	2.41	0.60
1:H:170:MET:HE3	1:H:178:LEU:HD22	1.84	0.59
1:E:158:CYS:SG	5:E:401:N2P:C5	2.91	0.59
1:M:245:PRO:HB2	1:M:273:MET:HE1	1.87	0.57
1:I:203[B]:MET:CE	1:I:216:ALA:HB2	2.35	0.57
1:K:245:PRO:CB	1:K:273:MET:HE1	2.35	0.56
1:P:108:ILE:HD12	1:P:108:ILE:N	2.21	0.56
1:G:287:LEU:HD22	1:H:142:GLY:HA2	1.87	0.55
1:H:120:ARG:HD2	1:H:141:THR:HG23	1.88	0.55
1:I:121:LYS:NZ	1:I:132:GLU:OE1	2.37	0.55
1:C:158:CYS:SG	5:C:401:N2P:N1	2.80	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:245:PRO:HB2	1:N:273:MET:HE1	1.88	0.55
1:H:118:ILE:HD13	1:H:118:ILE:N	2.21	0.54
1:K:245:PRO:CG	1:K:273:MET:HE1	2.36	0.54
1:K:9:VAL:HG21	1:K:267:LEU:HD11	1.89	0.54
1:N:158:CYS:SG	5:N:401:N2P:N1	2.80	0.54
1:D:211:LEU:HD23	1:D:273:MET:HE3	1.90	0.54
1:L:158:CYS:SG	5:L:401:N2P:N1	2.80	0.54
1:M:245:PRO:HG2	1:M:273:MET:HE1	1.90	0.54
1:G:109:ILE:HD13	1:G:115:ASP:HA	1.91	0.53
1:B:158:CYS:SG	5:B:401:N2P:N1	2.82	0.53
1:N:158:CYS:SG	5:N:401:N2P:C5	2.95	0.53
1:G:108:ILE:HG13	1:G:146:PHE:CD2	2.43	0.53
1:H:15:GLN:NE2	1:H:220:ILE:HD13	2.23	0.53
1:E:158:CYS:SG	5:E:401:N2P:N1	2.82	0.52
1:H:220:ILE:HD11	1:H:255:ASP:C	2.33	0.52
1:J:277:TRP:HD1	1:J:279:VAL:HG13	1.75	0.52
1:I:161:GLN:O	1:I:203[A]:MET:HE1	2.09	0.52
1:O:158:CYS:SG	5:O:401:N2P:C5	2.96	0.52
1:K:158:CYS:SG	5:K:401:N2P:N1	2.84	0.51
1:A:121:LYS:NZ	1:A:158:CYS:HB3	2.27	0.50
1:D:245:PRO:HG2	1:D:273:MET:HE1	1.93	0.50
1:D:48:GLU:HG2	1:D:49:LEU:HG	1.93	0.50
1:F:158:CYS:SG	5:F:402:N2P:N1	2.85	0.50
1:H:72:ASP:HA	1:H:77:MET:HE3	1.93	0.50
1:I:156:ALA:O	1:I:181:PRO:HD2	2.12	0.50
1:M:245:PRO:CB	1:M:273:MET:HE1	2.41	0.50
1:L:158:CYS:SG	5:L:401:N2P:C5	3.00	0.50
1:M:228:GLU:H	1:M:228:GLU:CD	2.19	0.50
1:O:158:CYS:SG	5:O:401:N2P:N1	2.85	0.49
1:G:185:GLY:HA2	1:G:237:TYR:CD2	2.47	0.49
1:H:29:ALA:O	1:H:33:VAL:HG23	2.12	0.49
1:D:158:CYS:SG	5:D:401:N2P:N1	2.85	0.49
1:J:55:PHE:HB2	1:J:63:PHE:CD2	2.47	0.49
1:P:180:TYR:HB2	1:P:203[B]:MET:HE1	1.94	0.49
1:N:245:PRO:CG	1:N:273:MET:HE1	2.43	0.48
1:I:203[B]:MET:CE	1:I:214:LEU:HD21	2.43	0.48
1:C:235:LYS:HE2	7:C:649:HOH:O	2.14	0.48
1:D:245:PRO:HB2	1:D:273:MET:HE1	1.95	0.48
1:P:285:PRO:HA	1:P:288:TYR:CD2	2.49	0.48
2:F:401:GOL:H31	7:F:642:HOH:O	2.13	0.48
1:J:156:ALA:O	1:J:181:PRO:HD2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:156:ALA:O	1:H:181:PRO:HD2	2.13	0.48
1:G:277:TRP:HD1	1:G:279:VAL:HG13	1.78	0.48
1:N:245:PRO:HG2	1:N:273:MET:HE1	1.96	0.48
1:E:235:LYS:HE2	7:E:506:HOH:O	2.12	0.47
1:H:22:VAL:O	1:H:26:VAL:HG23	2.14	0.47
1:I:161:GLN:HB2	1:I:203[A]:MET:SD	2.54	0.47
1:B:156:ALA:O	1:B:181:PRO:HD2	2.14	0.47
1:E:48:GLU:HG2	1:E:49:LEU:HG	1.96	0.47
1:G:80:GLN:HG2	1:G:109:ILE:HD12	1.96	0.47
1:H:141:THR:HB	1:H:144:LYS:HE3	1.96	0.47
1:H:17:ALA:N	1:H:220:ILE:HG22	2.30	0.47
1:L:245:PRO:CG	1:L:273:MET:HE1	2.41	0.47
1:G:158:CYS:SG	5:G:401:N2P:N1	2.88	0.47
1:A:273:MET:HE3	7:A:648:HOH:O	2.14	0.46
1:A:161:GLN:HB2	1:A:203[B]:MET:SD	2.55	0.46
1:K:228:GLU:H	1:K:228:GLU:CD	2.23	0.46
1:E:70:TYR:CZ	1:E:118:ILE:HG23	2.51	0.46
1:F:30:GLU:HG3	1:F:82:LEU:HD22	1.97	0.46
1:F:228:GLU:HG3	1:G:128:PRO:HG3	1.97	0.46
