



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:51 PM UTC

PDB ID : 1YUM / pdb_00001yum
Title : Crystal Structure of Nicotinic Acid Mononucleotide Adenylyltransferase from *Pseudomonas aeruginosa*
Authors : Yoon, H.J.; Kim, H.L.; Mikami, B.; Suh, S.W.
Deposited on : 2005-02-14
Resolution : 1.70 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

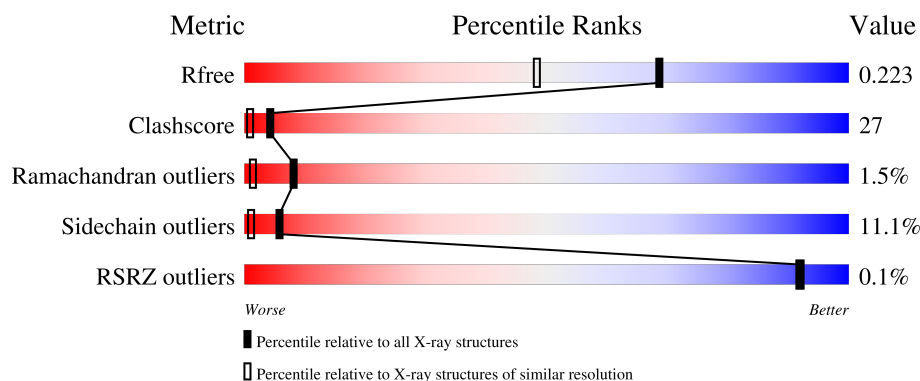
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	242	
1	B	242	
1	C	242	
1	D	242	

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	CIT	B	1301	-	X	-	-
2	CIT	D	3301	-	X	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7740 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 'Probable nicotinate-nucleotide adenylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	6	0
			1694	1073	313	302	6			
1	B	212	Total	C	N	O	S	0	5	0
			1682	1067	307	301	7			
1	C	212	Total	C	N	O	S	0	7	0
			1701	1078	315	302	6			
1	D	212	Total	C	N	O	S	0	8	0
			1710	1087	315	302	6			

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q9HX21
A	-18	GLY	-	expression tag	UNP Q9HX21
A	-17	SER	-	expression tag	UNP Q9HX21
A	-16	SER	-	expression tag	UNP Q9HX21
A	-15	HIS	-	expression tag	UNP Q9HX21
A	-14	HIS	-	expression tag	UNP Q9HX21
A	-13	HIS	-	expression tag	UNP Q9HX21
A	-12	HIS	-	expression tag	UNP Q9HX21
A	-11	HIS	-	expression tag	UNP Q9HX21
A	-10	HIS	-	expression tag	UNP Q9HX21
A	-9	SER	-	expression tag	UNP Q9HX21
A	-8	SER	-	expression tag	UNP Q9HX21
A	-7	GLY	-	expression tag	UNP Q9HX21
A	-6	LEU	-	expression tag	UNP Q9HX21
A	-5	VAL	-	expression tag	UNP Q9HX21
A	-4	PRO	-	expression tag	UNP Q9HX21
A	-3	ARG	-	expression tag	UNP Q9HX21
A	-2	GLY	-	expression tag	UNP Q9HX21
A	-1	SER	-	expression tag	UNP Q9HX21
A	0	HIS	-	expression tag	UNP Q9HX21
A	215	LEU	-	expression tag	UNP Q9HX21

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Chain	Residue	Modelled	Actual	Comment	Reference
A	216	GLU	-	expression tag	UNP Q9HX21
A	217	HIS	-	expression tag	UNP Q9HX21
A	218	HIS	-	expression tag	UNP Q9HX21
A	219	HIS	-	expression tag	UNP Q9HX21
A	220	HIS	-	expression tag	UNP Q9HX21
A	221	HIS	-	expression tag	UNP Q9HX21
A	222	HIS	-	expression tag	UNP Q9HX21
B	-19	MET	-	expression tag	UNP Q9HX21
B	-18	GLY	-	expression tag	UNP Q9HX21
B	-17	SER	-	expression tag	UNP Q9HX21
B	-16	SER	-	expression tag	UNP Q9HX21
B	-15	HIS	-	expression tag	UNP Q9HX21
B	-14	HIS	-	expression tag	UNP Q9HX21
B	-13	HIS	-	expression tag	UNP Q9HX21
B	-12	HIS	-	expression tag	UNP Q9HX21
B	-11	HIS	-	expression tag	UNP Q9HX21
B	-10	HIS	-	expression tag	UNP Q9HX21
B	-9	SER	-	expression tag	UNP Q9HX21
B	-8	SER	-	expression tag	UNP Q9HX21
B	-7	GLY	-	expression tag	UNP Q9HX21
B	-6	LEU	-	expression tag	UNP Q9HX21
B	-5	VAL	-	expression tag	UNP Q9HX21
B	-4	PRO	-	expression tag	UNP Q9HX21
B	-3	ARG	-	expression tag	UNP Q9HX21
B	-2	GLY	-	expression tag	UNP Q9HX21
B	-1	SER	-	expression tag	UNP Q9HX21
B	0	HIS	-	expression tag	UNP Q9HX21
B	215	LEU	-	expression tag	UNP Q9HX21
B	216	GLU	-	expression tag	UNP Q9HX21
B	217	HIS	-	expression tag	UNP Q9HX21
B	218	HIS	-	expression tag	UNP Q9HX21
B	219	HIS	-	expression tag	UNP Q9HX21
B	220	HIS	-	expression tag	UNP Q9HX21
B	221	HIS	-	expression tag	UNP Q9HX21
B	222	HIS	-	expression tag	UNP Q9HX21
C	-19	MET	-	expression tag	UNP Q9HX21
C	-18	GLY	-	expression tag	UNP Q9HX21
C	-17	SER	-	expression tag	UNP Q9HX21
C	-16	SER	-	expression tag	UNP Q9HX21
C	-15	HIS	-	expression tag	UNP Q9HX21
C	-14	HIS	-	expression tag	UNP Q9HX21
C	-13	HIS	-	expression tag	UNP Q9HX21

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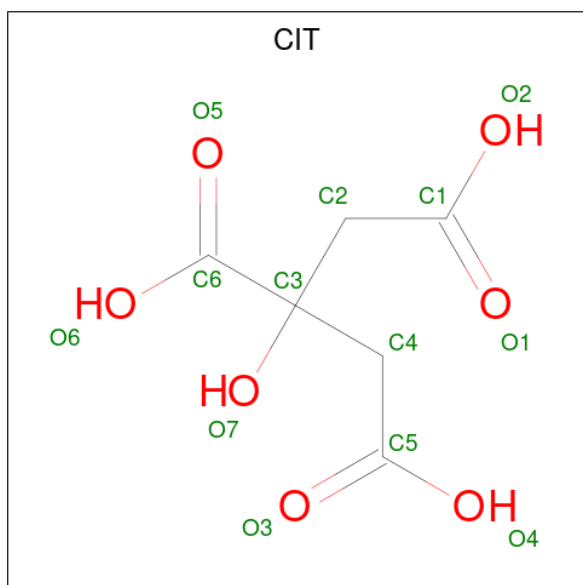
Chain	Residue	Modelled	Actual	Comment	Reference
C	-12	HIS	-	expression tag	UNP Q9HX21
C	-11	HIS	-	expression tag	UNP Q9HX21
C	-10	HIS	-	expression tag	UNP Q9HX21
C	-9	SER	-	expression tag	UNP Q9HX21
C	-8	SER	-	expression tag	UNP Q9HX21
C	-7	GLY	-	expression tag	UNP Q9HX21
C	-6	LEU	-	expression tag	UNP Q9HX21
C	-5	VAL	-	expression tag	UNP Q9HX21
C	-4	PRO	-	expression tag	UNP Q9HX21
C	-3	ARG	-	expression tag	UNP Q9HX21
C	-2	GLY	-	expression tag	UNP Q9HX21
C	-1	SER	-	expression tag	UNP Q9HX21
C	0	HIS	-	expression tag	UNP Q9HX21
C	215	LEU	-	expression tag	UNP Q9HX21
C	216	GLU	-	expression tag	UNP Q9HX21
C	217	HIS	-	expression tag	UNP Q9HX21
C	218	HIS	-	expression tag	UNP Q9HX21
C	219	HIS	-	expression tag	UNP Q9HX21
C	220	HIS	-	expression tag	UNP Q9HX21
C	221	HIS	-	expression tag	UNP Q9HX21
C	222	HIS	-	expression tag	UNP Q9HX21
D	-19	MET	-	expression tag	UNP Q9HX21
D	-18	GLY	-	expression tag	UNP Q9HX21
D	-17	SER	-	expression tag	UNP Q9HX21
D	-16	SER	-	expression tag	UNP Q9HX21
D	-15	HIS	-	expression tag	UNP Q9HX21
D	-14	HIS	-	expression tag	UNP Q9HX21
D	-13	HIS	-	expression tag	UNP Q9HX21
D	-12	HIS	-	expression tag	UNP Q9HX21
D	-11	HIS	-	expression tag	UNP Q9HX21
D	-10	HIS	-	expression tag	UNP Q9HX21
D	-9	SER	-	expression tag	UNP Q9HX21
D	-8	SER	-	expression tag	UNP Q9HX21
D	-7	GLY	-	expression tag	UNP Q9HX21
D	-6	LEU	-	expression tag	UNP Q9HX21
D	-5	VAL	-	expression tag	UNP Q9HX21
D	-4	PRO	-	expression tag	UNP Q9HX21
D	-3	ARG	-	expression tag	UNP Q9HX21
D	-2	GLY	-	expression tag	UNP Q9HX21
D	-1	SER	-	expression tag	UNP Q9HX21
D	0	HIS	-	expression tag	UNP Q9HX21
D	215	LEU	-	expression tag	UNP Q9HX21

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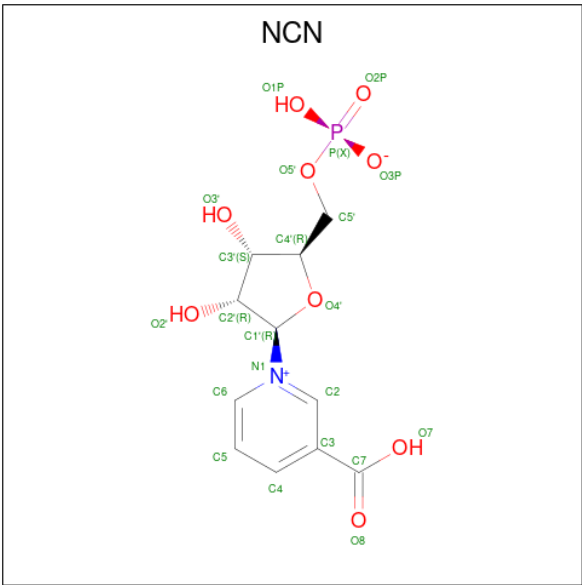
Chain	Residue	Modelled	Actual	Comment	Reference
D	216	GLU	-	expression tag	UNP Q9HX21
D	217	HIS	-	expression tag	UNP Q9HX21
D	218	HIS	-	expression tag	UNP Q9HX21
D	219	HIS	-	expression tag	UNP Q9HX21
D	220	HIS	-	expression tag	UNP Q9HX21
D	221	HIS	-	expression tag	UNP Q9HX21
D	222	HIS	-	expression tag	UNP Q9HX21

