



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:33 PM UTC

PDB ID : 5YQB / pdb_00005yqb
Title : Crystal structure of E.coli aminopeptidase N in complex with Puromycin
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Deposited on : 2017-11-06
Resolution : 1.56 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

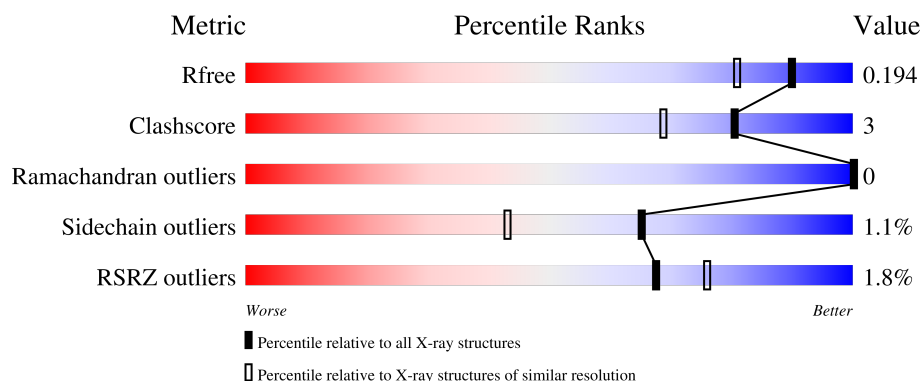
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.56 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	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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	891	2% 83% 13% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7949 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 Aminopeptidase N.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	866	7038	4460	1212	1337	29	0	19	0

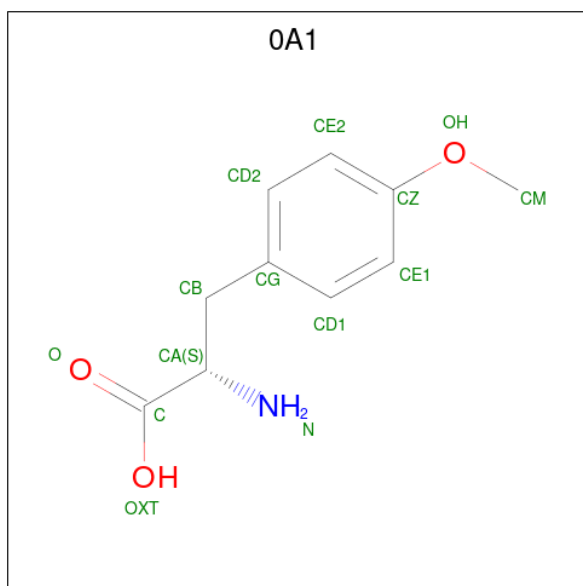
There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP P04825
A	-19	GLY	-	expression tag	UNP P04825
A	-18	SER	-	expression tag	UNP P04825
A	-17	SER	-	expression tag	UNP P04825
A	-16	HIS	-	expression tag	UNP P04825
A	-15	HIS	-	expression tag	UNP P04825
A	-14	HIS	-	expression tag	UNP P04825
A	-13	HIS	-	expression tag	UNP P04825
A	-12	HIS	-	expression tag	UNP P04825
A	-11	HIS	-	expression tag	UNP P04825
A	-10	SER	-	expression tag	UNP P04825
A	-9	SER	-	expression tag	UNP P04825
A	-8	GLY	-	expression tag	UNP P04825
A	-7	LEU	-	expression tag	UNP P04825
A	-6	VAL	-	expression tag	UNP P04825
A	-5	PRO	-	expression tag	UNP P04825
A	-4	ARG	-	expression tag	UNP P04825
A	-3	GLY	-	expression tag	UNP P04825
A	-2	SER	-	expression tag	UNP P04825
A	-1	HIS	-	expression tag	UNP P04825
A	0	MET	-	expression tag	UNP P04825