1:K:158:CYS:SG	5:K:401:N2P:C5	3.04	0.46
1:O:156:ALA:O	1:O:181:PRO:HD2	2.16	0.46
1:B:214:LEU:C	1:B:214:LEU:HD23	2.42	0.45
1:G:73:HIS:HD2	1:G:75:THR:H	1.64	0.45
1:D:45:LEU:HD11	1:D:181:PRO:HG3	1.99	0.45
1:A:228:GLU:H	1:A:228:GLU:CD	2.25	0.45
1:I:60:ARG:HB3	1:I:63:PHE:HD2	1.82	0.45
1:E:156:ALA:O	1:E:181:PRO:HD2	2.15	0.45
1:K:158:CYS:SG	5:K:401:N2P:H5C1	2.56	0.45
1:H:182:THR:HG21	1:H:199:TRP:CZ2	2.52	0.44
1:H:285:PRO:HA	1:H:288:TYR:CD2	2.53	0.44
1:M:211:LEU:HD23	1:M:273:MET:HE3	1.98	0.44
1:F:289:LYS:NZ	7:F:505:HOH:O	2.50	0.44
1:I:80:GLN:HA	1:I:109:ILE:HG21	2.00	0.44
1:F:273:MET:HE1	7:F:677:HOH:O	2.16	0.44
1:O:285:PRO:HG2	1:P:300:VAL:HG21	1.99	0.44
1:F:285:PRO:HA	1:F:288:TYR:CD2	2.52	0.44
1:N:22:VAL:HG23	7:N:674:HOH:O	2.17	0.44
1:A:156:ALA:O	1:A:181:PRO:HD2	2.18	0.44
1:E:12:SER:HA	1:E:261:LEU:O	2.17	0.44
1:B:228:GLU:CD	1:B:228:GLU:H	2.26	0.43
1:I:285:PRO:HA	1:I:288:TYR:CD2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:108:ILE:HG13	1:J:146:PHE:CD1	2.52	0.43
1:J:158:CYS:SG	5:J:401:N2P:N1	2.92	0.43
1:J:63:PHE:O	1:J:66:ARG:HG3	2.18	0.43
1:K:176:GLU:O	1:K:213:PRO:HD2	2.18	0.43
1:A:70:TYR:CZ	1:A:118:ILE:HG23	2.53	0.43
1:B:278:GLY:O	1:B:279:VAL:C	2.60	0.43
1:M:285:PRO:HG2	1:N:300:VAL:HG11	2.00	0.43
1:J:108:ILE:HG13	1:J:146:PHE:CE1	2.53	0.43
1:J:70:TYR:CZ	1:J:118:ILE:HG23	2.54	0.43
1:O:189:HIS:H	1:O:189:HIS:CD2	2.37	0.43
1:F:45:LEU:HD11	1:F:181:PRO:HG3	2.00	0.43
1:N:228:GLU:CD	1:N:228:GLU:H	2.21	0.43
1:B:68:LYS:NZ	7:B:503:HOH:O	2.52	0.43
1:M:214:LEU:C	1:M:214:LEU:HD23	2.44	0.43
1:A:92:VAL:C	1:A:106:ILE:HG13	2.44	0.42
1:I:70:TYR:CZ	1:I:118:ILE:HG23	2.54	0.42
1:P:108:ILE:N	1:P:108:ILE:CD1	2.82	0.42
2:B:402:GOL:H32	7:B:502:HOH:O	2.18	0.42
1:L:79:LEU:HD23	1:L:79:LEU:HA	1.88	0.42
1:O:12:SER:HA	1:O:261:LEU:O	2.19	0.42
1:B:108:ILE:HD12	1:B:108:ILE:N	2.34	0.42
1:D:70:TYR:CZ	1:D:118[A]:ILE:HG23	2.54	0.42
1:A:285:PRO:HA	1:A:288:TYR:CD2	2.55	0.42
1:D:133:LYS:HB3	1:E:133:LYS:HD3	2.01	0.42
1:K:118:ILE:HD13	1:K:118:ILE:HG21	1.84	0.42
1:M:185:GLY:HA2	1:M:237:TYR:CD2	2.54	0.42
1:D:185:GLY:HA2	1:D:237:TYR:CD2	2.55	0.42
1:D:245:PRO:CG	1:D:273:MET:HE1	2.49	0.42
1:G:82:LEU:HD22	1:G:86:LEU:HD22	2.02	0.41
1:P:214:LEU:C	1:P:214:LEU:HD23	2.45	0.41
1:M:108:ILE:HG13	1:M:146:PHE:CD1	2.55	0.41
1:G:279:VAL:O	1:G:283:ARG:HG3	2.20	0.41
1:L:118:ILE:HG21	1:L:118:ILE:HD13	1.81	0.41
1:A:48:GLU:HG2	1:A:49:LEU:HG	2.02	0.41
1:F:55:PHE:HB2	1:F:63:PHE:CD2	2.54	0.41
1:G:211:LEU:HA	1:G:245:PRO:O	2.21	0.41
1:C:299:PRO:HB3	7:C:638:HOH:O	2.20	0.41
1:E:108:ILE:HG13	1:E:146:PHE:CD2	2.56	0.41
1:G:48:GLU:HG2	1:G:49:LEU:HG	2.03	0.41
1:E:214:LEU:C	1:E:214:LEU:HD23	2.45	0.41
1:G:55:PHE:HB2	1:G:63:PHE:CD2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:108:ILE:HD12	1:H:108:ILE:N	2.36	0.41
1:O:30[A]:GLU:HG2	1:O:82:LEU:HD22	2.03	0.41
1:F:293:THR:HG22	1:F:298:ASN:C	2.46	0.40
1:L:128:PRO:HG3	1:M:228:GLU:HG3	2.03	0.40
1:O:211:LEU:HA	1:O:245:PRO:O	2.21	0.40
1:A:158:CYS:HA	1:A:182[B]:THR:OG1	2.22	0.40
1:F:61:GLU:HG3	1:H:276:CYS:HB2	2.03	0.40
1:G:80:GLN:HA	1:G:109:ILE:HG21	2.03	0.40