- Molecule 2 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		
2	C	1	Total	C	O	0	0
			13	6	7		
2	D	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is NICOTINATE MONONUCLEOTIDE (CCD ID: NCN) (formula: $C_{11}H_{14}NO_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			22	11	1	9	1		
3	B	1	Total	C	N	O	P	0	0
			22	11	1	9	1		
3	C	1	Total	C	N	O	P	0	0
			22	11	1	9	1		
3	D	1	Total	C	N	O	P	0	0
			22	11	1	9	1		

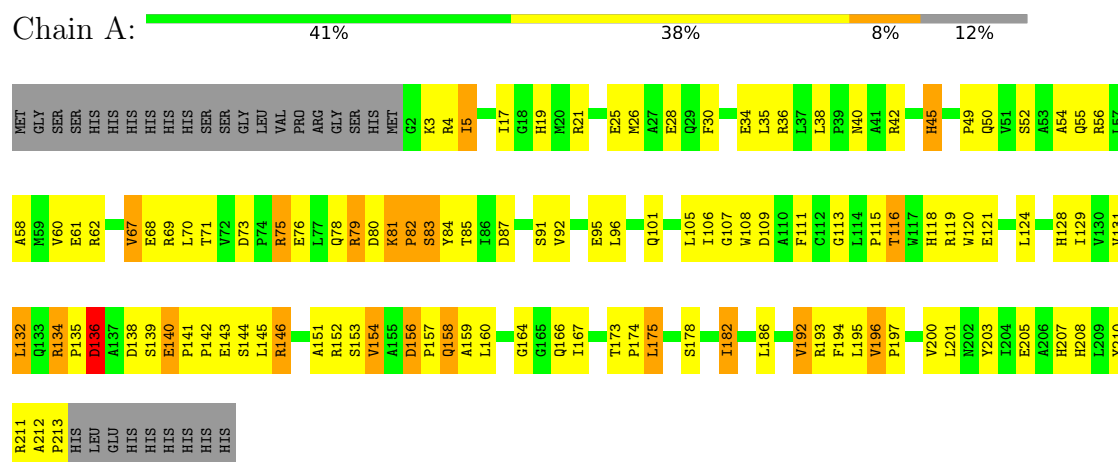
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	164	Total	O	0	0
			164	164		
4	B	215	Total	O	0	0
			215	215		
4	C	225	Total	O	0	0
			225	225		
4	D	209	Total	O	0	0
			209	209		

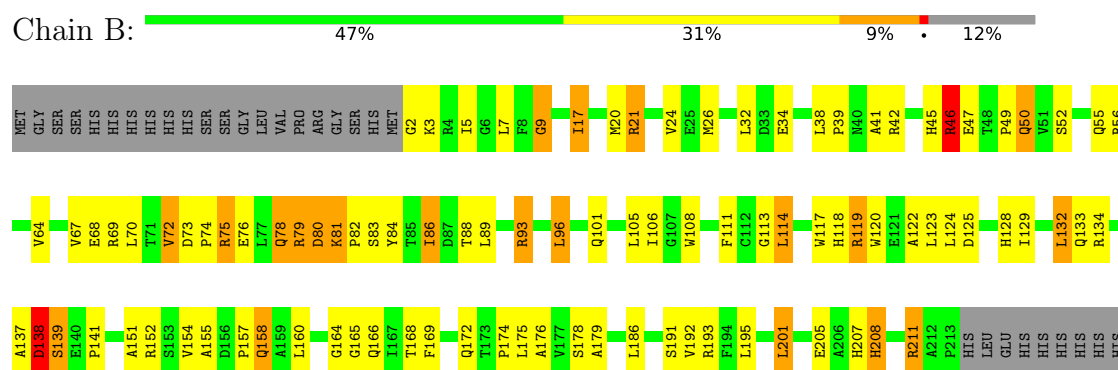
3 Residue-property plots

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.

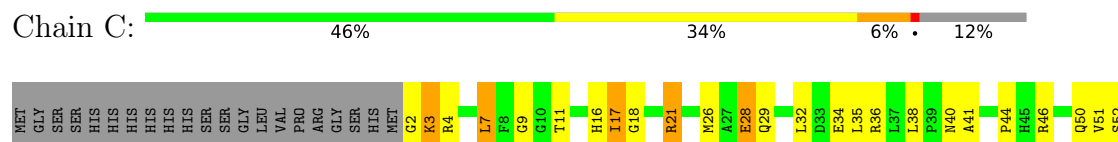
- Molecule 1: 'Probable nicotinate-nucleotide adenylyltransferase

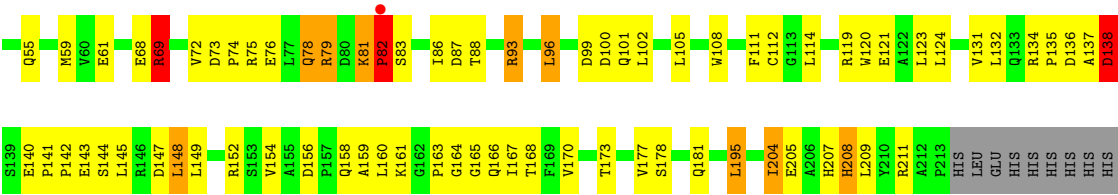


- Molecule 1: 'Probable nicotinate-nucleotide adenylyltransferase

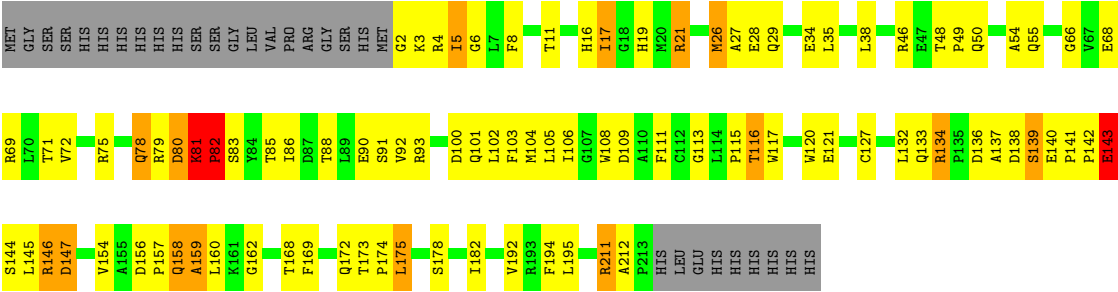


- Molecule 1: 'Probable nicotinate-nucleotide adenylyltransferase





● Molecule 1: 'Probable nicotinate-nucleotide adenylyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.18Å 110.66Å 65.20Å 90.00° 89.99° 90.00°	Depositor
Resolution (Å)	10.00 – 1.70 10.00 – 1.70	Depositor EDS
% Data completeness (in resolution range)	91.3 (10.00-1.70) 94.1 (10.00-1.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.83 (at 1.60Å)	Xtriage
Refinement program	SHELXL-97, SHELXL	Depositor
R, R_{free}	0.191 , 0.251 0.171 , 0.223	Depositor DCC
R_{free} test set	5982 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	18.5	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.487 for -l,k,h 0.488 for h,-k,-l 0.487 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7740	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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: NCN, CIT

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	0.52	0/1763	1.59	23/2402 (1.0%)
1	B	0.53	0/1743	1.62	28/2374 (1.2%)
1	C	0.52	0/1774	1.53	13/2415 (0.5%)
1	D	0.52	0/1791	1.57	20/2438 (0.8%)
All	All	0.52	0/7071	1.58	84/9629 (0.9%)

There are no bond length outliers.

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	30	PHE	CA-CB-CG	-8.20	105.60	113.80
1	B	49	PRO	CA-C-N	8.20	134.28	122.40
1	B	49	PRO	C-N-CA	8.20	134.28	122.40
1	B	164	GLY	CA-C-O	-7.71	114.31	121.63
1	C	46	ARG	CD-NE-CZ	7.19	134.46	124.40
1	A	131	VAL	CA-C-N	7.02	132.46	121.99
1	A	131	VAL	C-N-CA	7.02	132.46	121.99
1	D	143[A]	GLU	CA-C-N	6.99	131.89	120.63
1	D	143[A]	GLU	C-N-CA	6.99	131.89	120.63
1	D	143[B]	GLU	CA-C-N	6.99	131.89	120.63
1	D	143[B]	GLU	C-N-CA	6.99	131.89	120.63
1	B	207	HIS	CA-C-N	6.98	132.28	122.36
1	B	207	HIS	C-N-CA	6.98	132.28	122.36
1	D	21	ARG	NE-CZ-NH2	6.82	125.34	119.20
1	C	81	LYS	CA-C-N	6.76	128.29	119.84
1	C	81	LYS	C-N-CA	6.76	128.29	119.84
1	C	208	HIS	CA-CB-CG	-6.61	107.19	113.80
1	D	2	GLY	CA-C-N	6.56	135.10	121.32
1	D	2	GLY	C-N-CA	6.56	135.10	121.32
1	D	117	TRP	O-C-N	6.51	130.37	122.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	156	ASP	CA-C-O	6.49	123.51	119.29
1	B	208	HIS	CA-CB-CG	-6.47	107.33	113.80
1	C	87	ASP	CA-C-O	6.42	127.22	120.42
1	B	113	GLY	CA-C-N	6.27	128.86	120.58
1	B	113	GLY	C-N-CA	6.27	128.86	120.58
1	A	60	VAL	CA-C-O	6.26	127.23	120.47
1	A	134	ARG	CD-NE-CZ	6.25	133.16	124.40
1	B	137	ALA	CA-C-N	6.23	133.44	121.54
1	B	137	ALA	C-N-CA	6.23	133.44	121.54
1	B	120	TRP	CA-CB-CG	6.16	125.30	113.60
1	B	21	ARG	NE-CZ-NH1	-6.16	115.34	121.50
1	D	8	PHE	CA-CB-CG	6.12	119.92	113.80
1	A	158	GLN	CA-C-N	6.12	131.84	121.14
1	A	158	GLN	C-N-CA	6.12	131.84	121.14
1	A	164	GLY	CA-C-O	-6.12	115.51	121.75
1	D	117	TRP	CA-C-O	-6.08	115.18	121.87
1	A	207	HIS	CA-C-N	5.99	130.63	122.07
1	A	207	HIS	C-N-CA	5.99	130.63	122.07
1	D	147	ASP	CA-CB-CG	-5.86	106.74	112.60
1	A	192	VAL	N-CA-C	-5.86	106.70	111.91
1	D	54	ALA	O-C-N	5.78	128.25	122.12
1	A	128	HIS	CA-CB-CG	5.69	119.49	113.80
1	B	79	ARG	CA-C-O	-5.69	115.19	121.33
1	C	204	ILE	CB-CA-C	-5.63	104.67	112.04
1	A	50	GLN	CA-C-O	-5.62	114.05	121.08
1	B	101	GLN	CA-C-N	-5.60	114.54	122.94
1	B	101	GLN	C-N-CA	-5.60	114.54	122.94
1	A	56	ARG	CA-C-O	5.60	126.48	120.55
1	B	128	HIS	CA-C-N	-5.59	115.89	122.93
1	B	128	HIS	C-N-CA	-5.59	115.89	122.93
1	C	18	GLY	CA-C-O	5.57	127.01	121.00
1	A	21	ARG	NE-CZ-NH1	-5.47	116.03	121.50
1	D	71	THR	CA-C-N	-5.42	113.56	122.67
1	D	71	THR	C-N-CA	-5.42	113.56	122.67
1	C	195	LEU	CA-C-O	5.41	124.76	118.97
1	D	159	ALA	CA-C-N	-5.39	113.40	120.95
1	D	159	ALA	C-N-CA	-5.39	113.40	120.95
1	A	25	GLU	CA-C-N	-5.39	112.12	120.31
1	A	25	GLU	C-N-CA	-5.39	112.12	120.31
1	B	75	ARG	CD-NE-CZ	5.39	131.94	124.40
1	A	49	PRO	CA-C-N	5.36	129.96	122.08
1	A	49	PRO	C-N-CA	5.36	129.96	122.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	81	LYS	CA-C-N	5.33	126.50	119.84
1	D	81	LYS	C-N-CA	5.33	126.50	119.84
1	A	45	HIS	CA-CB-CG	-5.25	108.55	113.80
1	D	195	LEU	N-CA-C	-5.25	107.25	113.97
1	C	28	GLU	O-C-N	5.25	127.49	122.03
1	B	207	HIS	CA-CB-CG	5.25	119.05	113.80
1	B	39	PRO	CA-C-N	-5.24	113.74	122.64
1	B	39	PRO	C-N-CA	-5.24	113.74	122.64
1	C	69	ARG	CA-C-N	-5.23	115.50	122.72
1	C	69	ARG	C-N-CA	-5.23	115.50	122.72
1	C	164	GLY	CA-C-O	-5.22	115.87	121.35
1	B	9	GLY	CA-C-N	-5.17	113.16	121.32
1	B	9	GLY	C-N-CA	-5.17	113.16	121.32
1	B	191	SER	CA-C-N	-5.16	112.67	121.97
1	B	191	SER	C-N-CA	-5.16	112.67	121.97
1	B	24	VAL	O-C-N	5.16	126.94	121.89
1	B	46[A]	ARG	CA-C-O	5.15	125.67	119.59
1	B	46[B]	ARG	CA-C-O	5.15	125.67	119.59
1	C	21	ARG	NE-CZ-NH1	-5.05	116.45	121.50
1	A	35	LEU	CA-C-O	-5.02	114.93	120.30
1	D	66	GLY	CA-C-O	5.02	125.67	119.25
1	A	67	VAL	O-C-N	5.01	128.59	123.03

