



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:10 AM UTC

PDB ID : 8AT1 / pdb_00008at1
Title : CRYSTAL STRUCTURES OF ASPARTATE CARBAMOYLTRANSFERASE LIGATED WITH PHOSPHONOACETAMIDE, MALONATE, AND CTP OR ATP AT 2.8-ANGSTROMS RESOLUTION AND NEUTRAL P*H
Authors : Gouaux, J.E.; Stevens, R.C.; Lipscomb, W.N.
Deposited on : 1989-09-22
Resolution : 2.80 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)
Xtrriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

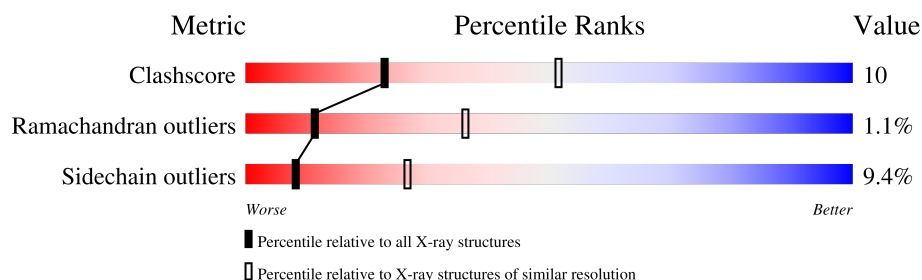
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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))
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	310	61% 31% 6% •
1	C	310	64% 31% 5% •
2	B	153	44% 39% 11% • 5%
2	D	153	47% 39% 8% • 5%
3	E	2	50% 50%
3	F	2	50% 50%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GLC	E	2	X	X	-	-
3	GLC	F	2	X	X	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7196 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 ASPARTATE CARBAMOYLTRANSFERASE (R STATE), CATALYTIC CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	310	Total	C	N	O	S	0	0	0
			2415	1527	423	456	9			
1	C	310	Total	C	N	O	S	0	0	0
			2415	1527	423	456	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	GLN	GLU	conflict	UNP P0A786
A	147	GLN	GLU	conflict	UNP P0A786
A	149	GLU	GLN	conflict	UNP P0A786
A	196	GLU	GLN	conflict	UNP P0A786
C	60	GLN	GLU	conflict	UNP P0A786
C	147	GLN	GLU	conflict	UNP P0A786
C	149	GLU	GLN	conflict	UNP P0A786
C	196	GLU	GLN	conflict	UNP P0A786

- Molecule 2 is a protein called ASPARTATE CARBAMOYLTRANSFERASE REGULATORY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1138	714	201	218	5			
2	D	146	Total	C	N	O	S	0	0	0
			1138	714	201	218	5			

There are 2 discrepancies between the modelled and reference sequences:

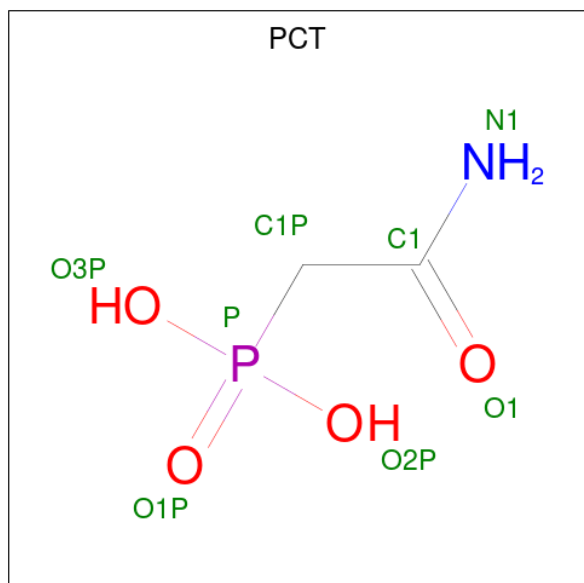
Chain	Residue	Modelled	Actual	Comment	Reference
B	8	GLY	GLN	conflict	UNP P0A7F3
D	8	GLY	GLN	conflict	UNP P0A7F3

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	E	2	Total	C	O	0	0	1
			7	4	3			
3	F	2	Total	C	O	0	0	1
			7	4	3			

- Molecule 4 is PHOSPHONOACETAMIDE (CCD ID: PCT) (formula: C₂H₆NO₄P).