1:J:54:TYR:CE2	1:J:56:CYS:HB2	2.55	0.40
1:E:285:PRO:HG2	1:F:300:VAL:HG11	2.04	0.40
1:G:288:TYR:HB3	1:G:291:LEU:HD12	2.03	0.40
1:I:300:VAL:HG21	1:J:285:PRO:HG2	2.04	0.40
1:N:245:PRO:CB	1:N:273:MET:HE1	2.49	0.40
1:M:45:LEU:HD11	1:M:181:PRO:HG3	2.02	0.40
1:N:21:ASP:HB2	7:N:713:HOH:O	2.21	0.40
1:O:237:TYR:O	1:O:238:GLY:C	2.64	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	297/304 (98%)	280 (94%)	16 (5%)	1 (0%)	36	42
1	B	295/304 (97%)	286 (97%)	8 (3%)	1 (0%)	36	42
1	C	300/304 (99%)	289 (96%)	11 (4%)	0	100	100
1	D	299/304 (98%)	292 (98%)	6 (2%)	1 (0%)	36	42
1	E	295/304 (97%)	284 (96%)	10 (3%)	1 (0%)	36	42
1	F	296/304 (97%)	287 (97%)	8 (3%)	1 (0%)	36	42
1	G	293/304 (96%)	281 (96%)	11 (4%)	1 (0%)	36	42

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	260/304 (86%)	245 (94%)	14 (5%)	1 (0%)	30	34
1	I	289/304 (95%)	279 (96%)	10 (4%)	0	100	100
1	J	297/304 (98%)	289 (97%)	7 (2%)	1 (0%)	36	42
1	K	299/304 (98%)	290 (97%)	8 (3%)	1 (0%)	36	42
1	L	298/304 (98%)	290 (97%)	7 (2%)	1 (0%)	36	42
1	M	296/304 (97%)	285 (96%)	10 (3%)	1 (0%)	36	42
1	N	297/304 (98%)	284 (96%)	12 (4%)	1 (0%)	36	42
1	O	297/304 (98%)	282 (95%)	14 (5%)	1 (0%)	36	42
1	P	272/304 (90%)	264 (97%)	8 (3%)	0	100	100
All	All	4680/4864 (96%)	4507 (96%)	160 (3%)	13 (0%)	36	42

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	158	CYS
1	N	158	CYS
1	E	158	CYS
1	F	158	CYS
1	G	158	CYS
1	K	158	CYS
1	M	158	CYS
1	O	158	CYS
1	A	158	CYS
1	B	158	CYS
1	D	158	CYS
1	J	158	CYS
1	L	158	CYS

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	249/252 (99%)	245 (98%)	4 (2%)	55	71
1	B	247/252 (98%)	246 (100%)	1 (0%)	84	92
1	C	251/252 (100%)	250 (100%)	1 (0%)	84	92
1	D	251/252 (100%)	250 (100%)	1 (0%)	84	92
1	E	247/252 (98%)	246 (100%)	1 (0%)	84	92
1	F	248/252 (98%)	248 (100%)	0	100	100
1	G	246/252 (98%)	245 (100%)	1 (0%)	84	92
1	H	222/252 (88%)	220 (99%)	2 (1%)	70	84
1	I	243/252 (96%)	241 (99%)	2 (1%)	73	85
1	J	249/252 (99%)	248 (100%)	1 (0%)	84	92
1	K	250/252 (99%)	248 (99%)	2 (1%)	73	85
1	L	250/252 (99%)	250 (100%)	0	100	100
1	M	248/252 (98%)	246 (99%)	2 (1%)	73	85
1	N	249/252 (99%)	248 (100%)	1 (0%)	84	92
1	O	249/252 (99%)	249 (100%)	0	100	100
1	P	232/252 (92%)	229 (99%)	3 (1%)	61	76
All	All	3931/4032 (98%)	3909 (99%)	22 (1%)	81	89

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	ASP
1	A	182[A]	THR
1	A	182[B]	THR
1	A	226	GLU
1	B	228	GLU
1	C	237	TYR
1	D	228	GLU
1	E	5	LYS
1	G	82	LEU
1	H	100	ASN
1	H	118	ILE
1	I	229	HIS
1	I	237	TYR
1	J	237	TYR

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Mol	Chain	Res	Type
1	K	228	GLU
1	K	237	TYR
1	M	237	TYR
1	M	251	SER
1	N	228	GLU
1	P	57	GLN
1	P	65	GLN
1	P	88	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (57) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	204	GLN
1	B	65	GLN
1	C	73	HIS
1	C	204	GLN
1	D	39	GLN
1	D	65	GLN
1	D	204	GLN
1	E	204	GLN
1	F	37	HIS
1	F	39	GLN
1	F	59	GLN
1	F	102	HIS
1	F	204	GLN
1	G	102	HIS
1	G	147	GLN
1	G	189	HIS
1	H	73	HIS
1	H	80	GLN
1	H	99	ASN
1	H	100	ASN
1	H	102	HIS
1	H	204	GLN
1	H	239	ASN
1	H	298	ASN