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	1694	0	1692	104	0
1	B	1682	0	1678	82	0
1	C	1701	0	1704	112	0
1	D	1710	0	1725	80	0
2	A	13	0	5	3	0
2	B	13	0	5	4	0
2	C	13	0	5	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	13	0	5	6	0
3	A	22	0	12	1	0
3	B	22	0	12	1	0
3	C	22	0	12	2	0
3	D	22	0	11	1	0
4	A	164	0	0	15	0
4	B	215	0	0	16	0
4	C	225	0	0	21	0
4	D	209	0	0	17	0
All	All	7740	0	6866	375	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 27.

All (375) 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:79[C]:ARG:NH2	1:A:83:SER:CB	1.90	1.31
1:A:79[C]:ARG:NH2	1:A:83:SER:HB3	1.01	1.30
1:A:79[C]:ARG:CZ	1:A:83:SER:HB3	1.62	1.30
1:B:208:HIS:HB3	1:B:211:ARG:HD3	1.37	1.02
1:B:17[A]:ILE:HG23	1:D:194:PHE:HB3	1.45	0.98
1:B:52:SER:H	1:B:55:GLN:HE21	1.10	0.97
1:A:79[C]:ARG:HH21	1:A:83:SER:CB	1.65	0.96
1:D:26:MET:HE2	1:D:105:LEU:HD22	1.47	0.95
1:A:79[C]:ARG:CZ	1:A:83:SER:CB	2.34	0.94
1:A:182:ILE:HG23	1:A:192:VAL:HG11	1.47	0.93
1:A:79[C]:ARG:HH21	1:A:83:SER:HB3	1.15	0.92
1:C:145:LEU:HD22	1:C:148:LEU:HD22	1.51	0.91
1:C:26:MET:HE2	1:C:105:LEU:HD22	1.62	0.80
1:B:154:VAL:HG11	1:B:166:GLN:HE21	1.48	0.79
1:C:52:SER:H	1:C:55:GLN:HE21	1.31	0.78
1:D:158:GLN:HG3	4:D:962:HOH:O	1.84	0.78
1:A:85:THR:OG1	3:A:302:NCN:H2	1.84	0.78
1:C:38:LEU:HD11	1:C:75[B]:ARG:HH12	1.48	0.77
1:C:11[B]:THR:HG23	4:C:778:HOH:O	1.85	0.77
1:A:156:ASP:HB3	1:A:159:ALA:HB2	1.68	0.76
1:B:105:LEU:HD11	1:B:132:LEU:HD23	1.67	0.74
1:A:211:ARG:HD2	4:A:1073:HOH:O	1.86	0.74
1:D:85:THR:OG1	3:D:3302:NCN:H2	1.88	0.74
1:B:151:ALA:HB1	4:B:856:HOH:O	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:5[A]:ILE:HD13	1:D:34:GLU:HB3	1.70	0.73
1:C:178:SER:HB2	2:C:2301:CIT:O1	1.89	0.73
1:B:17[B]:ILE:HG12	1:B:195:LEU:O	1.89	0.72
1:C:28:GLU:HA	1:C:69:ARG:HH22	1.53	0.72
1:A:113:GLY:O	1:A:116:THR:HG23	1.89	0.71
1:C:161:LYS:H	1:C:166:GLN:NE2	1.89	0.70
1:A:79[C]:ARG:NE	1:A:83:SER:HB2	2.06	0.70
1:A:67:VAL:HG11	1:A:70:LEU:HD12	1.74	0.69
1:A:151:ALA:HB1	4:A:1046:HOH:O	1.91	0.69
1:A:158:GLN:HG2	4:A:590:HOH:O	1.92	0.69
1:A:152[B]:ARG:HG2	1:A:152[B]:ARG:HH21	1.55	0.69
1:A:196[A]:VAL:HG23	1:A:200:VAL:HB	1.74	0.69
1:D:4:ARG:HH11	1:D:101:GLN:HE22	1.40	0.69
1:D:143[A]:GLU:HB3	4:D:564:HOH:O	1.93	0.69
1:B:201:LEU:O	1:B:205:GLU:HG3	1.92	0.68
1:C:86:ILE:HD13	1:C:123:LEU:HD23	1.76	0.68
1:C:152[A]:ARG:HD2	1:C:165:GLY:O	1.94	0.68
1:B:2:GLY:HA2	4:B:617:HOH:O	1.93	0.68
1:A:124:LEU:HD23	1:A:129:ILE:HD13	1.74	0.68
3:B:1302:NCN:O3P	3:B:1302:NCN:H3'	1.93	0.68
1:A:79[C]:ARG:NE	1:A:83:SER:CB	2.57	0.68
1:B:86:ILE:HD11	1:B:122:ALA:HB3	1.76	0.67
1:A:208:HIS:O	1:A:211:ARG:HG3	1.95	0.67
1:A:118:HIS:HA	1:C:143:GLU:OE2	1.95	0.67
1:D:104:MET:HE3	1:D:106:ILE:HD11	1.77	0.66
1:D:143[B]:GLU:HB2	4:D:564:HOH:O	1.96	0.66
1:D:182[B]:ILE:HG23	1:D:192:VAL:HG11	1.76	0.66
1:D:75:ARG:HD3	4:D:1066:HOH:O	1.96	0.65
1:B:83:SER:HB3	4:B:548:HOH:O	1.97	0.65
1:A:19:HIS:HE2	2:A:301:CIT:H41	1.62	0.64
1:C:112:CYS:HB3	1:C:142:PRO:HD3	1.78	0.64
1:A:61:GLU:HG2	4:A:442:HOH:O	1.98	0.64
1:C:50[A]:GLN:NE2	1:D:158:GLN:HG2	2.12	0.64
1:B:75:ARG:HD2	1:B:88:THR:HG23	1.79	0.64
1:D:113:GLY:O	1:D:116:THR:HG23	1.98	0.64
1:B:152:ARG:HD2	1:B:165:GLY:O	1.99	0.63
1:A:58:ALA:O	1:A:62:ARG:HG2	1.98	0.63
1:C:38:LEU:HD11	1:C:75[B]:ARG:HH22	1.63	0.63
1:B:42:ARG:HA	1:B:82:PRO:HG3	1.80	0.62
1:D:19:HIS:HE2	2:D:3301:CIT:H41	1.62	0.62
1:C:81:LYS:HD2	1:C:82:PRO:HD2	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:LEU:HD11	1:C:167:ILE:HD11	1.81	0.62
1:A:109[A]:ASP:HB2	4:A:409:HOH:O	2.00	0.62
3:C:2302:NCN:O3P	3:C:2302:NCN:H3'	1.99	0.62
1:A:79[C]:ARG:HE	1:A:83:SER:HB2	1.63	0.62
1:B:141:PRO:HG2	4:B:689:HOH:O	1.99	0.62
1:A:79[C]:ARG:HH21	1:A:83:SER:CA	2.13	0.62
1:C:204:ILE:HD13	1:C:209:LEU:HD12	1.81	0.62
1:A:26:MET:HE3	1:A:105:LEU:HD22	1.81	0.61
1:D:38:LEU:HD21	1:D:88:THR:HG21	1.81	0.61
1:B:178:SER:HB2	2:B:1301:CIT:O2	2.00	0.61
1:A:157:PRO:HG2	4:A:590:HOH:O	1.99	0.61
1:D:115:PRO:HA	1:D:120:TRP:CD2	2.35	0.61
1:D:142:PRO:HB2	1:D:144:SER:OG	2.00	0.61
1:D:121:GLU:HG3	4:D:1004:HOH:O	2.00	0.60
1:C:16:HIS:HE2	2:C:2301:CIT:C6	2.15	0.60
1:C:79:ARG:HH11	1:C:79:ARG:HG2	1.67	0.60
1:D:11[B]:THR:HG23	4:D:888:HOH:O	2.02	0.60
1:D:75:ARG:HA	1:D:78:GLN:HG3	1.83	0.60
1:C:38:LEU:HD11	1:C:75[B]:ARG:NH1	2.17	0.60
1:C:161:LYS:H	1:C:166:GLN:HE22	1.48	0.60
1:A:119:ARG:HG3	1:C:143:GLU:OE1	2.02	0.59
1:B:69:ARG:HG3	1:B:69:ARG:HH11	1.68	0.59
1:D:3[A]:LYS:HD2	4:D:1006:HOH:O	2.01	0.59
1:D:172:GLN:HB3	4:D:938:HOH:O	2.02	0.58
1:A:182:ILE:HD12	1:A:192:VAL:HG12	1.85	0.58
1:B:32:LEU:O	1:B:69:ARG:HD2	2.02	0.58
1:B:155:ALA:O	1:B:157:PRO:HD3	2.03	0.58
1:A:124:LEU:CD2	1:A:129:ILE:HD13	2.33	0.58
1:C:32:LEU:O	1:C:69:ARG:HD2	2.03	0.58
1:A:134:ARG:CZ	1:A:174:PRO:HG2	2.34	0.58
1:C:96:LEU:HD13	1:C:100:ASP:CB	2.33	0.58
1:C:7[B]:LEU:HD11	1:C:38:LEU:HD22	1.85	0.58
1:D:3[C]:LYS:HE2	1:D:100:ASP:OD2	2.04	0.58
1:A:142:PRO:HB2	1:A:144:SER:OG	2.04	0.57
1:D:16:HIS:HE2	2:D:3301:CIT:C6	2.17	0.57
1:C:17[B]:ILE:HG22	1:C:195:LEU:O	2.03	0.57
3:C:2302:NCN:H6	3:C:2302:NCN:H5'1	1.86	0.57
1:B:73:ASP:OD2	1:B:75:ARG:HB3	2.05	0.57
1:B:133:GLN:HE22	1:B:139:SER:CB	2.17	0.57
1:C:44:PRO:HD2	4:C:843:HOH:O	2.05	0.57
1:B:26[A]:MET:HE2	1:B:105:LEU:HB2	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:108:TRP:O	1:C:111:PHE:HB3	2.04	0.57
1:C:152[C]:ARG:HG2	1:C:152[C]:ARG:HH21	1.68	0.57
1:D:154:VAL:HG21	1:D:160:LEU:HD23	1.86	0.57
1:D:172:GLN:O	1:D:174:PRO:HD3	2.04	0.57
1:C:75[B]:ARG:NH1	1:C:88:THR:HG23	2.20	0.57
1:D:4:ARG:HG2	1:D:101:GLN:NE2	2.20	0.56
1:C:7[B]:LEU:HG	1:C:38:LEU:HD22	1.87	0.56