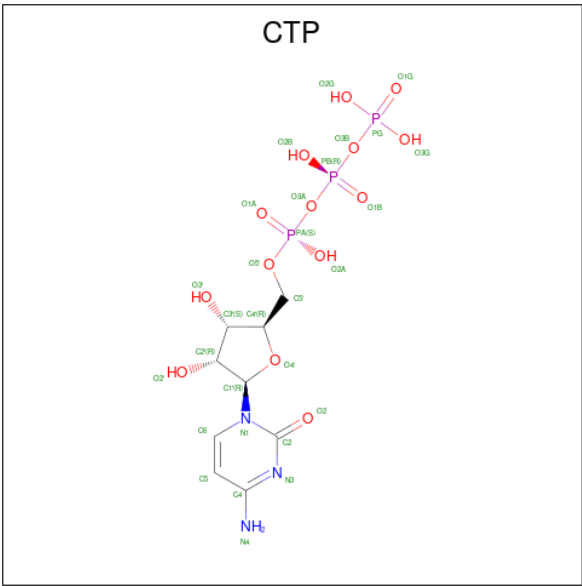


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			8	2	1	4	1		
4	C	1	Total	C	N	O	P	0	0
			8	2	1	4	1		

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Zn	0	0
			1	1		
5	D	1	Total	Zn	0	0
			1	1		

- Molecule 6 is CYTIDINE-5'-TRIPHOSPHATE (CCD ID: CTP) (formula: C₉H₁₆N₃O₁₄P₃).



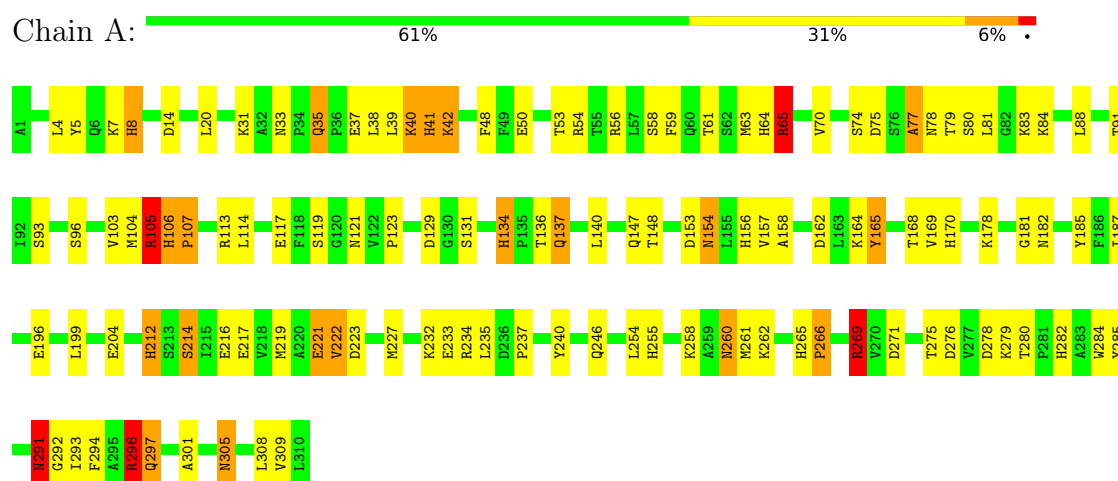
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	P	0	0
			29	9	3	14	3		
6	D	1	Total	C	N	O	P	0	0
			29	9	3	14	3		