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

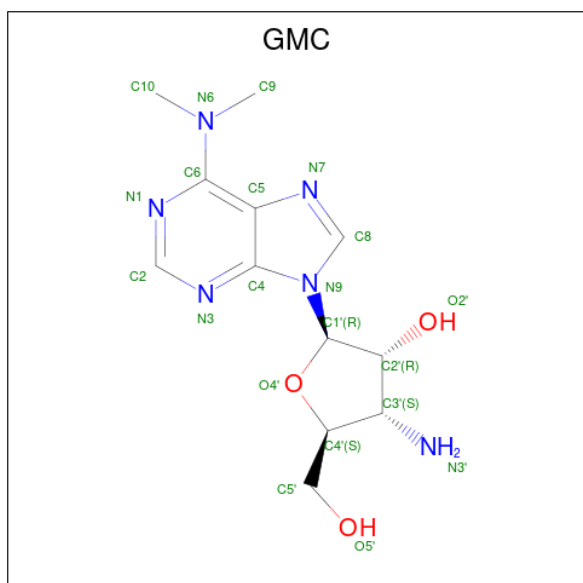
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is O-methyl-L-tyrosine (CCD ID: 0A1) (formula: $C_{10}H_{13}NO_3$).



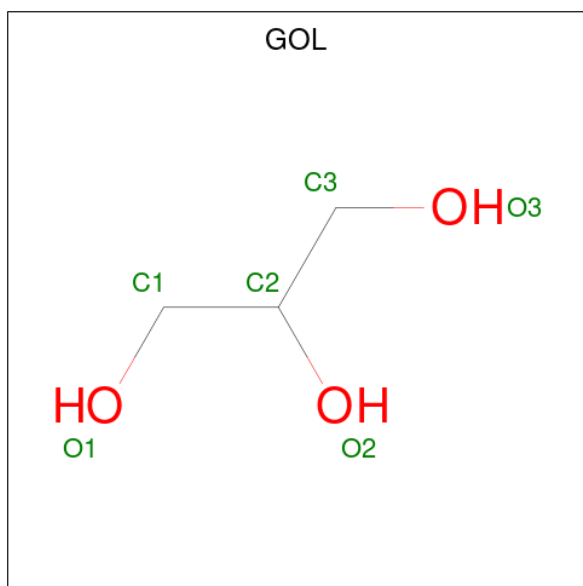
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	10	1	3		

- Molecule 4 is (2R,3R,4S,5S)-4-AMINO-2-[6-(DIMETHYLAMINO)-9H-PURIN-9-YL]-5-(HYDROXYMETHYL)TETRAHYDRO-3-FURANOL (CCD ID: GMC) (formula: $C_{12}H_{18}N_6O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			21	12	6	3		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

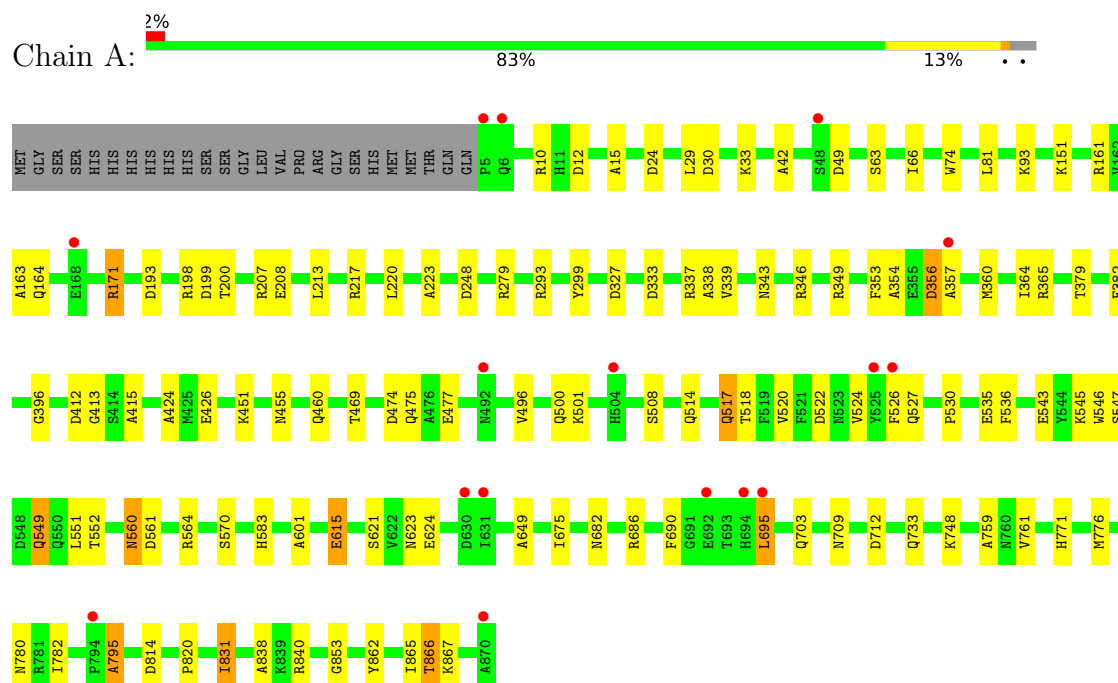
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	869	Total	O	0	0
			869	869		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Aminopeptidase N