1	I	39	GLN
1	I	65	GLN
1	I	102	HIS
1	I	204	GLN
1	J	37	HIS
1	J	59	GLN

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Mol	Chain	Res	Type
1	K	39	GLN
1	K	65	GLN
1	K	73	HIS
1	K	102	HIS
1	L	37	HIS
1	L	39	GLN
1	L	65	GLN
1	L	102	HIS
1	L	204	GLN
1	M	59	GLN
1	M	204	GLN
1	N	65	GLN
1	N	73	HIS
1	N	204	GLN
1	O	37	HIS
1	O	102	HIS
1	O	173	GLN
1	O	189	HIS
1	O	222	ASN
1	O	298	ASN
1	P	39	GLN
1	P	65	GLN
1	P	73	HIS
1	P	100	ASN
1	P	137	ASN
1	P	204	GLN
1	P	298	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

50 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	GOL	M	404	-	5,5,5	0.42	0	5,5,5	0.11	0
4	EDO	K	404	-	3,3,3	0.52	0	2,2,2	0.21	0
5	N2P	F	402	-	6,6,6	0.48	0	5,5,5	0.69	0
2	GOL	B	402	-	5,5,5	0.50	0	5,5,5	0.46	0
2	GOL	F	401	-	5,5,5	0.30	0	5,5,5	0.49	0
3	PEG	L	402	-	6,6,6	0.60	0	5,5,5	0.25	0
2	GOL	K	402	-	5,5,5	0.24	0	5,5,5	0.21	0
2	GOL	O	402	-	5,5,5	0.19	0	5,5,5	0.75	0
5	N2P	L	401	-	6,6,6	0.28	0	5,5,5	0.59	0
2	GOL	F	403	-	5,5,5	0.50	0	5,5,5	0.22	0
2	GOL	D	402	-	5,5,5	0.81	0	5,5,5	0.50	0
2	GOL	M	402	-	5,5,5	0.38	0	5,5,5	0.30	0
4	EDO	I	404	-	3,3,3	0.54	0	2,2,2	0.25	0
2	GOL	I	401	-	5,5,5	0.37	0	5,5,5	0.32	0
2	GOL	G	403	-	5,5,5	0.33	0	5,5,5	0.60	0
2	GOL	I	403	-	5,5,5	0.47	0	5,5,5	0.26	0
2	GOL	C	402	-	5,5,5	0.57	0	5,5,5	0.29	0
4	EDO	N	403	-	3,3,3	0.53	0	2,2,2	0.16	0
2	GOL	C	403	-	5,5,5	0.45	0	5,5,5	0.27	0
5	N2P	G	401	-	6,6,6	0.35	0	5,5,5	0.80	0
4	EDO	L	403	-	3,3,3	0.72	0	2,2,2	0.03	0
4	EDO	D	403	-	3,3,3	0.59	0	2,2,2	0.19	0
2	GOL	H	401	-	5,5,5	0.67	0	5,5,5	0.58	0
5	N2P	J	401	-	6,6,6	0.23	0	5,5,5	0.84	0
5	N2P	C	401	-	6,6,6	0.20	0	5,5,5	0.93	0
4	EDO	J	402	-	3,3,3	0.41	0	2,2,2	0.44	0
5	N2P	B	401	-	6,6,6	0.24	0	5,5,5	1.04	0
5	N2P	E	401	-	6,6,6	0.25	0	5,5,5	0.41	0
2	GOL	E	402	-	5,5,5	0.54	0	5,5,5	0.34	0
3	PEG	G	402	-	6,6,6	0.63	0	5,5,5	0.37	0
2	GOL	E	403	-	5,5,5	0.37	0	5,5,5	0.27	0
2	GOL	I	402	-	5,5,5	0.52	0	5,5,5	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GOL	N	402	-	5,5,5	0.49	0	5,5,5	0.24	0
6	PGE	B	403	-	9,9,9	0.59	0	8,8,8	0.28	0
4	EDO	F	404	-	3,3,3	0.58	0	2,2,2	0.22	0
5	N2P	N	401	-	6,6,6	0.47	0	5,5,5	0.39	0
4	EDO	A	405	-	3,3,3	0.69	0	2,2,2	0.17	0
5	N2P	K	401	-	6,6,6	0.30	0	5,5,5	0.88	0
2	GOL	K	403	-	5,5,5	1.05	0	5,5,5	0.90	0
2	GOL	C	404	-	5,5,5	0.53	0	5,5,5	0.22	0
2	GOL	A	404	-	5,5,5	0.56	0	5,5,5	0.38	0
2	GOL	A	402	-	5,5,5	0.39	0	5,5,5	0.50	0
4	EDO	C	405	-	3,3,3	0.59	0	2,2,2	0.16	0
2	GOL	M	403	-	5,5,5	1.33	0	5,5,5	0.76	0
2	GOL	A	401	-	5,5,5	0.76	0	5,5,5	0.43	0
5	N2P	D	401	-	6,6,6	0.29	0	5,5,5	0.49	0
5	N2P	O	401	-	6,6,6	0.55	0	5,5,5	0.81	0
2	GOL	O	403	-	5,5,5	0.47	0	5,5,5	0.12	0
3	PEG	A	403	-	6,6,6	0.64	0	5,5,5	0.27	0
5	N2P	M	401	-	6,6,6	0.43	0	5,5,5	1.05	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	GOL	M	404	-	-	4/4/4/4	-
4	EDO	K	404	-	-	1/1/1/1	-
5	N2P	F	402	-	-	1/4/4/4	-
2	GOL	B	402	-	-	2/4/4/4	-
2	GOL	F	401	-	-	2/4/4/4	-
3	PEG	L	402	-	-	0/4/4/4	-
2	GOL	K	402	-	-	2/4/4/4	-
2	GOL	O	402	-	-	2/4/4/4	-
5	N2P	L	401	-	-	1/4/4/4	-
2	GOL	F	403	-	-	2/4/4/4	-
2	GOL	D	402	-	-	2/4/4/4	-
2	GOL	M	402	-	-	2/4/4/4	-