1:C:205:GLU:HG2	4:C:783:HOH:O	2.04	0.56
1:C:7[B]:LEU:CG	1:C:38:LEU:HD22	2.35	0.56
1:C:69:ARG:HA	4:C:758:HOH:O	2.05	0.56
1:D:88:THR:O	1:D:92:VAL:HG23	2.06	0.56
1:B:67:VAL:HG11	1:B:70:LEU:HD12	1.86	0.56
1:C:142:PRO:HB2	1:C:144:SER:OG	2.06	0.56
1:A:105:LEU:O	1:A:106:ILE:HD13	2.06	0.56
1:B:42:ARG:CA	1:B:82:PRO:HG3	2.36	0.55
1:A:68:GLU:O	1:A:68:GLU:HG2	2.06	0.55
1:C:38:LEU:CD1	1:C:73:ASP:HB3	2.37	0.55
1:D:157:PRO:HG2	4:D:715:HOH:O	2.05	0.55
1:B:176:ALA:HB3	4:D:1053:HOH:O	2.06	0.55
1:C:17[B]:ILE:CG2	1:C:177:VAL:HG21	2.35	0.55
1:D:19:HIS:NE2	2:D:3301:CIT:H41	2.21	0.55
1:A:182:ILE:HD11	1:A:195:LEU:HD13	1.88	0.55
1:C:143:GLU:HG2	4:C:466:HOH:O	2.06	0.55
1:C:83:SER:HB2	4:C:740:HOH:O	2.06	0.55
1:C:59:MET:SD	1:C:204:ILE:HD11	2.46	0.55
2:D:3301:CIT:H42	4:D:888:HOH:O	2.06	0.55
1:B:134:ARG:CZ	1:B:174:PRO:HG2	2.37	0.55
1:D:46:ARG:HG3	1:D:46:ARG:HH11	1.70	0.55
1:B:178:SER:HB2	2:B:1301:CIT:O6	2.08	0.54
1:A:108:TRP:HH2	1:A:141:PRO:HD3	1.72	0.54
1:A:111:PHE:CE2	1:A:145:LEU:HD13	2.43	0.54
1:D:86:ILE:O	1:D:90:GLU:HG3	2.08	0.54
1:C:7[B]:LEU:CD1	1:C:38:LEU:HD22	2.38	0.53
1:C:41:ALA:HB2	1:C:76:GLU:O	2.08	0.53
1:C:154:VAL:HG22	1:C:168:THR:HB	1.91	0.53
1:A:201:LEU:O	1:A:205:GLU:HG3	2.09	0.53
1:C:34:GLU:CD	1:C:36:ARG:HE	2.17	0.53
1:D:4:ARG:NH1	1:D:101:GLN:HE22	2.06	0.53
1:D:109:ASP:HB2	4:D:1023:HOH:O	2.07	0.53
1:B:138:ASP:O	1:B:139:SER:HB2	2.09	0.53
1:A:115:PRO:HB3	1:A:120:TRP:CZ2	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17[B]:ILE:HD12	1:B:175:LEU:HD13	1.91	0.53
1:A:119:ARG:H	1:C:143:GLU:CD	2.17	0.53
1:C:38:LEU:CD1	1:C:75[B]:ARG:HH22	2.22	0.52
1:B:86:ILE:HD12	1:B:123:LEU:HD23	1.91	0.52
1:C:148:LEU:HD11	1:C:167:ILE:HD11	1.91	0.52
1:A:79[C]:ARG:NH2	1:A:83:SER:CA	2.69	0.52
1:B:124:LEU:HD21	1:B:129:ILE:HG13	1.92	0.52
1:D:156:ASP:O	1:D:159:ALA:HB3	2.10	0.51
1:A:26:MET:HE2	1:A:105:LEU:HB2	1.91	0.51
1:B:7:LEU:HD11	1:B:38:LEU:HD13	1.93	0.51
1:A:182:ILE:HD12	1:A:192:VAL:CG1	2.41	0.51
1:A:146:ARG:NE	1:A:146:ARG:HA	2.25	0.51
1:D:3[C]:LYS:HE2	1:D:100:ASP:CG	2.36	0.51
1:B:26[B]:MET:SD	1:B:105:LEU:HD13	2.51	0.51
1:A:116:THR:HA	4:A:862:HOH:O	2.11	0.50
1:B:52:SER:H	1:B:55:GLN:NE2	1.94	0.50
1:C:120:TRP:CE2	1:C:121:GLU:HG3	2.46	0.50
1:B:106:ILE:HD11	1:B:111:PHE:HD1	1.76	0.50
1:C:51:VAL:HA	1:C:55:GLN:NE2	2.27	0.50
1:A:203:TYR:HA	4:A:468:HOH:O	2.10	0.50
1:D:111:PHE:CZ	1:D:145:LEU:HD13	2.46	0.50
1:D:156:ASP:HB3	1:D:159:ALA:HB2	1.93	0.50
1:C:156:ASP:OD2	1:C:158[B]:GLN:OE1	2.30	0.50
1:C:38:LEU:HD11	1:C:75[B]:ARG:NH2	2.27	0.49
1:C:81:LYS:O	1:C:82:PRO:O	2.29	0.49
1:B:26[B]:MET:SD	1:B:105:LEU:HD22	2.52	0.49
1:C:9:GLY:HA2	1:C:38:LEU:HB3	1.95	0.49
1:A:146:ARG:HA	1:A:146:ARG:HE	1.78	0.49
1:B:74:PRO:HG2	1:B:78:GLN:OE1	2.13	0.49
1:C:7[A]:LEU:HD12	1:C:36:ARG:HB2	1.94	0.49
1:C:163:PRO:HG2	4:C:846:HOH:O	2.12	0.49
1:A:156:ASP:O	1:A:159:ALA:HB3	2.13	0.49
1:B:9:GLY:HA2	1:B:38:LEU:HB3	1.95	0.49
1:A:19:HIS:NE2	2:A:301:CIT:H41	2.26	0.49
1:B:68:GLU:O	1:B:69:ARG:HB2	2.13	0.49
1:C:40:ASN:HB3	4:C:869:HOH:O	2.12	0.49
1:C:96:LEU:HD13	1:C:100:ASP:HB2	1.95	0.48
1:D:19:HIS:HE2	2:D:3301:CIT:C4	2.26	0.48
1:C:17[B]:ILE:HD13	1:C:17[B]:ILE:HG21	1.54	0.48
1:A:121:GLU:HG3	4:A:425:HOH:O	2.14	0.48
1:C:147:ASP:HB3	4:C:836:HOH:O	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:137:ALA:O	1:C:138:ASP:O	2.32	0.48
1:A:166:GLN:NE2	4:A:1206:HOH:O	2.46	0.48
1:D:3[C]:LYS:HG3	1:D:100:ASP:OD1	2.13	0.48
1:A:107:GLY:HA2	1:A:132:LEU:HB3	1.96	0.48
1:D:139:SER:O	1:D:139:SER:OG	2.31	0.48
1:C:166:GLN:HG3	4:C:1210:HOH:O	2.13	0.48
1:D:68:GLU:O	1:D:69:ARG:HB2	2.14	0.48
1:B:208:HIS:CB	1:B:211:ARG:HH21	2.26	0.47
1:C:38:LEU:HD12	1:C:73:ASP:HB3	1.96	0.47
1:C:61:GLU:HG2	1:C:72:VAL:HG11	1.95	0.47
1:C:156:ASP:OD2	1:C:158[B]:GLN:HG3	2.13	0.47
1:D:17[A]:ILE:HD11	4:D:889:HOH:O	2.12	0.47
1:A:28:GLU:HG2	1:A:69:ARG:HH22	1.79	0.47
1:C:7[A]:LEU:HD22	1:C:102:LEU:HD11	1.95	0.47
1:C:207:HIS:HD2	4:C:788:HOH:O	1.96	0.47
1:D:136:ASP:OD1	1:D:136:ASP:O	2.32	0.47
1:B:125:ASP:HB3	4:B:622:HOH:O	2.14	0.47
1:B:93:ARG:HG2	4:B:557:HOH:O	2.13	0.47
1:B:193:ARG:O	1:D:21:ARG:HD3	2.14	0.47
1:D:93:ARG:HA	1:D:93:ARG:HD3	1.73	0.47
2:B:1301:CIT:O4	2:B:1301:CIT:O7	2.28	0.47
1:A:4:ARG:C	1:A:5:ILE:HD13	2.40	0.47
1:A:36:ARG:HG2	1:A:71:THR:OG1	2.15	0.47
1:C:135:PRO:CG	1:C:173:THR:HG22	2.45	0.47
1:D:48:THR:HG23	1:D:49:PRO:HD2	1.97	0.47
1:B:41:ALA:HB2	1:B:76:GLU:O	2.15	0.46
1:B:179:ALA:N	2:B:1301:CIT:O6	2.47	0.46
1:A:160:LEU:HA	1:A:166:GLN:OE1	2.16	0.46
1:D:211[B]:ARG:NH1	1:D:212:ALA:O	2.49	0.46
1:C:96:LEU:HD13	1:C:100:ASP:HB3	1.97	0.46
1:D:81:LYS:HG3	1:D:82:PRO:N	2.23	0.46
1:B:5:ILE:HD12	1:B:96:LEU:CD1	2.45	0.46
1:C:156:ASP:CG	1:C:158[B]:GLN:HG3	2.41	0.46
1:A:79[A]:ARG:NH2	1:A:80:ASP:OD1	2.49	0.46
1:D:5[A]:ILE:HD13	1:D:34:GLU:CB	2.41	0.46
1:A:186:LEU:HD13	1:A:210:TYR:HB2	1.98	0.46
1:B:89:LEU:HD23	1:B:89:LEU:HA	1.66	0.46
1:B:108:TRP:CZ2	1:B:139:SER:HA	2.51	0.46
1:A:81:LYS:O	1:A:83:SER:N	2.49	0.46
1:A:134:ARG:NH2	1:A:174:PRO:HG2	2.31	0.46
1:A:135:PRO:O	1:A:136:ASP:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:ASN:HD22	1:A:40:ASN:N	2.13	0.46
1:D:79:ARG:NH1	4:D:1100:HOH:O	2.49	0.46
1:D:108:TRP:NE1	1:D:138:ASP:O	2.49	0.46
1:A:19:HIS:HE2	2:A:301:CIT:C4	2.28	0.45
1:A:105:LEU:C	1:A:106:ILE:HD13	2.41	0.45
1:C:3:LYS:NZ	4:C:1183:HOH:O	2.49	0.45
1:C:29:GLN:NE2	4:C:822:HOH:O	2.50	0.45
1:C:75[B]:ARG:HG2	1:C:75[B]:ARG:HH11	1.80	0.45
1:C:152[B]:ARG:NH1	4:C:607:HOH:O	2.49	0.45
1:A:17[B]:ILE:HG12	1:A:195:LEU:O	2.16	0.45
1:A:84:TYR:HB2	1:A:87:ASP:OD2	2.15	0.45
1:B:69:ARG:NH2	4:B:510:HOH:O	2.49	0.45