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.29Å 120.29Å 170.12Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 1.56 50.00 – 1.56	Depositor EDS
% Data completeness (in resolution range)	98.2 (50.00-1.56) 98.2 (50.00-1.56)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.76 (at 1.56Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.161 , 0.182 0.175 , 0.194	Depositor DCC
R_{free} test set	9844 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	22.4	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.006 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7949	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GMC, ZN, GOL, 0A1

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	1.75	69/7247 (1.0%)	1.55	57/9835 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (69) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	866	THR	CB-OG1	-9.93	1.27	1.43
1	A	207	ARG	CD-NE	9.08	1.58	1.46
1	A	357	ALA	N-CA	8.95	1.56	1.46
1	A	349	ARG	CD-NE	-8.63	1.34	1.46
1	A	518	THR	C-O	8.51	1.34	1.23
1	A	564	ARG	CD-NE	-8.42	1.34	1.46
1	A	354	ALA	N-CA	7.66	1.55	1.46
1	A	530	PRO	CA-CB	-7.60	1.45	1.54
1	A	501	LYS	N-CA	-7.56	1.37	1.46
1	A	865	ILE	N-CA	7.52	1.55	1.46
1	A	460	GLN	N-CA	7.50	1.55	1.46
1	A	451	LYS	N-CA	7.09	1.54	1.45
1	A	560	ASN	CG-OD1	6.97	1.36	1.23
1	A	29	LEU	CB-CG	-6.74	1.40	1.53
1	A	354	ALA	CA-CB	-6.70	1.42	1.53
1	A	415	ALA	CA-CB	-6.69	1.44	1.53
1	A	338	ALA	CA-C	-6.61	1.44	1.52
1	A	551	LEU	CA-C	-6.48	1.44	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	761	VAL	N-CA	6.30	1.54	1.46
1	A	601	ALA	CA-CB	-6.28	1.43	1.53
1	A	552	THR	CA-CB	6.05	1.63	1.53
1	A	831	ILE	CB-CG1	-6.01	1.41	1.53
1	A	349	ARG	NE-CZ	6.00	1.39	1.33
1	A	649	ALA	CA-CB	-6.00	1.44	1.53
1	A	615	GLU	CD-OE2	5.92	1.36	1.25
1	A	198	ARG	NE-CZ	-5.90	1.26	1.33
1	A	74	TRP	C-O	5.87	1.30	1.23
1	A	63	SER	CA-C	-5.82	1.45	1.52
1	A	382	GLU	CG-CD	-5.78	1.37	1.52
1	A	15	ALA	CA-CB	-5.73	1.44	1.53
1	A	524	VAL	CA-C	5.64	1.59	1.52
1	A	771	HIS	C-O	5.63	1.30	1.23
1	A	349	ARG	CZ-NH2	-5.63	1.26	1.33
1	A	690	PHE	N-CA	5.58	1.52	1.46
1	A	66	ILE	C-O	5.53	1.30	1.24
1	A	733	GLN	N-CA	-5.51	1.39	1.46
1	A	759	ALA	CA-C	-5.50	1.45	1.52
1	A	820	PRO	CA-CB	-5.50	1.46	1.53
1	A	349	ARG	CG-CD	5.49	1.69	1.52
1	A	530	PRO	C-O	-5.49	1.17	1.23
1	A	477	GLU	N-CA	-5.49	1.39	1.46
1	A	364	ILE	CA-CB	-5.44	1.46	1.54
1	A	200	THR	CA-C	-5.43	1.46	1.52
1	A	42	ALA	CA-C	-5.42	1.46	1.52
1	A	520	VAL	C-O	5.39	1.30	1.24
1	A	337	ARG	CA-CB	-5.38	1.45	1.53
1	A	10	ARG	CZ-NH2	5.36	1.40	1.33
1	A	49	ASP	C-O	-5.35	1.16	1.24
1	A	500	GLN	CA-C	-5.34	1.46	1.52
1	A	840	ARG	CA-C	-5.34	1.45	1.52
1	A	853	GLY	C-O	5.33	1.29	1.24
1	A	12	ASP	C-O	5.32	1.31	1.24
1	A	469	THR	C-N	-5.28	1.26	1.33
1	A	12	ASP	N-CA	5.20	1.52	1.46
1	A	475	GLN	CA-C	-5.19	1.47	1.52
1	A	508	SER	CA-CB	-5.19	1.44	1.53
1	A	547	SER	CA-C	-5.17	1.46	1.53
1	A	333	ASP	C-O	5.17	1.30	1.24
1	A	12	ASP	CA-CB	-5.16	1.44	1.53
1	A	424	ALA	N-CA	-5.16	1.40	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	379	THR	N-CA	-5.13	1.40	1.46
1	A	193	ASP	CG-OD1	-5.12	1.15	1.25
1	A	543	GLU	CD-OE1	5.11	1.35	1.25
1	A	546	TRP	CA-CB	-5.09	1.45	1.53
1	A	163	ALA	CA-CB	-5.08	1.45	1.53
1	A	93	LYS	C-O	5.07	1.29	1.24
1	A	526	PHE	CA-C	-5.07	1.46	1.52
1	A	213	LEU	CA-C	-5.04	1.46	1.52
1	A	171	ARG	C-O	5.02	1.29	1.23