3 Residue-property plots

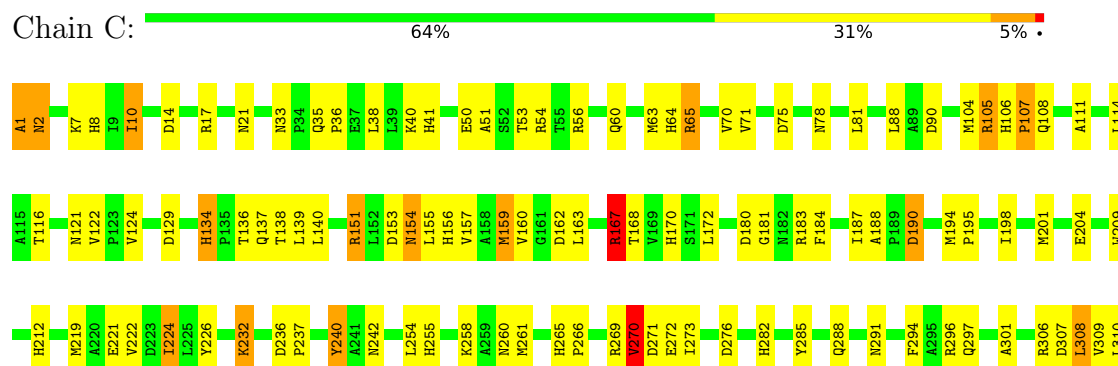
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

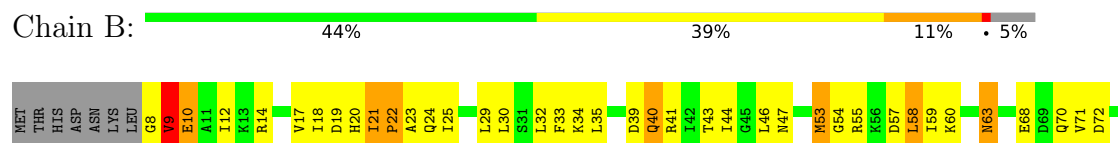
• Molecule 1: ASPARTATE CARBAMOYLTRANSFERASE (R STATE), CATALYTIC CHAIN

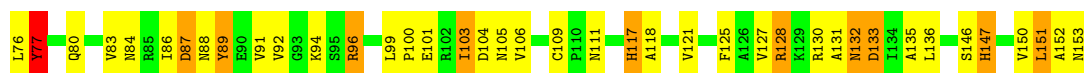


• Molecule 1: ASPARTATE CARBAMOYLTRANSFERASE (R STATE), CATALYTIC CHAIN



• Molecule 2: ASPARTATE CARBAMOYLTRANSFERASE REGULATORY CHAIN





- Molecule 2: ASPARTATE CARBAMOYLTRANSFERASE REGULATORY CHAIN



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	122.20Å 122.20Å 155.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.80	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.167 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7196	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PCT, GLC, CTP, ZN

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	1.10	19/2461 (0.8%)	1.95	74/3339 (2.2%)
1	C	1.14	16/2461 (0.7%)	1.93	71/3339 (2.1%)
2	B	1.00	5/1155 (0.4%)	1.92	29/1561 (1.9%)
2	D	0.98	5/1155 (0.4%)	1.82	27/1561 (1.7%)
All	All	1.08	45/7232 (0.6%)	1.92	201/9800 (2.1%)

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	3
1	C	0	1
2	B	0	2
2	D	0	1
All	All	0	7