1:D:101:GLN:HG2	1:D:162:GLY:O	2.17	0.45
1:B:108:TRP:HZ3	1:B:169:PHE:HE2	1.63	0.45
1:B:133:GLN:NE2	1:B:139:SER:OG	2.50	0.45
1:C:2:GLY:N	1:C:99:ASP:O	2.50	0.45
1:C:26:MET:HE2	1:C:105:LEU:CD2	2.40	0.45
1:C:34:GLU:OE2	1:C:36:ARG:NE	2.50	0.45
1:C:152[C]:ARG:HG2	1:C:152[C]:ARG:NH2	2.28	0.45
1:D:29:GLN:NE2	4:D:1013:HOH:O	2.49	0.45
1:B:132:LEU:HD21	1:B:172:GLN:CD	2.42	0.45
1:A:118:HIS:ND1	1:C:143:GLU:OE1	2.50	0.45
1:B:46[A]:ARG:NH2	4:B:657:HOH:O	2.50	0.45
1:B:114:LEU:HA	1:B:117:TRP:CD1	2.51	0.45
1:D:26:MET:HE3	1:D:35:LEU:HD21	1.99	0.45
1:C:211:ARG:HG2	4:C:731:HOH:O	2.16	0.45
1:D:102:LEU:HD23	1:D:127:CYS:HB3	1.99	0.45
1:D:143[B]:GLU:OE2	1:D:144:SER:N	2.49	0.45
1:A:101:GLN:NE2	4:A:433:HOH:O	2.49	0.45
1:B:45:HIS:HD2	4:B:648:HOH:O	1.99	0.45
1:B:118:HIS:CD2	1:B:119:ARG:HG3	2.52	0.45
1:B:174:PRO:HB3	4:B:545:HOH:O	2.17	0.45
1:C:93:ARG:HA	1:C:93:ARG:HD3	1.65	0.45
1:D:134[A]:ARG:HE	1:D:134[A]:ARG:HB2	1.29	0.45
1:A:52:SER:H	1:A:55:GLN:HE21	1.65	0.44
1:D:21:ARG:HD2	1:D:21:ARG:HA	1.75	0.44
1:B:47:GLU:HB2	4:B:594:HOH:O	2.17	0.44
1:D:46:ARG:HG3	1:D:46:ARG:NH1	2.32	0.44
1:C:75[B]:ARG:CZ	1:C:88:THR:HG23	2.47	0.44
1:C:152[C]:ARG:NH1	4:C:607:HOH:O	2.49	0.44
1:A:160:LEU:HB2	4:A:1094:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:73:ASP:OD1	1:C:75[B]:ARG:NH2	2.50	0.44
1:A:42:ARG:HA	1:A:82:PRO:HG3	1.98	0.44
1:A:115:PRO:HA	1:A:120:TRP:CD2	2.53	0.44
1:B:138:ASP:OD1	1:B:138:ASP:N	2.50	0.44
1:C:38:LEU:HD21	1:C:88:THR:HG21	1.99	0.44
1:A:196[A]:VAL:CG2	1:A:200:VAL:HB	2.46	0.44
1:C:4:ARG:HH11	1:C:101:GLN:NE2	2.16	0.44
1:C:140:GLU:HA	1:C:141:PRO:HD3	1.85	0.44
1:D:46:ARG:HA	1:D:46:ARG:HD2	1.72	0.44
1:D:75:ARG:NH1	1:D:91:SER:OG	2.50	0.44
1:B:41:ALA:HB3	1:B:82:PRO:HA	2.00	0.44
1:D:146:ARG:HG3	1:D:147:ASP:N	2.27	0.44
1:A:45:HIS:H	1:A:45:HIS:CD2	2.36	0.43
1:A:175:LEU:H	1:A:175:LEU:HG	1.70	0.43
1:B:20:MET:HB3	1:D:194:PHE:CE1	2.53	0.43
1:C:119:ARG:NH2	4:C:851:HOH:O	2.50	0.43
1:D:4:ARG:HG2	1:D:101:GLN:HE21	1.83	0.43
1:B:50:GLN:NE2	4:B:1141:HOH:O	2.50	0.43
1:B:83:SER:OG	1:B:84:TYR:N	2.50	0.43
1:B:21:ARG:NH2	4:B:516:HOH:O	2.50	0.43
1:C:204:ILE:HD12	1:C:209:LEU:HB2	2.00	0.43
1:D:172:GLN:C	1:D:174:PRO:HD3	2.44	0.43
1:A:108:TRP:CZ2	1:A:139:SER:HA	2.53	0.43
1:B:69:ARG:HG3	1:B:69:ARG:NH1	2.33	0.43
1:B:81:LYS:HB2	1:B:82:PRO:HD2	1.99	0.43
1:C:160:LEU:HA	1:C:166:GLN:NE2	2.34	0.43
1:C:178:SER:HA	2:C:2301:CIT:O6	2.19	0.43
1:A:115:PRO:HA	1:A:120:TRP:CG	2.54	0.43
1:A:140:GLU:HA	1:A:141:PRO:HD2	1.78	0.43
1:D:17[A]:ILE:CG2	1:D:175:LEU:HD13	2.48	0.43
1:A:134:ARG:HB2	1:A:135:PRO:HD2	1.99	0.43
1:A:153:SER:HA	1:A:167:ILE:O	2.19	0.43
1:A:62:ARG:HD2	1:A:62:ARG:HA	1.72	0.43
1:A:193:ARG:HG2	1:A:194:PHE:CD1	2.54	0.43
1:B:106:ILE:HD11	1:B:111:PHE:CD1	2.54	0.43
1:A:81:LYS:HG3	1:A:82:PRO:HD2	2.00	0.42
1:C:78:GLN:HE21	1:C:78:GLN:HB2	1.58	0.42
1:C:135:PRO:O	1:C:136:ASP:HB2	2.19	0.42
1:A:54:ALA:HB3	4:A:421:HOH:O	2.18	0.42
1:B:84:TYR:CD2	1:B:118:HIS:HB2	2.54	0.42
1:B:138:ASP:HB2	1:B:139:SER:H	1.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:138:ASP:OD1	1:C:138:ASP:N	2.50	0.42
1:A:139:SER:O	1:A:139:SER:OG	2.30	0.42
1:B:78:GLN:NE2	4:B:682:HOH:O	2.50	0.42
1:B:119:ARG:HE	1:B:119:ARG:HB3	1.62	0.42
1:B:133:GLN:HE22	1:B:139:SER:HB2	1.83	0.42
1:A:212:ALA:HA	1:A:213:PRO:HD2	1.85	0.42
1:A:92:VAL:O	1:A:96:LEU:HG	2.20	0.42
1:C:26:MET:HG3	1:C:105:LEU:HD13	2.02	0.42
1:C:154:VAL:CG2	1:C:168:THR:HB	2.48	0.42
1:D:21:ARG:HG3	1:D:175:LEU:HD11	2.02	0.42
1:D:28:GLU:HG2	4:D:936:HOH:O	2.19	0.42
1:A:154:VAL:HA	4:A:418:HOH:O	2.20	0.42
1:D:19:HIS:CE1	2:D:3301:CIT:H41	2.55	0.42
1:D:79:ARG:NE	1:D:83:SER:OG	2.49	0.42
1:C:134:ARG:O	1:C:137:ALA:HB3	2.19	0.42
1:D:134[A]:ARG:NH2	1:D:174:PRO:HD2	2.35	0.42
1:C:86:ILE:CD1	1:C:123:LEU:HD23	2.48	0.41
1:A:73:ASP:OD2	1:A:75:ARG:HB3	2.20	0.41
1:A:95:GLU:C	1:A:96:LEU:HD23	2.45	0.41
1:B:3:LYS:NZ	4:B:620:HOH:O	2.48	0.41
1:A:38:LEU:HG	1:A:76:GLU:OE1	2.20	0.41
1:C:119:ARG:HE	1:C:119:ARG:HB3	1.59	0.41
1:C:156:ASP:O	1:C:159:ALA:HB3	2.20	0.41
1:B:56:ARG:HH11	1:B:56:ARG:HD2	1.69	0.41
1:B:64:VAL:HG21	1:B:72:VAL:HG23	2.03	0.41
1:D:158:GLN:OE1	1:D:158:GLN:N	2.53	0.41
1:A:96:LEU:HD23	1:A:96:LEU:N	2.35	0.41
1:A:115:PRO:HA	1:A:120:TRP:CD1	2.55	0.41
1:C:111:PHE:O	1:C:114:LEU:HB2	2.20	0.41
1:C:131:VAL:O	1:C:170:VAL:HG22	2.20	0.41
1:D:168:THR:OG1	1:D:169:PHE:N	2.52	0.41
1:C:72:VAL:HG12	4:C:732:HOH:O	2.21	0.41
1:B:79:ARG:HD3	4:B:548:HOH:O	2.21	0.41
1:D:6:GLY:HA2	1:D:103:PHE:O	2.21	0.41
1:A:5:ILE:HD13	1:A:5:ILE:N	2.35	0.41
1:A:115:PRO:HB3	1:A:120:TRP:CE2	2.56	0.41
1:C:4:ARG:HH11	1:C:101:GLN:HE22	1.68	0.41
1:C:211:ARG:HD3	4:C:883:HOH:O	2.20	0.41
1:D:4:ARG:NH1	1:D:101:GLN:NE2	2.69	0.41
1:B:17[A]:ILE:HG22	1:B:175:LEU:HD12	2.03	0.40
1:B:186:LEU:HD21	1:B:192:VAL:HG22	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:149:LEU:HD23	1:C:149:LEU:HA	1.85	0.40
1:C:166:GLN:NE2	4:C:1210:HOH:O	2.54	0.40
1:A:151:ALA:HB3	1:A:152[B]:ARG:NH2	2.36	0.40
1:C:156:ASP:OD1	1:C:158[B]:GLN:HG3	2.21	0.40
1:D:27:ALA:HB1	1:D:69:ARG:NH1	2.37	0.40
1:D:68:GLU:H	1:D:68:GLU:CD	2.27	0.40
1:A:17[A]:ILE:HG22	1:A:175:LEU:CD1	2.52	0.40
1:A:115:PRO:HA	1:A:120:TRP:CE2	2.56	0.40
1:A:151:ALA:O	1:A:152[B]:ARG:HD3	2.21	0.40
1:D:86:ILE:HG22	4:D:945:HOH:O	2.21	0.40
1:A:197:PRO:HA	4:A:713:HOH:O	2.22	0.40
1:B:158:GLN:H	1:B:158:GLN:HG2	1.65	0.40
1:B:160:LEU:HD11	1:B:168:THR:HG22	2.03	0.40
1:C:208:HIS:ND1	1:C:211:ARG:NH2	2.70	0.40
1:D:17[A]:ILE:HG22	1:D:175:LEU:HD13	2.04	0.40
1:B:93:ARG:HA	1:B:93:ARG:HD3	1.71	0.40
1:C:26:MET:HE1	1:C:35:LEU:HD11	2.03	0.40
1:C:28:GLU:HG3	4:C:1196:HOH:O	2.21	0.40