4	EDO	I	404	-	-	1/1/1/1	-
2	GOL	I	401	-	-	2/4/4/4	-
2	GOL	G	403	-	-	1/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	I	403	-	-	2/4/4/4	-
2	GOL	C	402	-	-	2/4/4/4	-
4	EDO	N	403	-	-	1/1/1/1	-
2	GOL	C	403	-	-	2/4/4/4	-
5	N2P	G	401	-	-	1/4/4/4	-
4	EDO	L	403	-	-	1/1/1/1	-
4	EDO	D	403	-	-	0/1/1/1	-
2	GOL	H	401	-	-	3/4/4/4	-
5	N2P	J	401	-	-	0/4/4/4	-
5	N2P	C	401	-	-	0/4/4/4	-
4	EDO	J	402	-	-	1/1/1/1	-
5	N2P	B	401	-	-	0/4/4/4	-
5	N2P	E	401	-	-	0/4/4/4	-
2	GOL	E	402	-	-	4/4/4/4	-
3	PEG	G	402	-	-	2/4/4/4	-
2	GOL	E	403	-	-	2/4/4/4	-
2	GOL	I	402	-	-	0/4/4/4	-
2	GOL	N	402	-	-	2/4/4/4	-
6	PGE	B	403	-	-	4/7/7/7	-
4	EDO	F	404	-	-	1/1/1/1	-
5	N2P	N	401	-	-	0/4/4/4	-
4	EDO	A	405	-	-	1/1/1/1	-
5	N2P	K	401	-	-	0/4/4/4	-
2	GOL	K	403	-	-	2/4/4/4	-
2	GOL	C	404	-	-	4/4/4/4	-
2	GOL	A	404	-	-	2/4/4/4	-
2	GOL	A	402	-	-	2/4/4/4	-
4	EDO	C	405	-	-	1/1/1/1	-
2	GOL	M	403	-	-	4/4/4/4	-
2	GOL	A	401	-	-	2/4/4/4	-
5	N2P	D	401	-	-	0/4/4/4	-
5	N2P	O	401	-	-	1/4/4/4	-
2	GOL	O	403	-	-	0/4/4/4	-
3	PEG	A	403	-	-	3/4/4/4	-
5	N2P	M	401	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (75) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	402	GOL	O1-C1-C2-O2
2	A	404	GOL	O2-C2-C3-O3
2	B	402	GOL	C1-C2-C3-O3
2	C	404	GOL	O1-C1-C2-O2
2	C	404	GOL	O1-C1-C2-C3
2	D	402	GOL	O1-C1-C2-C3
2	H	401	GOL	O1-C1-C2-O2
2	K	402	GOL	C1-C2-C3-O3
2	K	403	GOL	O1-C1-C2-C3
2	M	403	GOL	O1-C1-C2-C3
2	M	403	GOL	C1-C2-C3-O3
2	M	403	GOL	O2-C2-C3-O3
2	M	404	GOL	O1-C1-C2-O2
2	M	404	GOL	O1-C1-C2-C3
2	O	402	GOL	O1-C1-C2-O2
2	O	402	GOL	O1-C1-C2-C3
2	D	402	GOL	O1-C1-C2-O2
2	E	402	GOL	O2-C2-C3-O3
2	K	402	GOL	O2-C2-C3-O3
2	N	402	GOL	O2-C2-C3-O3
3	A	403	PEG	O2-C3-C4-O4
3	A	403	PEG	O1-C1-C2-O2
3	G	402	PEG	O2-C3-C4-O4
2	A	402	GOL	O1-C1-C2-C3
2	A	404	GOL	C1-C2-C3-O3
2	C	402	GOL	O1-C1-C2-C3
2	C	403	GOL	O1-C1-C2-C3
2	C	404	GOL	C1-C2-C3-O3
2	E	402	GOL	O1-C1-C2-C3
2	E	402	GOL	C1-C2-C3-O3
2	E	403	GOL	O1-C1-C2-C3
2	F	401	GOL	O1-C1-C2-C3
2	F	403	GOL	O1-C1-C2-C3
2	H	401	GOL	O1-C1-C2-C3
2	I	401	GOL	C1-C2-C3-O3
2	M	402	GOL	C1-C2-C3-O3
2	M	404	GOL	C1-C2-C3-O3
2	N	402	GOL	C1-C2-C3-O3
2	B	402	GOL	O2-C2-C3-O3
2	C	402	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	F	401	GOL	O1-C1-C2-O2
2	F	403	GOL	O1-C1-C2-O2
2	I	401	GOL	O2-C2-C3-O3
2	K	403	GOL	O1-C1-C2-O2
2	M	402	GOL	O2-C2-C3-O3
2	M	403	GOL	O1-C1-C2-O2
4	I	404	EDO	O1-C1-C2-O2
6	B	403	PGE	O2-C3-C4-O3
2	E	402	GOL	O1-C1-C2-O2
4	A	405	EDO	O1-C1-C2-O2
4	K	404	EDO	O1-C1-C2-O2
4	N	403	EDO	O1-C1-C2-O2
2	A	401	GOL	O2-C2-C3-O3
2	I	403	GOL	O2-C2-C3-O3
2	M	404	GOL	O2-C2-C3-O3
5	L	401	N2P	C2-C3-C4-C5
4	C	405	EDO	O1-C1-C2-O2
4	F	404	EDO	O1-C1-C2-O2
4	L	403	EDO	O1-C1-C2-O2
5	O	401	N2P	C2-C3-C4-C5
6	B	403	PGE	C6-C5-O3-C4
6	B	403	PGE	C1-C2-O2-C3
2	C	403	GOL	O1-C1-C2-O2
2	C	404	GOL	O2-C2-C3-O3
2	E	403	GOL	O1-C1-C2-O2
2	I	403	GOL	O1-C1-C2-C3
3	G	402	PEG	O1-C1-C2-O2
5	G	401	N2P	C2-C3-C4-C5
2	A	401	GOL	C1-C2-C3-O3
5	F	402	N2P	C2-C3-C4-C5
3	A	403	PEG	C4-C3-O2-C2
2	G	403	GOL	O2-C2-C3-O3
4	J	402	EDO	O1-C1-C2-O2
6	B	403	PGE	O1-C1-C2-O2
2	H	401	GOL	O2-C2-C3-O3

There are no ring outliers.