All (45) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	64	HIS	CD2-NE2	-7.72	1.29	1.37
1	A	64	HIS	CD2-NE2	-7.37	1.29	1.37
1	C	170	HIS	CD2-NE2	-7.34	1.29	1.37
1	C	156	HIS	CD2-NE2	-7.24	1.29	1.37
2	D	117	HIS	CD2-NE2	-7.09	1.30	1.37
1	A	265	HIS	CD2-NE2	-7.06	1.30	1.37
1	A	41	HIS	CD2-NE2	-7.05	1.30	1.37
1	C	212	HIS	CD2-NE2	-6.87	1.30	1.37
1	C	222	VAL	CA-CB	6.81	1.64	1.54
2	B	117	HIS	CD2-NE2	-6.80	1.30	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	265	HIS	CD2-NE2	-6.69	1.30	1.37
1	A	134	HIS	CD2-NE2	-6.65	1.30	1.37
1	A	212	HIS	CD2-NE2	-6.65	1.30	1.37
1	A	265	HIS	CG-ND1	-6.44	1.31	1.38
1	C	41	HIS	CD2-NE2	-6.41	1.30	1.37
1	A	8	HIS	CD2-NE2	-6.41	1.30	1.37
2	D	147	HIS	CD2-NE2	-6.40	1.30	1.37
1	C	8	HIS	CD2-NE2	-6.31	1.30	1.37
1	A	106	HIS	CD2-NE2	-6.31	1.30	1.37
1	C	106	HIS	CD2-NE2	-6.29	1.30	1.37
1	A	156	HIS	CD2-NE2	-6.24	1.30	1.37
2	B	147	HIS	CD2-NE2	-6.22	1.31	1.37
1	A	282	HIS	CD2-NE2	-6.19	1.31	1.37
2	D	20	HIS	CD2-NE2	-6.19	1.31	1.37
1	C	134	HIS	CD2-NE2	-6.15	1.31	1.37
1	C	8	HIS	CG-ND1	-6.09	1.31	1.38
1	A	8	HIS	CG-ND1	-6.02	1.31	1.38
1	A	64	HIS	CG-ND1	-5.96	1.31	1.38
2	D	117	HIS	CG-ND1	-5.82	1.31	1.38
1	A	41	HIS	CG-ND1	-5.79	1.31	1.38
1	A	282	HIS	CG-ND1	-5.76	1.31	1.38
1	A	255	HIS	CD2-NE2	-5.74	1.31	1.37
1	A	170	HIS	CD2-NE2	-5.61	1.31	1.37
1	C	212	HIS	CG-ND1	-5.56	1.32	1.38
1	C	282	HIS	CD2-NE2	-5.55	1.31	1.37
1	A	106	HIS	CG-ND1	-5.50	1.32	1.38
1	A	212	HIS	CG-ND1	-5.39	1.32	1.38
2	D	20	HIS	CG-ND1	-5.36	1.32	1.38
1	C	255	HIS	CD2-NE2	-5.35	1.31	1.37
2	B	106	VAL	CA-CB	5.32	1.61	1.54
1	C	106	HIS	CG-ND1	-5.32	1.32	1.38
2	B	20	HIS	CD2-NE2	-5.26	1.32	1.37
1	A	134	HIS	CG-ND1	-5.21	1.32	1.38
1	C	170	HIS	CG-ND1	-5.19	1.32	1.38
2	B	20	HIS	CG-ND1	-5.09	1.32	1.38