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	217/242 (90%)	202 (93%)	8 (4%)	7 (3%)	3	0
1	B	215/242 (89%)	205 (95%)	7 (3%)	3 (1%)	9	2
1	C	218/242 (90%)	202 (93%)	14 (6%)	2 (1%)	14	4
1	D	220/242 (91%)	208 (94%)	9 (4%)	3 (1%)	9	2
All	All	870/968 (90%)	817 (94%)	38 (4%)	15 (2%)	8	1

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	136	ASP
1	A	138	ASP
1	B	138	ASP
1	B	139	SER
1	C	82	PRO
1	C	138	ASP
1	D	80	ASP
1	D	82	PRO
1	D	137	ALA
1	A	83	SER
1	A	79[A]	ARG
1	A	79[B]	ARG
1	A	79[C]	ARG
1	A	82	PRO
1	B	80	ASP

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	180/200 (90%)	160 (89%)	20 (11%)	6	1
1	B	178/200 (89%)	158 (89%)	20 (11%)	6	1
1	C	181/200 (90%)	163 (90%)	18 (10%)	7	1
1	D	183/200 (92%)	153 (84%)	30 (16%)	2	0
All	All	722/800 (90%)	634 (88%)	88 (12%)	6	1

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	5	ILE
1	A	34	GLU
1	A	75	ARG
1	A	78	GLN
1	A	81	LYS
1	A	91	SER

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Mol	Chain	Res	Type
1	A	116	THR
1	A	132	LEU
1	A	136	ASP
1	A	140	GLU
1	A	143	GLU
1	A	146	ARG
1	A	154	VAL
1	A	173	THR
1	A	175	LEU
1	A	178	SER
1	A	182	ILE
1	A	196[A]	VAL
1	A	196[B]	VAL
1	B	17[A]	ILE
1	B	17[B]	ILE
1	B	34	GLU
1	B	46[A]	ARG
1	B	46[B]	ARG
1	B	50	GLN
1	B	72	VAL
1	B	78	GLN
1	B	80	ASP
1	B	81	LYS
1	B	86	ILE
1	B	93	ARG
1	B	96	LEU
1	B	114	LEU
1	B	119	ARG
1	B	132	LEU
1	B	138	ASP
1	B	158	GLN
1	B	201	LEU
1	B	211	ARG
1	C	3	LYS
1	C	7[A]	LEU
1	C	7[B]	LEU
1	C	17[A]	ILE
1	C	17[B]	ILE
1	C	21	ARG
1	C	68	GLU
1	C	69	ARG
1	C	74	PRO

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Mol	Chain	Res	Type
1	C	78	GLN
1	C	79	ARG
1	C	82	PRO
1	C	93	ARG
1	C	96	LEU
1	C	132	LEU
1	C	138	ASP
1	C	148	LEU
1	C	181	GLN
1	D	5[A]	ILE
1	D	5[B]	ILE
1	D	17[A]	ILE
1	D	17[B]	ILE
1	D	26	MET
1	D	50	GLN
1	D	55	GLN
1	D	72	VAL
1	D	78	GLN
1	D	80	ASP
1	D	81	LYS
1	D	82	PRO
1	D	116	THR
1	D	132	LEU
1	D	133	GLN
1	D	134[A]	ARG
1	D	134[B]	ARG
1	D	134[C]	ARG
1	D	139	SER
1	D	140	GLU
1	D	141	PRO
1	D	143[A]	GLU
1	D	143[B]	GLU
1	D	146	ARG
1	D	158	GLN
1	D	173	THR
1	D	175	LEU
1	D	178	SER
1	D	211[A]	ARG
1	D	211[B]	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	40	ASN
1	A	45	HIS
1	A	55	GLN
1	A	101	GLN
1	A	126	HIS
1	A	207	HIS
1	B	45	HIS
1	B	55	GLN
1	B	101	GLN
1	B	118	HIS
1	B	133	GLN
1	B	166	GLN
1	B	207	HIS
1	C	29	GLN
1	C	55	GLN
1	C	78	GLN
1	C	101	GLN
1	C	118	HIS
1	C	133	GLN
1	C	166	GLN
1	C	207	HIS
1	D	29	GLN
1	D	40	ASN
1	D	45	HIS
1	D	50	GLN
1	D	55	GLN
1	D	101	GLN
1	D	133	GLN
1	D	207	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

8 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	CIT	D	3301	-	12,12,12	1.76	5 (41%)	17,17,17	3.38	9 (52%)
3	NCN	B	1302	-	21,23,23	2.50	10 (47%)	27,34,34	3.58	13 (48%)
2	CIT	A	301	-	12,12,12	1.86	4 (33%)	17,17,17	3.03	8 (47%)
3	NCN	C	2302	-	21,23,23	2.53	9 (42%)	27,34,34	3.65	14 (51%)
3	NCN	D	3302	-	21,23,23	2.57	12 (57%)	27,34,34	3.66	15 (55%)
2	CIT	B	1301	-	12,12,12	1.79	3 (25%)	17,17,17	3.60	8 (47%)
3	NCN	A	302	-	21,23,23	2.44	9 (42%)	27,34,34	3.34	10 (37%)
2	CIT	C	2301	-	12,12,12	1.80	3 (25%)	17,17,17	2.87	7 (41%)

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	CIT	D	3301	-	-	11/16/16/16	-
3	NCN	B	1302	-	-	3/14/30/30	0/2/2/2
2	CIT	A	301	-	-	4/16/16/16	-
3	NCN	C	2302	-	-	3/14/30/30	0/2/2/2
3	NCN	D	3302	-	-	4/14/30/30	0/2/2/2
2	CIT	B	1301	-	-	9/16/16/16	-
3	NCN	A	302	-	-	3/14/30/30	0/2/2/2
2	CIT	C	2301	-	-	6/16/16/16	-

All (55) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	302	NCN	C4-C3	5.54	1.47	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1302	NCN	C4-C3	5.49	1.47	1.39
3	D	3302	NCN	C4-C3	5.31	1.47	1.39
3	C	2302	NCN	O8-C7	4.77	1.36	1.22
3	C	2302	NCN	C4-C3	4.73	1.46	1.39
3	D	3302	NCN	O8-C7	4.52	1.35	1.22
3	B	1302	NCN	O8-C7	4.50	1.35	1.22
3	D	3302	NCN	C6-N1	4.45	1.45	1.35
3	A	302	NCN	O8-C7	4.33	1.35	1.22
3	A	302	NCN	C6-N1	4.28	1.45	1.35
3	C	2302	NCN	C6-N1	4.27	1.45	1.35
2	A	301	CIT	C3-C6	4.24	1.57	1.53
3	C	2302	NCN	C2-N1	4.14	1.39	1.35
3	B	1302	NCN	C6-N1	4.14	1.44	1.35
2	C	2301	CIT	O2-C1	-3.70	1.18	1.30
2	C	2301	CIT	C3-C6	3.69	1.57	1.53
3	C	2302	NCN	C3'-C4'	-3.57	1.43	1.53
2	B	1301	CIT	C3-C6	3.54	1.57	1.53
2	D	3301	CIT	O2-C1	-3.47	1.19	1.30
3	C	2302	NCN	O7-C7	-3.32	1.20	1.30
3	B	1302	NCN	C2-N1	3.17	1.38	1.35
2	B	1301	CIT	O2-C1	-3.12	1.20	1.30
3	A	302	NCN	O7-C7	-3.06	1.21	1.30
3	D	3302	NCN	O7-C7	-3.05	1.21	1.30
2	A	301	CIT	O2-C1	-3.02	1.20	1.30
3	A	302	NCN	C3'-C4'	-3.01	1.45	1.53
3	B	1302	NCN	O7-C7	-2.99	1.21	1.30
3	D	3302	NCN	C3'-C4'	-2.97	1.45	1.53
3	D	3302	NCN	C2-N1	2.85	1.38	1.35
3	D	3302	NCN	C5'-C4'	-2.85	1.43	1.51
3	A	302	NCN	C2-N1	2.83	1.38	1.35
3	B	1302	NCN	C3'-C4'	-2.70	1.46	1.53
3	B	1302	NCN	C3-C7	2.66	1.55	1.49
3	C	2302	NCN	C3-C7	2.63	1.55	1.49
3	D	3302	NCN	C5-C4	2.59	1.43	1.38
3	A	302	NCN	C3-C7	2.57	1.54	1.49
3	A	302	NCN	C5-C4	2.51	1.43	1.38
2	D	3301	CIT	C3-C6	2.48	1.56	1.53
3	C	2302	NCN	C5-C4	2.47	1.43	1.38
3	B	1302	NCN	O4'-C4'	2.47	1.50	1.45
3	D	3302	NCN	C3-C7	2.47	1.54	1.49
3	A	302	NCN	O4'-C4'	2.45	1.50	1.45
3	D	3302	NCN	O4'-C1'	2.43	1.44	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1302	NCN	C5-C4	2.42	1.43	1.38
3	D	3302	NCN	O4'-C4'	2.37	1.50	1.45
2	D	3301	CIT	C4-C5	2.35	1.57	1.50
2	B	1301	CIT	O1-C1	2.32	1.29	1.22
3	C	2302	NCN	O4'-C4'	2.27	1.50	1.45
2	C	2301	CIT	O1-C1	2.12	1.29	1.22
3	D	3302	NCN	C2'-C3'	2.11	1.59	1.53
2	D	3301	CIT	O1-C1	2.09	1.28	1.22
3	B	1302	NCN	C2'-C3'	2.04	1.58	1.53
2	D	3301	CIT	C4-C3	2.03	1.56	1.54
2	A	301	CIT	O1-C1	2.02	1.28	1.22
2	A	301	CIT	C4-C5	2.00	1.56	1.50