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	866	THR	OG1-CB-CG2	-10.35	88.60	109.30
1	A	686	ARG	NE-CZ-NH2	10.04	128.24	119.20
1	A	349	ARG	NE-CZ-NH2	-9.05	111.05	119.20
1	A	686	ARG	NE-CZ-NH1	-8.42	113.08	121.50
1	A	675	ILE	O-C-N	7.99	129.62	121.87
1	A	795	ALA	N-CA-C	7.98	119.98	111.28
1	A	349	ARG	NH1-CZ-NH2	7.46	129.00	119.30
1	A	356	ASP	O-C-N	7.45	132.41	122.43
1	A	496	VAL	CA-C-N	7.28	127.83	122.59
1	A	496	VAL	C-N-CA	7.28	127.83	122.59
1	A	365	ARG	CB-CA-C	7.06	116.79	110.44
1	A	838	ALA	N-CA-C	7.03	118.94	111.28
1	A	469	THR	O-C-N	6.84	127.48	121.39
1	A	624	GLU	OE1-CD-OE2	6.74	139.07	122.90
1	A	615	GLU	CB-CG-CD	6.67	123.94	112.60
1	A	564	ARG	NE-CZ-NH2	-6.42	113.43	119.20
1	A	413	GLY	O-C-N	-6.38	114.44	122.41
1	A	675	ILE	CA-C-O	-6.31	114.38	120.95
1	A	570	SER	N-CA-C	-6.31	104.40	111.28
1	A	621	SER	CA-C-O	-6.30	114.81	121.99
1	A	396	GLY	CA-C-O	-6.24	116.58	122.39
1	A	248	ASP	N-CA-CB	-6.18	101.45	110.53
1	A	527	GLN	CA-C-N	-6.10	114.01	120.66
1	A	527	GLN	C-N-CA	-6.10	114.01	120.66
1	A	353	PHE	O-C-N	6.05	128.53	122.12
1	A	161	ARG	CA-C-N	5.94	128.92	121.85
1	A	161	ARG	C-N-CA	5.94	128.92	121.85
1	A	623	ASN	O-C-N	-5.83	115.79	122.09
1	A	474	ASP	N-CA-C	5.82	120.10	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	333	ASP	CA-C-O	-5.76	114.31	120.42
1	A	426	GLU	CB-CG-CD	-5.76	102.80	112.60
1	A	357	ALA	N-CA-C	-5.76	106.16	112.72
1	A	339	VAL	N-CA-CB	5.68	117.20	110.55
1	A	560	ASN	CB-CG-OD1	5.65	132.09	120.80
1	A	475	GLN	CB-CA-C	-5.63	102.96	110.79
1	A	522	ASP	CA-C-O	5.54	128.93	121.89
1	A	624	GLU	O-C-N	5.39	127.84	122.12
1	A	455	ASN	CA-C-O	5.39	124.55	119.59
1	A	10	ARG	NE-CZ-NH1	5.38	126.88	121.50
1	A	81	LEU	O-C-N	-5.35	117.01	123.27
1	A	199	ASP	CA-CB-CG	5.33	117.93	112.60
1	A	223	ALA	N-CA-C	5.30	121.53	109.81
1	A	299	TYR	O-C-N	-5.30	116.50	122.12
1	A	561	ASP	N-CA-CB	5.26	117.85	110.12
1	A	207	ARG	O-C-N	-5.25	116.35	122.96
1	A	63	SER	N-CA-C	5.24	116.81	108.79
1	A	522	ASP	O-C-N	-5.24	116.76	122.84
1	A	865	ILE	CA-C-N	-5.21	112.38	120.31
1	A	865	ILE	C-N-CA	-5.21	112.38	120.31
1	A	560	ASN	CB-CG-ND2	-5.20	108.61	116.40
1	A	583	HIS	O-C-N	5.18	127.62	122.12
1	A	549	GLN	CA-C-O	-5.16	114.03	119.97
1	A	217	ARG	O-C-N	5.14	129.23	123.06
1	A	549	GLN	CB-CG-CD	5.13	121.32	112.60
1	A	536	PHE	CA-CB-CG	5.06	118.86	113.80
1	A	353	PHE	CA-C-O	-5.02	115.23	120.55
1	A	279	ARG	CD-NE-CZ	5.02	131.43	124.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	412	ASP	Sidechain