14 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	402	N2P	3	0
2	B	402	GOL	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	401	GOL	1	0
5	L	401	N2P	3	0
5	G	401	N2P	2	0
5	J	401	N2P	3	0
5	C	401	N2P	3	0
5	B	401	N2P	3	0
5	E	401	N2P	3	0
5	N	401	N2P	3	0
5	K	401	N2P	3	0
5	D	401	N2P	3	0
5	O	401	N2P	3	0
5	M	401	N2P	2	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	296/304 (97%)	0.38	20 (6%)	23 20	26, 55, 98, 126	3 (1%)
1	B	297/304 (97%)	-0.28	4 (1%)	75 73	29, 41, 61, 93	0
1	C	301/304 (99%)	-0.38	3 (0%)	79 77	25, 40, 57, 109	1 (0%)
1	D	298/304 (98%)	-0.47	1 (0%)	90 88	20, 37, 49, 85	3 (1%)
1	E	297/304 (97%)	-0.37	1 (0%)	90 88	29, 41, 56, 83	0
1	F	298/304 (98%)	-0.02	2 (0%)	84 82	30, 51, 71, 92	0
1	G	295/304 (97%)	1.26	54 (18%)	3 2	52, 76, 101, 114	0
1	H	268/304 (88%)	2.09	131 (48%)	0 0	59, 97, 136, 155	0
1	I	292/304 (96%)	0.22	13 (4%)	38 35	24, 49, 101, 126	1 (0%)
1	J	298/304 (98%)	-0.30	3 (1%)	79 77	27, 41, 59, 113	1 (0%)
1	K	301/304 (99%)	-0.29	3 (0%)	79 77	28, 42, 60, 116	0
1	L	298/304 (98%)	-0.47	1 (0%)	90 88	27, 38, 52, 100	2 (0%)
1	M	298/304 (98%)	-0.42	2 (0%)	84 82	28, 38, 54, 92	0
1	N	298/304 (98%)	-0.31	2 (0%)	84 82	27, 40, 55, 92	1 (0%)
1	O	298/304 (98%)	0.05	5 (1%)	69 66	34, 52, 73, 121	1 (0%)
1	P	279/304 (91%)	0.72	28 (10%)	12 10	33, 69, 116, 140	1 (0%)
All	All	4712/4864 (96%)	0.07	273 (5%)	29 25	20, 46, 98, 155	14 (0%)

All (273) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	193	ILE	6.2
1	A	129	GLY	6.2
1	H	185	GLY	6.1
1	H	234	ILE	6.0
1	H	135	TYR	5.6

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Mol	Chain	Res	Type	RSRZ
1	H	134	PHE	5.6
1	H	49	LEU	5.4
1	H	63	PHE	5.4
1	I	128	PRO	5.3
1	H	58	ALA	5.3
1	H	53	TYR	5.3
1	I	130	TYR	5.3
1	H	22	VAL	5.3
1	H	159	TRP	5.2
1	H	118	ILE	5.1
1	I	301	LEU	5.1
1	H	232	SER	5.0
1	H	19	THR	5.0
1	H	98	ALA	4.8
1	P	300	VAL	4.5
1	H	55	PHE	4.5
1	A	128	PRO	4.5
1	H	70	TYR	4.4
1	G	82	LEU	4.4
1	G	116	LEU	4.3
1	H	64	ILE	4.3
1	G	273	MET	4.2
1	H	28	THR	4.1
1	H	184	ILE	4.1
1	O	301	LEU	4.1
1	G	35	ALA	4.0
1	H	50	PHE	4.0
1	H	89	VAL	4.0
1	O	277	TRP	4.0
1	H	116	LEU	4.0
1	H	76	ILE	4.0
1	A	126	ASP	4.0
1	G	8	LYS	3.9
1	G	9	VAL	3.9
1	H	220	ILE	3.9
1	F	301	LEU	3.9
1	G	86	LEU	3.9
1	G	277	TRP	3.9
1	P	225	ILE	3.9
1	H	24	THR	3.9
1	B	277	TRP	3.8
1	G	107	ALA	3.8

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Mol	Chain	Res	Type	RSRZ
1	G	263	ALA	3.7
1	P	56	CYS	3.6
1	H	236	PHE	3.6
1	A	127	GLY	3.6
1	H	21	ASP	3.6
1	K	2	ALA	3.6
1	H	5	LYS	3.6
1	H	162	TRP	3.5
1	H	45	LEU	3.5
1	K	301	LEU	3.5
1	C	1	MET	3.5
1	I	57	GLN	3.5
1	H	23	SER	3.5
1	H	121	LYS	3.5
1	H	57	GLN	3.5
1	H	43	ILE	3.4
1	H	46	ILE	3.4
1	M	301	LEU	3.4
1	P	58	ALA	3.4
1	E	301	LEU	3.4
1	G	301	LEU	3.4
1	P	231	LYS	3.4
1	A	130	TYR	3.4
1	H	27	THR	3.4
1	I	300	VAL	3.4
1	H	92	VAL	3.3
1	H	60	ARG	3.3
1	G	27	THR	3.3
1	I	127	GLY	3.3
1	H	235	LYS	3.3
1	H	93	SER	3.3