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2302	NCN	C5-C4-C3	-8.59	111.92	120.36
3	A	302	NCN	C4-C3-C7	-8.56	103.44	120.37
3	B	1302	NCN	C4-C3-C7	-8.06	104.43	120.37
3	D	3302	NCN	C5-C4-C3	-7.85	112.64	120.36
2	A	301	CIT	O6-C6-C3	-7.70	98.37	113.14
3	A	302	NCN	C5-C4-C3	-7.60	112.89	120.36
3	D	3302	NCN	C4-C3-C7	-7.56	105.40	120.37
3	B	1302	NCN	C5-C4-C3	-7.55	112.94	120.36
2	B	1301	CIT	O7-C3-C6	-7.55	98.25	108.96
3	B	1302	NCN	C2-C3-C7	7.54	132.72	119.78
3	C	2302	NCN	O7-C7-C3	7.25	133.44	114.84
3	D	3302	NCN	C2-C3-C7	7.10	131.97	119.78
2	D	3301	CIT	O6-C6-C3	-6.93	99.83	113.14
3	A	302	NCN	C5-C6-N1	-6.86	111.02	120.38
2	B	1301	CIT	C2-C3-C6	6.85	125.19	110.03
2	D	3301	CIT	C2-C3-C6	6.64	124.72	110.03
2	B	1301	CIT	O6-C6-C3	-6.49	100.68	113.14
3	C	2302	NCN	C6-C5-C4	6.30	128.53	119.45
2	C	2301	CIT	O6-C6-C3	-6.29	101.07	113.14
2	C	2301	CIT	O7-C3-C6	-6.25	100.10	108.96
3	C	2302	NCN	C4-C3-C7	-5.94	108.62	120.37
3	D	3302	NCN	O7-C7-C3	5.93	130.06	114.84
2	D	3301	CIT	O7-C3-C6	-5.76	100.79	108.96
3	C	2302	NCN	O8-C7-C3	-5.55	106.91	121.46
3	B	1302	NCN	O7-C7-C3	5.39	128.66	114.84
3	D	3302	NCN	C5-C6-N1	-5.33	113.12	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2302	NCN	C2-C3-C7	5.18	128.67	119.78
3	C	2302	NCN	C5-C6-N1	-5.14	113.38	120.38
3	A	302	NCN	C2-C3-C7	5.13	128.58	119.78
2	A	301	CIT	O5-C6-C3	4.94	131.67	122.09
2	A	301	CIT	O7-C3-C6	-4.91	102.00	108.96
3	A	302	NCN	O8-C7-C3	-4.89	108.64	121.46
3	D	3302	NCN	C6-C5-C4	4.87	126.47	119.45
3	B	1302	NCN	O8-C7-C3	-4.83	108.80	121.46
3	D	3302	NCN	C2'-C3'-C4'	-4.76	93.41	102.61
2	B	1301	CIT	C4-C3-C6	-4.67	99.71	110.03
3	B	1302	NCN	C5-C6-N1	-4.63	114.06	120.38
3	D	3302	NCN	C4'-O4'-C1'	-4.62	105.69	109.92
3	A	302	NCN	O7-C7-C3	4.55	126.50	114.84
3	B	1302	NCN	C2'-C3'-C4'	-4.41	94.08	102.61
2	B	1301	CIT	O5-C6-C3	4.34	130.50	122.09
3	D	3302	NCN	O8-C7-C3	-4.29	110.21	121.46
2	D	3301	CIT	O5-C6-C3	4.09	130.01	122.09
3	B	1302	NCN	C2-C3-C4	3.90	122.79	118.26
3	B	1302	NCN	C6-C5-C4	3.75	124.85	119.45
3	B	1302	NCN	O3'-C3'-C4'	-3.71	100.42	111.08
3	C	2302	NCN	O2'-C2'-C3'	3.69	123.64	111.82
3	C	2302	NCN	C2-N1-C1'	-3.57	111.26	119.13
2	C	2301	CIT	O2-C1-O1	-3.55	114.20	123.33
2	B	1301	CIT	O2-C1-O1	-3.55	114.20	123.33
2	A	301	CIT	O2-C1-C2	3.54	125.55	114.35
3	C	2302	NCN	C2-C3-C4	3.40	122.22	118.26
3	A	302	NCN	C6-C5-C4	3.38	124.32	119.45
2	C	2301	CIT	C2-C3-C6	3.36	117.46	110.03
2	D	3301	CIT	C4-C3-C6	-3.28	102.78	110.03
3	B	1302	NCN	C4'-O4'-C1'	-3.21	106.99	109.92
2	D	3301	CIT	C3-C4-C5	-3.05	105.59	113.92
3	A	302	NCN	O3'-C3'-C4'	-3.01	102.43	111.08
2	C	2301	CIT	O5-C6-C3	3.01	127.92	122.09
3	D	3302	NCN	O3'-C3'-C4'	-2.97	102.54	111.08
2	A	301	CIT	O1-C1-C2	-2.90	114.74	122.95
2	A	301	CIT	C4-C3-C6	-2.87	103.69	110.03
2	D	3301	CIT	O4-C5-C4	-2.86	105.30	114.35
2	B	1301	CIT	O1-C1-C2	-2.82	114.97	122.95
2	C	2301	CIT	O2-C1-C2	2.79	123.18	114.35
3	C	2302	NCN	O4'-C4'-C3'	2.61	110.34	105.15
3	A	302	NCN	O3'-C3'-C2'	2.60	120.16	111.82
3	C	2302	NCN	O3'-C3'-C4'	-2.59	103.63	111.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1302	NCN	C6-N1-C1'	-2.58	114.66	119.73
2	A	301	CIT	C4-C3-C2	2.35	115.35	109.31
2	A	301	CIT	C2-C3-C6	2.34	115.21	110.03
3	D	3302	NCN	C6-N1-C1'	-2.26	115.29	119.73
3	C	2302	NCN	O7-C7-O8	-2.23	118.56	123.35
3	C	2302	NCN	O3'-C3'-C2'	2.22	118.94	111.82
3	D	3302	NCN	C2-C3-C4	2.17	120.79	118.26
3	D	3302	NCN	O7-C7-O8	-2.12	118.79	123.35
2	B	1301	CIT	O4-C5-O3	-2.12	117.89	123.33
3	B	1302	NCN	O7-C7-O8	-2.11	118.82	123.35
3	D	3302	NCN	C6-N1-C2	2.08	123.65	121.88
2	D	3301	CIT	O1-C1-C2	-2.08	117.05	122.95
2	C	2301	CIT	C4-C3-C2	2.08	114.64	109.31
3	A	302	NCN	C2'-C3'-C4'	-2.05	98.65	102.61
3	D	3302	NCN	O5'-C5'-C4'	2.04	115.93	108.99
2	D	3301	CIT	O7-C3-C4	-2.02	104.78	109.38

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	CIT	C2-C3-C4-C5
2	A	301	CIT	O7-C3-C4-C5
2	A	301	CIT	C6-C3-C4-C5
2	B	1301	CIT	C1-C2-C3-O7
2	B	1301	CIT	C2-C3-C4-C5
2	B	1301	CIT	C6-C3-C4-C5
2	C	2301	CIT	C2-C3-C4-C5
2	C	2301	CIT	O7-C3-C4-C5
2	C	2301	CIT	C6-C3-C4-C5
2	D	3301	CIT	C6-C3-C4-C5
2	D	3301	CIT	C2-C3-C6-O5
3	D	3302	NCN	C5'-O5'-P-O1P
3	D	3302	NCN	C5'-O5'-P-O2P
3	D	3302	NCN	C5'-O5'-P-O3P
3	B	1302	NCN	O4'-C4'-C5'-O5'
2	D	3301	CIT	C1-C2-C3-O7
2	D	3301	CIT	C1-C2-C3-C4
2	D	3301	CIT	C1-C2-C3-C6
2	D	3301	CIT	C2-C3-C4-C5
3	B	1302	NCN	C3'-C4'-C5'-O5'
2	B	1301	CIT	C1-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
3	C	2302	NCN	O4'-C4'-C5'-O5'
2	B	1301	CIT	O7-C3-C4-C5
2	D	3301	CIT	O7-C3-C4-C5
3	C	2302	NCN	C3'-C4'-C5'-O5'
3	A	302	NCN	C4-C3-C7-O7
2	C	2301	CIT	C2-C3-C6-O5
2	A	301	CIT	C1-C2-C3-C6
3	C	2302	NCN	C4'-C5'-O5'-P
2	D	3301	CIT	C3-C4-C5-O3
3	B	1302	NCN	C4'-C5'-O5'-P
2	C	2301	CIT	O7-C3-C6-O5
2	B	1301	CIT	C4-C3-C6-O5
2	B	1301	CIT	C1-C2-C3-C6
3	A	302	NCN	O4'-C4'-C5'-O5'
2	B	1301	CIT	C3-C4-C5-O4
3	D	3302	NCN	C3'-C4'-C5'-O5'
2	D	3301	CIT	C4-C3-C6-O5
2	C	2301	CIT	C1-C2-C3-O7
2	B	1301	CIT	C3-C4-C5-O3
2	D	3301	CIT	O7-C3-C6-O5
3	A	302	NCN	C3'-C4'-C5'-O5'
2	D	3301	CIT	C3-C4-C5-O4

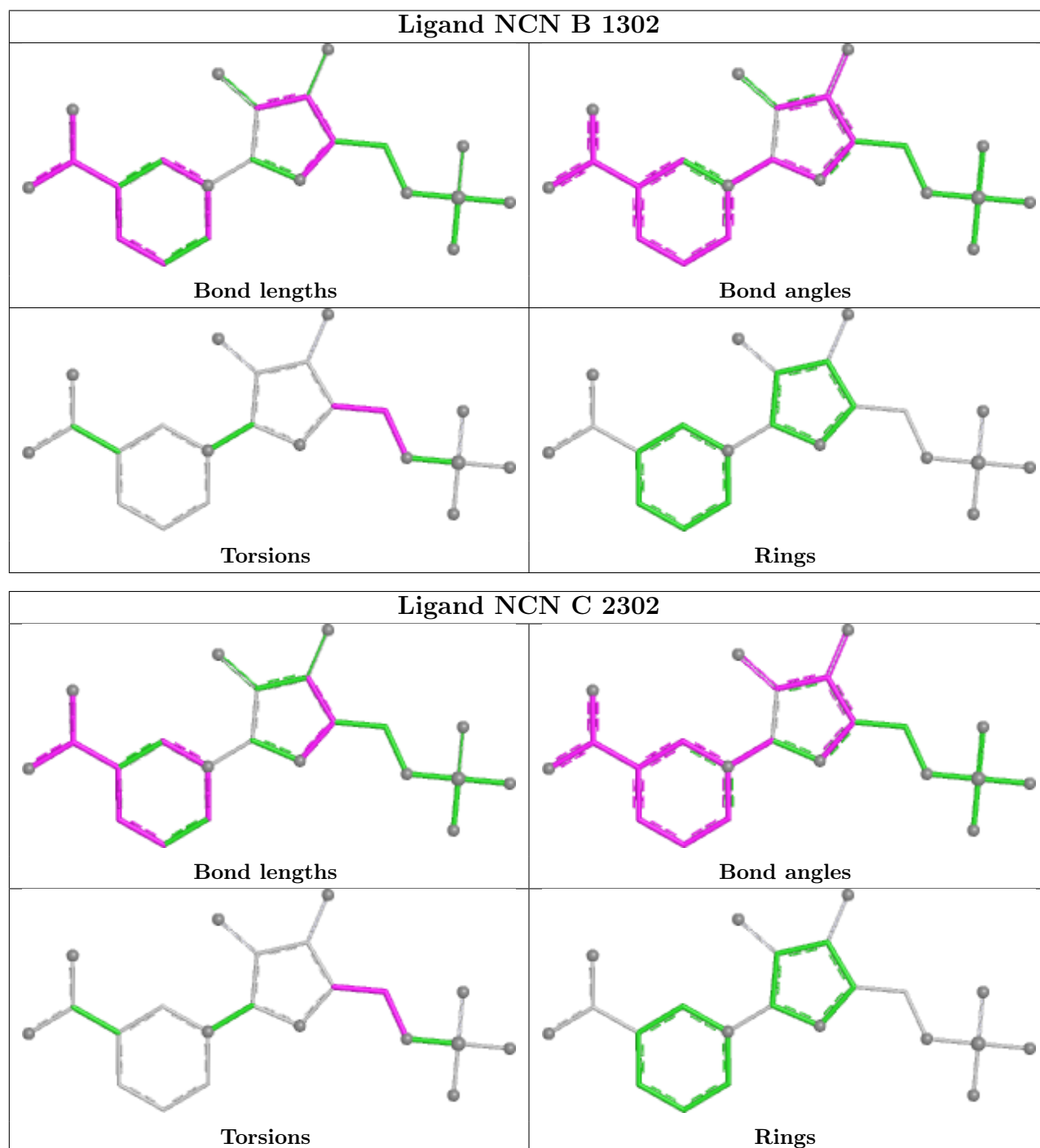
There are no ring outliers.