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	7038	0	6911	33	1
2	A	1	0	0	0	0
3	A	14	0	12	1	0
4	A	21	0	18	3	0
5	A	6	0	8	0	0
6	A	869	0	0	14	1
All	All	7949	0	6949	37	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (37) 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:346[B]:ARG:NH1	1:A:615:GLU:OE1	1.71	1.23
1:A:24[B]:ASP:CG	6:A:1001:HOH:O	2.29	0.74
1:A:514:GLN:H	1:A:517[A]:GLN:HE22	1.34	0.73
1:A:862:TYR:O	1:A:866:THR:HG23	1.89	0.71
1:A:24[B]:ASP:OD2	6:A:1001:HOH:O	2.07	0.70
1:A:346[B]:ARG:CZ	1:A:615:GLU:OE1	2.43	0.67
4:A:903:GMC:N7	4:A:903:GMC:H101	2.09	0.66
1:A:517[B]:GLN:HG2	6:A:1542:HOH:O	1.97	0.64
4:A:903:GMC:H2'	4:A:903:GMC:N3	2.14	0.61
1:A:360[B]:MET:HG2	1:A:831:ILE:CD1	2.32	0.60
1:A:171:ARG:NH1	6:A:1003:HOH:O	2.28	0.59
1:A:24[B]:ASP:OD1	6:A:1001:HOH:O	2.17	0.58
1:A:709:ASN:ND2	1:A:712:ASP:H	2.03	0.56
1:A:535[B]:GLU:OE2	1:A:560:ASN:CG	2.50	0.55
4:A:903:GMC:N3	4:A:903:GMC:C2'	2.69	0.55
1:A:360[B]:MET:HE1	1:A:867:LYS:HD2	1.90	0.54
1:A:360[B]:MET:HG2	1:A:831:ILE:HD11	1.91	0.52
3:A:902:OA1:HE1	6:A:1223:HOH:O	2.08	0.52
1:A:780:ASN:ND2	6:A:1014:HOH:O	2.45	0.49
1:A:709:ASN:HD21	1:A:712:ASP:H	1.58	0.49
1:A:360[B]:MET:CG	1:A:831:ILE:CD1	2.92	0.47
1:A:695:LEU:CD2	6:A:1868:HOH:O	2.62	0.47
1:A:360[B]:MET:CE	1:A:867:LYS:HD2	2.44	0.46
1:A:549:GLN:HG2	6:A:1594:HOH:O	2.15	0.46
1:A:695:LEU:HD21	6:A:1859:HOH:O	2.15	0.45
1:A:535[B]:GLU:OE1	6:A:1002:HOH:O	2.21	0.44
1:A:795:ALA:HB3	6:A:1061:HOH:O	2.17	0.44
1:A:682:ASN:HD21	1:A:703:GLN:HE22	1.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:ASP:HB3	1:A:33:LYS:O	2.19	0.43
1:A:695:LEU:HD22	6:A:1868:HOH:O	2.17	0.43
1:A:360[B]:MET:HG2	1:A:831:ILE:HD12	1.99	0.42
1:A:535[B]:GLU:OE2	1:A:560:ASN:ND2	2.52	0.42
1:A:682:ASN:ND2	1:A:703:GLN:HE22	2.17	0.42
1:A:814:ASP:C	1:A:814:ASP:OD1	2.63	0.41
1:A:748:LYS:HG3	6:A:1019:HOH:O	2.21	0.40
1:A:776:MET:O	1:A:782:ILE:HD11	2.22	0.40
1:A:343:ASN:OD1	1:A:346[B]:ARG:NH2	2.55	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:1297:HOH:O	6:A:1297:HOH:O[6_554]	1.07	1.13
1:A:208:GLU:OE1	1:A:545[B]:LYS:NZ[5_544]	2.16	0.04