All (201) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	309	VAL	O-C-N	13.70	137.10	122.45
1	C	105	ARG	NE-CZ-NH2	9.67	127.91	119.20
1	C	14	ASP	CA-CB-CG	9.22	121.82	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	41	HIS	ND1-CG-CD2	8.78	114.88	106.10
1	C	75	ASP	CA-CB-CG	8.73	121.33	112.60
1	A	105	ARG	NE-CZ-NH2	8.69	127.02	119.20
1	A	269	ARG	NE-CZ-NH2	8.68	127.01	119.20
1	C	10	ILE	N-CA-C	8.55	118.91	111.56
1	A	222	VAL	N-CA-C	8.51	119.32	110.05
1	A	106	HIS	ND1-CG-CD2	8.44	114.54	106.10
1	C	105	ARG	NE-CZ-NH1	-8.39	113.11	121.50
1	C	269	ARG	CA-CB-CG	8.38	130.87	114.10
2	B	133	ASP	CA-CB-CG	8.27	120.87	112.60
1	C	297	GLN	OE1-CD-NE2	-8.23	114.37	122.60
2	B	118	ALA	N-CA-C	8.02	120.02	111.28
2	B	153	ASN	CA-CB-CG	7.97	120.57	112.60
2	B	47	ASN	CA-CB-CG	7.89	120.49	112.60
2	B	70	GLN	OE1-CD-NE2	-7.80	114.80	122.60
1	A	276	ASP	CA-CB-CG	7.76	120.36	112.60
1	A	282	HIS	ND1-CG-CD2	7.71	113.81	106.10
1	A	75	ASP	CA-CB-CG	7.69	120.29	112.60
2	B	96	ARG	CA-CB-CG	-7.69	98.72	114.10
2	B	152	ALA	N-CA-C	7.49	126.76	110.80
1	C	242	ASN	N-CA-C	7.42	122.51	113.17
2	D	147	HIS	ND1-CG-CD2	7.41	113.51	106.10
2	B	63	ASN	CA-CB-CG	7.39	119.99	112.60
1	C	154	ASN	CB-CG-ND2	7.37	127.45	116.40
1	C	51	ALA	N-CA-C	7.34	120.63	110.35
2	B	53	MET	N-CA-C	-7.26	99.47	110.14
1	C	129	ASP	CA-CB-CG	7.10	119.70	112.60
1	C	60	GLN	CA-CB-CG	-7.08	99.94	114.10
2	B	101	GLU	CA-CB-CG	7.02	128.13	114.10
1	A	309	VAL	CA-C-N	6.97	134.24	121.70
1	A	309	VAL	C-N-CA	6.97	134.24	121.70
1	C	276	ASP	CA-CB-CG	6.95	119.55	112.60
1	C	265	HIS	CA-C-N	6.85	127.13	119.32
1	C	265	HIS	C-N-CA	6.85	127.13	119.32
1	A	269	ARG	NE-CZ-NH1	-6.84	114.66	121.50
1	C	151	ARG	NE-CZ-NH2	6.84	125.35	119.20
1	C	282	HIS	ND1-CG-CD2	6.75	112.85	106.10
1	C	8	HIS	ND1-CG-CD2	6.72	112.82	106.10
1	A	77	ALA	N-CA-C	6.68	119.39	111.71
1	C	154	ASN	OD1-CG-ND2	-6.66	115.94	122.60
2	B	47	ASN	OD1-CG-ND2	-6.66	115.94	122.60
1	A	297	GLN	OE1-CD-NE2	-6.65	115.95	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	104	ASP	CA-CB-CG	6.65	119.25	112.60
1	A	8	HIS	ND1-CG-CD2	6.65	112.75	106.10
2	B	33	PHE	N-CA-C	-6.64	98.42	109.24
1	A	33	ASN	CA-C-N	6.60	126.62	119.89
1	A	33	ASN	C-N-CA	6.60	126.62	119.89
1	C	53	THR	N-CA-C	-6.59	103.10	112.13
1	C	190	ASP	CA-CB-CG	6.58	119.18	112.60
1	C	291	ASN	OD1-CG-ND2	-6.58	116.02	122.60
2	B	21	ILE	N-CA-C	-6.52	103.02	108.63
1	C	134	HIS	ND1-CG-CD2	6.52	112.62	106.10
2	D	78	ALA	CA-C-N	6.49	126.17	119.56
2	D	78	ALA	C-N-CA	6.49	126.17	119.56
2	B	21	ILE	N-CA-CB	6.48	117.17	110.23
1	A	14	ASP	CA-CB-CG	6.47	119.07	112.60
1	A	35	GLN	CA-C-N	6.39	126.31	119.28
1	A	35	GLN	C-N-CA	6.39	126.31	119.28
1	C	54	ARG	CA-C-O	-6.38	114.33	120.90
2	B	18	ILE	N-CA-C	-6.37	99.12	108.54