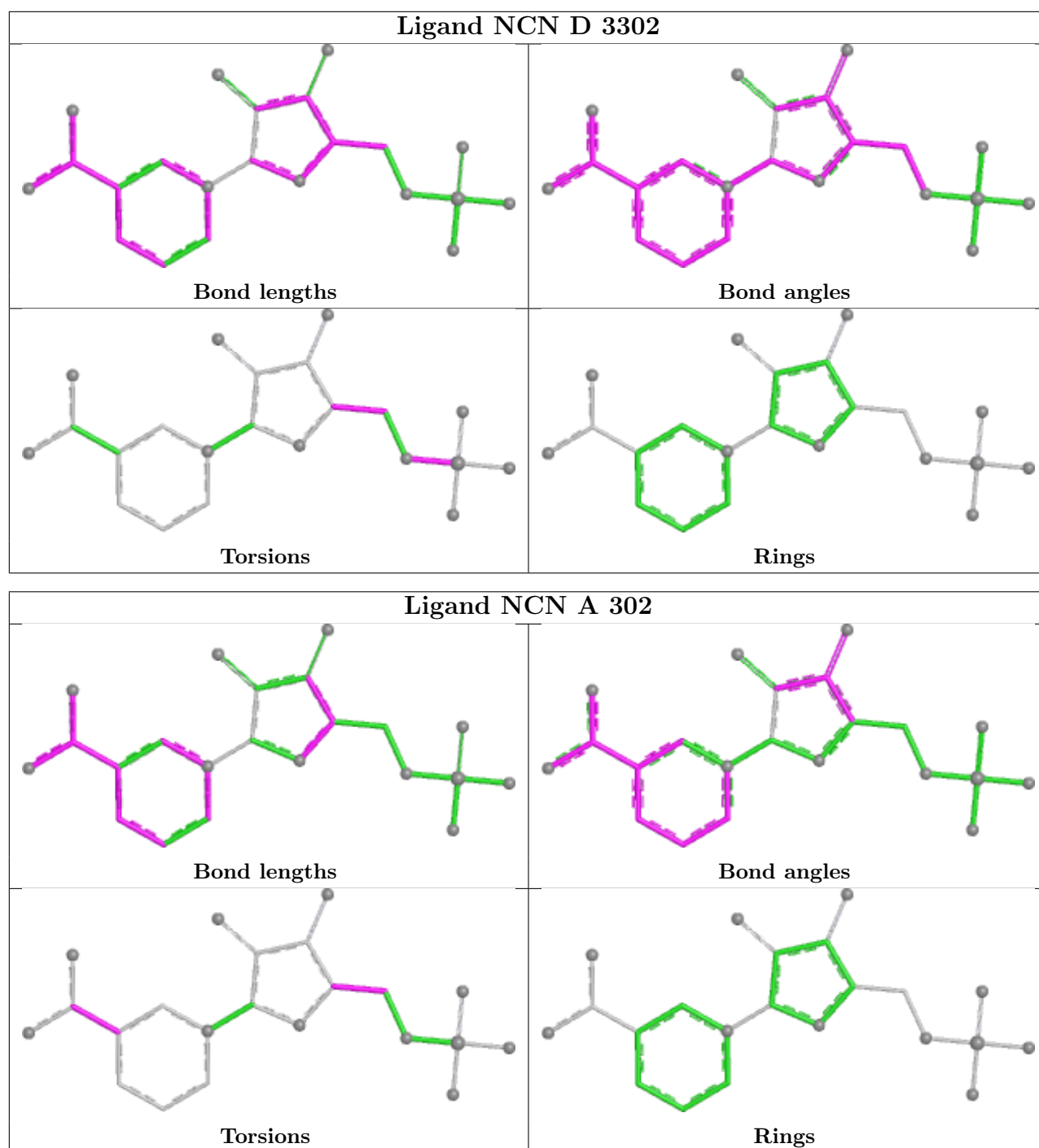
8 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	3301	CIT	6	0
3	B	1302	NCN	1	0
2	A	301	CIT	3	0
3	C	2302	NCN	2	0
3	D	3302	NCN	1	0
2	B	1301	CIT	4	0
3	A	302	NCN	1	0
2	C	2301	CIT	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	212/242 (87%)	-0.61	0 100 100	9, 24, 48, 88	7 (3%)
1	B	212/242 (87%)	-0.59	0 100 100	12, 23, 47, 73	5 (2%)
1	C	212/242 (87%)	-0.55	1 (0%) 87 89	10, 23, 53, 87	7 (3%)
1	D	212/242 (87%)	-0.56	0 100 100	8, 24, 52, 97	8 (3%)
All	All	848/968 (87%)	-0.58	1 (0%) 92 92	8, 24, 51, 97	27 (3%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	82	PRO	2.4

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	NCN	D	3302	22/22	0.96	0.09	25,47,75,95	0

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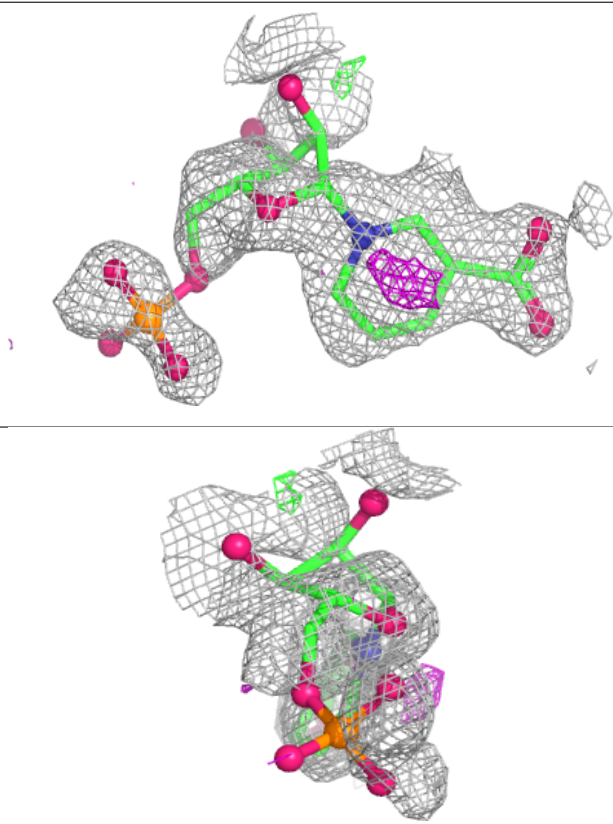
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NCN	B	1302	22/22	0.97	0.09	32,57,83,98	0
3	NCN	A	302	22/22	0.97	0.09	25,56,83,89	0
2	CIT	D	3301	13/13	0.98	0.07	18,38,70,76	0
2	CIT	A	301	13/13	0.98	0.07	14,45,67,75	0
2	CIT	B	1301	13/13	0.98	0.09	27,49,79,90	0
3	NCN	C	2302	22/22	0.98	0.07	18,42,74,81	0
2	CIT	C	2301	13/13	0.98	0.09	15,35,75,79	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.

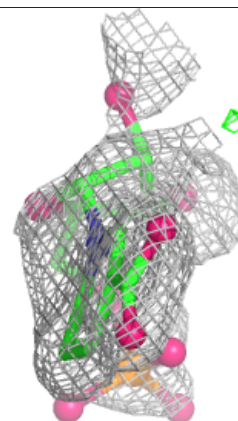
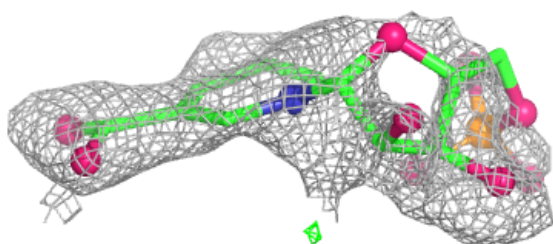
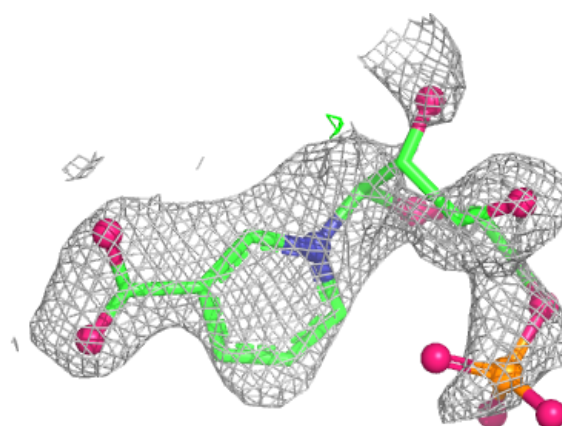
Electron density around NCN D 3302:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

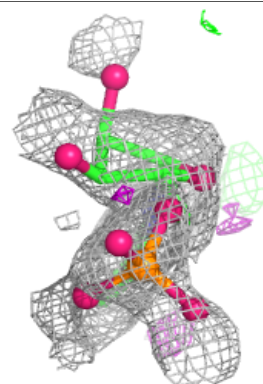
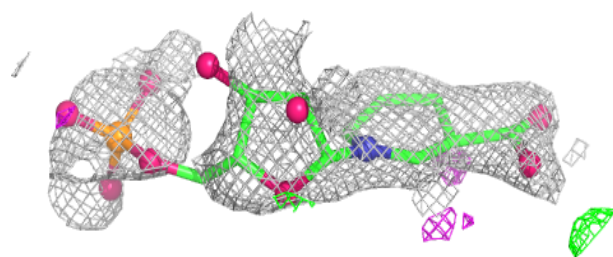
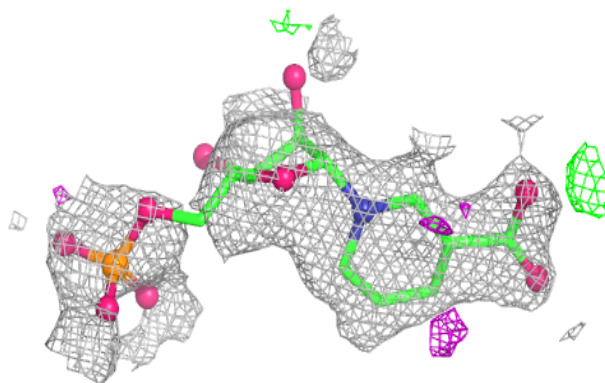


Electron density around NCN B 1302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

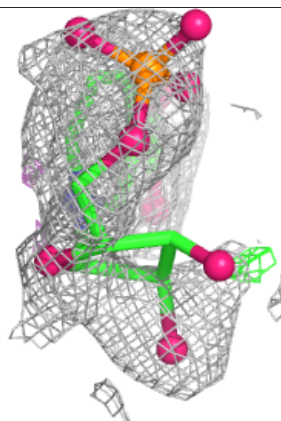
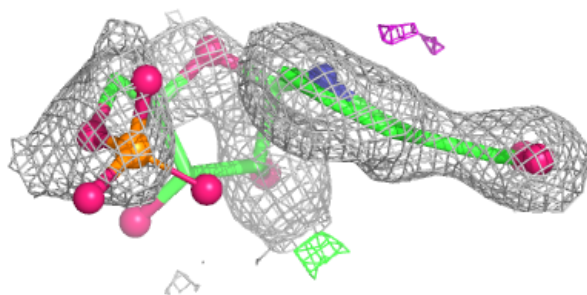
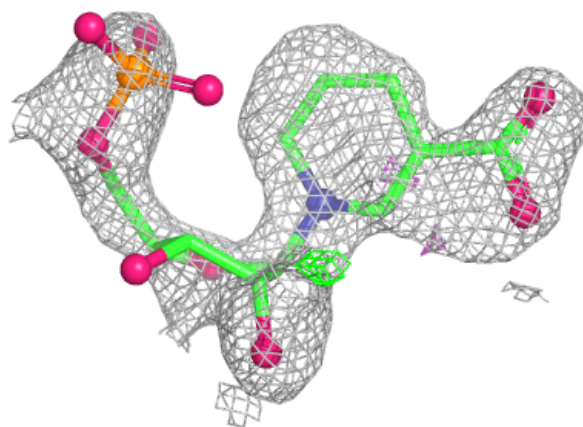
**Electron density around NCN A 302:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around NCN C 2302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.