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	883/891 (99%)	870 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	760/763 (100%)	751 (99%)	9 (1%)	63	40

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	151	LYS
1	A	164	GLN
1	A	220	LEU
1	A	293	ARG
1	A	327	ASP
1	A	356	ASP
1	A	517[A]	GLN
1	A	517[B]	GLN
1	A	695	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	164	GLN
1	A	340	ASN
1	A	584	GLN
1	A	682	ASN
1	A	709	ASN
1	A	724	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	0A1	A	902	2	13,14,14	1.12	0	14,18,18	1.65	3 (21%)
4	GMC	A	903	-	23,23,23	2.25	3 (13%)	31,34,34	3.83	15 (48%)
5	GOL	A	904	-	5,5,5	0.77	0	5,5,5	1.53	1 (20%)

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	0A1	A	902	2	-	0/10/10/10	0/1/1/1
4	GMC	A	903	-	-	0/10/26/26	0/3/3/3
5	GOL	A	904	-	-	2/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	903	GMC	C2'-C3'	-6.84	1.44	1.53
4	A	903	GMC	O4'-C4'	-6.24	1.31	1.45
4	A	903	GMC	C3'-N3'	-2.87	1.43	1.47

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	903	GMC	C5'-C4'-C3'	-11.68	98.31	115.42
4	A	903	GMC	O4'-C4'-C3'	10.62	119.67	104.22
4	A	903	GMC	O4'-C4'-C5'	6.96	123.94	109.22
4	A	903	GMC	C4'-O4'-C1'	-4.95	98.55	109.47
4	A	903	GMC	O4'-C1'-C2'	4.60	116.45	106.62
4	A	903	GMC	N1-C2-N3	-4.23	122.18	128.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	903	GMC	C2-N1-C6	3.55	120.51	111.83
3	A	902	0A1	CG-CB-CA	3.50	121.28	114.13
4	A	903	GMC	C2'-C1'-N9	-3.32	105.06	113.30
4	A	903	GMC	O2'-C2'-C1'	3.04	120.57	110.10
4	A	903	GMC	C9-N6-C6	-2.97	112.96	120.52
4	A	903	GMC	C4-C5-C6	-2.84	112.97	115.91
3	A	902	0A1	CM-OH-CZ	-2.64	111.84	117.50
4	A	903	GMC	C4-N9-C1'	-2.61	120.53	126.63
4	A	903	GMC	C5-C6-N6	-2.50	121.38	125.33
5	A	904	GOL	O2-C2-C3	2.33	118.82	109.18
4	A	903	GMC	O5'-C5'-C4'	-2.30	103.50	111.33
3	A	902	0A1	CE2-CZ-CE1	2.23	123.41	120.16
4	A	903	GMC	C1'-N9-C8	2.07	131.68	127.09

There are no chirality outliers.