1	G	11	VAL	3.3
1	H	79	LEU	3.3
1	H	69	PRO	3.2
1	H	35	ALA	3.2
1	H	299	PRO	3.2
1	G	22	VAL	3.2
1	H	233	GLU	3.2
1	H	29	ALA	3.2
1	L	301	LEU	3.1
1	G	249	ILE	3.1
1	I	129	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	G	10	VAL	3.1
1	H	44	VAL	3.1
1	H	32	LEU	3.1
1	H	74	PRO	3.0
1	H	88	VAL	3.0
1	H	20	ASP	3.0
1	H	59	GLN	3.0
1	J	277	TRP	3.0
1	H	33	VAL	3.0
1	J	4	ASP	3.0
1	P	133	LYS	3.0
1	H	237	TYR	3.0
1	K	1	MET	3.0
1	G	211	LEU	2.9
1	J	301	LEU	2.9
1	H	125	PRO	2.9
1	G	89	VAL	2.9
1	G	252	ILE	2.9
1	H	106	ILE	2.9
1	H	157	ILE	2.9
1	A	191	GLN	2.9
1	H	94	PHE	2.9
1	P	134	PHE	2.9
1	I	133	LYS	2.9
1	B	301	LEU	2.9
1	H	102	HIS	2.9
1	H	158	CYS	2.9
1	H	26	VAL	2.9
1	H	273	MET	2.9
1	A	133	LYS	2.9
1	H	11	VAL	2.8
1	H	51	GLU	2.8
1	P	54	TYR	2.8
1	A	125	PRO	2.8
1	H	109	ILE	2.8
1	H	12	SER	2.8
1	G	106	ILE	2.8
1	H	103	TYR	2.8
1	A	299	PRO	2.8
1	H	101	ALA	2.7
1	C	301	LEU	2.7
1	H	246	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	P	186	SER	2.7
1	P	5	LYS	2.7
1	H	167	ALA	2.7
1	A	57	GLN	2.7
1	P	132	GLU	2.7
1	P	64	ILE	2.7
1	G	19	THR	2.7
1	N	301	LEU	2.7
1	G	88	VAL	2.7
1	F	4	ASP	2.7
1	H	149	LYS	2.7
1	C	2	ALA	2.6
1	H	243	ALA	2.6
1	G	16	PHE	2.6
1	A	188	PRO	2.6
1	H	119	TYR	2.6
1	H	82	LEU	2.6
1	H	178	LEU	2.6
1	P	4	ASP	2.6
1	H	56	CYS	2.6
1	H	17	ALA	2.6
1	H	75	THR	2.6
1	P	55	PHE	2.6
1	H	83	ALA	2.6
1	H	141	THR	2.6
1	H	136	PHE	2.5
1	H	163	PHE	2.5
1	A	131	GLU	2.5
1	G	250	VAL	2.5
1	G	38	LYS	2.5
1	G	152	LYS	2.5
1	P	57	GLN	2.5
1	G	260	VAL	2.5
1	A	229	HIS	2.5
1	H	62	ASP	2.5
1	H	194	ASP	2.5
1	H	183	ALA	2.5
1	I	232	SER	2.5
1	H	123	HIS	2.5
1	G	118	ILE	2.5
1	H	66	ARG	2.5
1	A	190	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	13	ALA	2.5
1	H	95	PHE	2.5
1	A	300	VAL	2.4
1	H	124	ILE	2.5
1	H	86	LEU	2.4
1	G	146	PHE	2.4
1	G	33	VAL	2.4
1	G	153	ILE	2.4
1	H	9	VAL	2.4
1	H	108	ILE	2.4
1	H	180	TYR	2.4
1	H	36	ALA	2.4
1	H	297	LYS	2.4
1	G	37	HIS	2.4
1	H	117	GLY	2.4
1	P	63	PHE	2.4
1	P	236	PHE	2.4
1	P	194	ASP	2.4
1	H	259	ALA	2.4
1	G	267	LEU	2.4
1	H	16	PHE	2.4
1	H	143	PHE	2.4
1	G	193	ILE	2.4
1	G	224	ILE	2.4
1	G	279	VAL	2.4
1	B	5	LYS	2.4
1	H	67	ALA	2.3
1	I	187	GLU	2.3
1	H	14	LEU	2.3
1	H	65	GLN	2.3
1	G	215	VAL	2.3
1	H	212	VAL	2.3
1	G	259	ALA	2.3
1	G	261	LEU	2.3
1	G	46	ILE	2.3
1	H	252	ILE	2.3
1	G	32	LEU	2.3
1	H	256	LYS	2.3
1	A	232	SER	2.3
1	H	262	ILE	2.3
1	P	100	ASN	2.3
1	H	156	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	H	166	ALA	2.3
1	H	54	TYR	2.3
1	H	181	PRO	2.2
1	H	38	LYS	2.2
1	H	52	GLY	2.2
1	D	301	LEU	2.2
1	I	132	GLU	2.2
1	G	90	ILE	2.2
1	G	177	ILE	2.2
1	H	99	ASN	2.2
1	H	250	VAL	2.2
1	M	4	ASP	2.2
1	A	132	GLU	2.2
1	O	276	CYS	2.2
1	I	229	HIS	2.2
1	P	53	TYR	2.2
1	H	182	THR	2.2
1	H	241	PHE	2.2
1	G	92	VAL	2.2
1	H	244	GLY	2.2
1	G	36	ALA	2.2
1	B	275	HIS	2.2
1	H	155	VAL	2.1
1	P	60	ARG	2.1
1	H	195	SER	2.1
1	O	5	LYS	2.1
1	G	150	TYR	2.1
1	O	273	MET	2.1
1	H	48	GLU	2.1
1	P	135	TYR	2.1
1	H	249	ILE	2.1
1	G	179	PHE	2.1
1	G	145	VAL	2.1
1	H	202	VAL	2.1
1	H	261	LEU	2.1
1	P	273	MET	2.1
1	P	62	ASP	2.1
1	G	7	ARG	2.1
1	P	125	PRO	2.1
1	G	129	GLY	2.1
1	A	5	LYS	2.1
1	G	262	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	H	260	VAL	2.1
1	G	287	LEU	2.0
1	P	99	ASN	2.0
1	H	18	CYS	2.0
1	A	189	HIS	2.0
1	H	153	ILE	2.0
1	I	225	ILE	2.0
1	H	10	VAL	2.0
1	G	29	ALA	2.0
1	H	107	ALA	2.0
1	H	111	ALA	2.0
1	H	263	ALA	2.0
1	N	21	ASP	2.0
1	G	149	LYS	2.0
1	P	24	THR	2.0
1	P	69	PRO	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 oligosaccharides 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, 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	PEG	A	403	7/7	0.66	0.22	81,85,92,92	0
3	PEG	G	402	7/7	0.77	0.18	72,76,84,86	0
2	GOL	H	401	6/6	0.79	0.25	94,97,99,101	0
3	PEG	L	402	7/7	0.79	0.22	75,83,90,90	0
2	GOL	A	404	6/6	0.84	0.14	69,74,76,77	0
4	EDO	A	405	4/4	0.84	0.20	60,60,61,62	0
2	GOL	I	403	6/6	0.85	0.17	69,74,82,84	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	D	403	4/4	0.85	0.16	74,75,78,79	0
6	PGE	B	403	10/10	0.85	0.18	63,78,88,94	0
2	GOL	B	402	6/6	0.86	0.17	62,68,71,78	0
2	GOL	O	403	6/6	0.86	0.15	66,71,75,75	0
4	EDO	C	405	4/4	0.86	0.21	68,68,72,73	0
2	GOL	F	403	6/6	0.86	0.15	67,80,81,83	0
4	EDO	I	404	4/4	0.86	0.19	64,64,65,69	0
4	EDO	L	403	4/4	0.86	0.20	59,61,61,66	0
2	GOL	A	401	6/6	0.86	0.18	80,83,85,86	0
4	EDO	F	404	4/4	0.88	0.21	74,77,77,78	0
2	GOL	C	404	6/6	0.88	0.18	72,78,80,81	0
5	N2P	M	401	7/7	0.89	0.15	39,40,41,43	0
4	EDO	J	402	4/4	0.90	0.14	68,68,69,70	0
2	GOL	M	403	6/6	0.90	0.14	54,60,61,62	0
4	EDO	K	404	4/4	0.91	0.16	65,66,70,74	0
2	GOL	E	403	6/6	0.91	0.11	54,59,61,63	0
4	EDO	N	403	4/4	0.91	0.19	73,74,75,75	0
2	GOL	N	402	6/6	0.91	0.12	54,65,68,68	0
2	GOL	E	402	6/6	0.91	0.17	70,78,83,89	0
2	GOL	M	404	6/6	0.92	0.13	59,65,67,79	0
2	GOL	F	401	6/6	0.92	0.13	56,63,66,66	0
2	GOL	O	402	6/6	0.92	0.11	62,63,63,64	0
2	GOL	I	402	6/6	0.92	0.11	58,64,67,69	0
5	N2P	E	401	7/7	0.92	0.14	42,42,43,44	0
2	GOL	C	403	6/6	0.92	0.10	58,66,71,71	0
5	N2P	N	401	7/7	0.92	0.13	45,45,46,46	0
2	GOL	G	403	6/6	0.92	0.10	60,64,66,67	0
5	N2P	C	401	7/7	0.93	0.14	41,41,43,43	0
2	GOL	M	402	6/6	0.93	0.12	54,59,64,67	0
5	N2P	J	401	7/7	0.93	0.13	45,46,47,48	0
5	N2P	K	401	7/7	0.93	0.13	45,45,47,48	0
2	GOL	A	402	6/6	0.93	0.14	58,62,68,72	0
2	GOL	D	402	6/6	0.93	0.13	55,57,62,65	0
5	N2P	B	401	7/7	0.93	0.11	42,42,43,44	0
2	GOL	K	403	6/6	0.94	0.11	53,58,61,61	0
5	N2P	F	402	7/7	0.94	0.13	61,61,63,63	0
5	N2P	G	401	7/7	0.94	0.14	71,73,74,75	0
5	N2P	O	401	7/7	0.94	0.11	50,50,51,52	0
5	N2P	D	401	7/7	0.94	0.14	42,43,44,44	0
2	GOL	C	402	6/6	0.95	0.10	52,57,64,67	0
2	GOL	I	401	6/6	0.95	0.10	49,53,57,60	0
2	GOL	K	402	6/6	0.95	0.10	59,66,69,72	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	N2P	L	401	7/7	0.95	0.13	42,42,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.