All (2) torsion outliers are listed below:

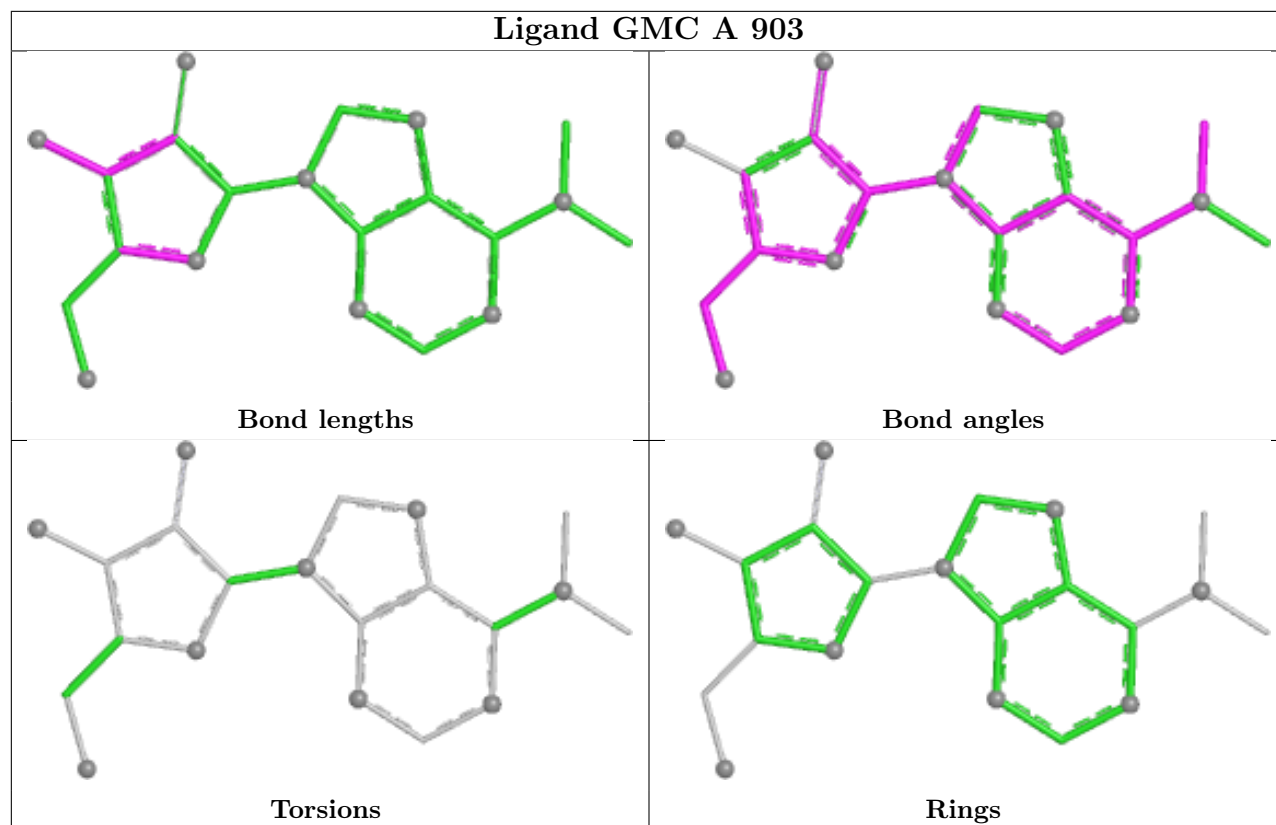
Mol	Chain	Res	Type	Atoms
5	A	904	GOL	O1-C1-C2-C3
5	A	904	GOL	O2-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	902	0A1	1	0
4	A	903	GMC	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	866/891 (97%)	0.01	16 (1%) 67 76	12, 21, 34, 57	20 (2%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	PRO	5.5
1	A	526	PHE	4.8
1	A	694	HIS	3.3
1	A	504	HIS	3.0
1	A	48	SER	2.9
1	A	695	LEU	2.8
1	A	870	ALA	2.6
1	A	794	PRO	2.6
1	A	492	ASN	2.5
1	A	168	GLU	2.5
1	A	525	TYR	2.4
1	A	692	GLU	2.3
1	A	631	ILE	2.2
1	A	630	ASP	2.2
1	A	357	ALA	2.1
1	A	6	GLN	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](#)

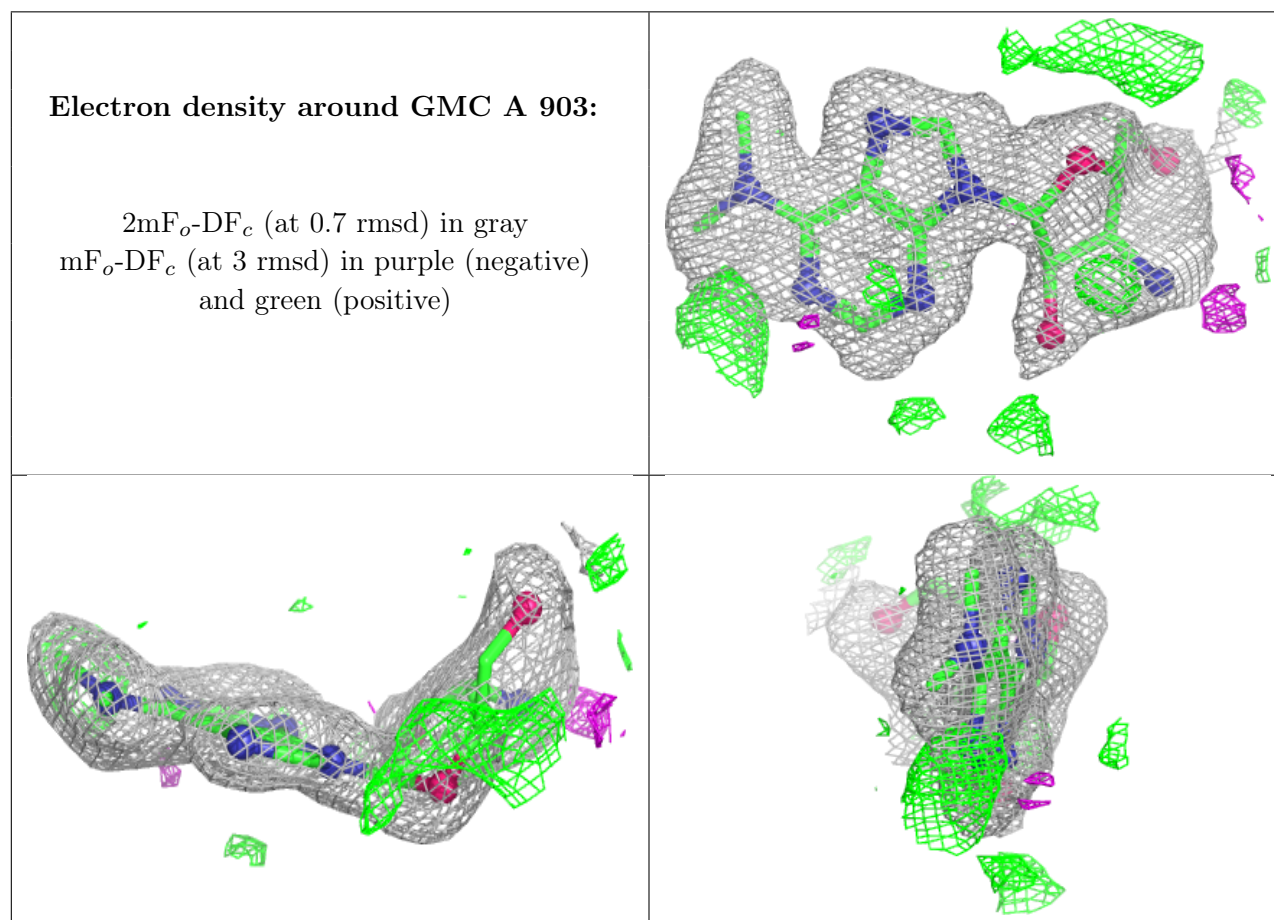
There are no oligosaccharides in this entry.

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
5	GOL	A	904	6/6	0.67	0.23	46,52,57,60	0
4	GMC	A	903	21/21	0.87	0.16	25,32,47,57	21
3	0A1	A	902	14/14	0.96	0.07	16,20,25,25	0
2	ZN	A	901	1/1	1.00	0.02	15,15,15,15	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.