1	C	10	ILE	N-CA-CB	-6.36	104.41	111.41
1	A	56	ARG	NE-CZ-NH2	6.32	124.88	119.20
1	A	106	HIS	CG-CD2-NE2	-6.29	100.91	107.20
2	B	70	GLN	CG-CD-NE2	6.28	125.82	116.40
1	C	260	ASN	OD1-CG-ND2	-6.27	116.33	122.60
1	A	37	GLU	N-CA-C	-6.24	104.83	112.88
1	C	282	HIS	CG-CD2-NE2	-6.22	100.98	107.20
1	A	129	ASP	CA-CB-CG	6.20	118.80	112.60
1	C	106	HIS	ND1-CG-CD2	6.18	112.28	106.10
1	A	79	THR	N-CA-C	-6.17	101.65	110.59
1	A	305	ASN	CA-CB-CG	-6.16	106.44	112.60
2	D	152	ALA	CA-C-O	6.16	127.95	120.62
1	C	138	THR	CA-CB-OG1	-6.16	100.37	109.60
1	A	282	HIS	N-CA-C	6.15	118.96	111.82
1	A	232	LYS	N-CA-CB	-6.15	99.76	109.78
1	A	296	ARG	NE-CZ-NH2	6.08	124.67	119.20
1	A	53	THR	N-CA-C	-6.03	102.35	111.34
2	D	153	ASN	OD1-CG-ND2	-6.03	116.57	122.60
1	A	182	ASN	OD1-CG-ND2	-6.03	116.57	122.60
1	C	124	VAL	N-CA-C	-6.03	99.73	108.17
2	B	77	TYR	N-CA-C	-6.02	107.76	114.62
1	A	169	VAL	CA-CB-CG1	-5.95	100.28	110.40
1	A	134	HIS	ND1-CG-CD2	5.95	112.05	106.10
1	C	294	PHE	CA-CB-CG	-5.93	107.86	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	102	ARG	NE-CZ-NH2	5.91	124.52	119.20
1	A	162	ASP	CA-CB-CG	5.90	118.50	112.60
2	D	14	ARG	NE-CZ-NH2	5.88	124.50	119.20
2	B	9	VAL	CA-C-O	-5.88	113.43	120.78
1	A	221	GLU	N-CA-CB	-5.87	101.53	110.33
2	D	88	ASN	CA-CB-CG	-5.86	106.75	112.60
1	C	159	MET	CA-CB-CG	-5.83	102.44	114.10
2	D	18	ILE	N-CA-C	-5.83	100.03	108.36
1	A	222	VAL	CB-CA-C	-5.80	105.21	111.59
2	B	151	LEU	CA-C-N	5.79	132.60	121.54
2	B	151	LEU	C-N-CA	5.79	132.60	121.54
2	D	9	VAL	O-C-N	5.79	129.80	122.57
1	A	93	SER	CA-C-O	-5.75	114.75	120.90
1	A	107	PRO	N-CA-C	5.75	121.32	113.84
1	C	35	GLN	O-C-N	-5.74	116.20	121.37
2	D	96	ARG	N-CA-C	-5.74	101.65	110.32
1	C	56	ARG	NE-CZ-NH2	5.73	124.36	119.20
1	C	270	VAL	O-C-N	-5.72	115.42	122.57
2	D	87	ASP	CA-CB-CG	-5.71	106.89	112.60
2	D	9	VAL	CA-C-N	5.70	132.43	121.54
2	D	9	VAL	C-N-CA	5.70	132.43	121.54
2	D	128	ARG	CB-CG-CD	5.70	124.42	111.30
2	D	74	LEU	N-CA-C	-5.69	105.08	111.28
1	C	78	ASN	N-CA-CB	-5.67	102.08	110.53
2	B	151	LEU	N-CA-CB	-5.67	102.74	110.56
1	C	122	VAL	CA-C-N	5.66	125.95	119.83
1	C	122	VAL	C-N-CA	5.66	125.95	119.83
2	D	90	GLU	CA-CB-CG	5.65	125.39	114.10
1	C	291	ASN	CB-CG-ND2	5.64	124.87	116.40
1	A	282	HIS	CG-CD2-NE2	-5.63	101.57	107.20
1	C	107	PRO	N-CA-C	5.62	120.88	113.86
1	A	154	ASN	OD1-CG-ND2	-5.61	116.99	122.60
1	C	265	HIS	ND1-CG-CD2	5.61	111.71	106.10
2	D	105	ASN	N-CA-C	5.60	122.73	110.80
2	D	139	LYS	CB-CG-CD	-5.58	98.46	111.30
1	C	17	ARG	N-CA-C	-5.58	105.20	111.28
2	D	118	ALA	N-CA-C	5.57	118.28	111.82
1	C	272	GLU	N-CA-C	5.56	118.27	111.82
1	A	103	VAL	N-CA-C	-5.55	100.40	108.17
1	C	153	ASP	CA-CB-CG	5.55	118.15	112.60
1	A	104	MET	CG-SD-CE	-5.54	88.70	100.90
1	C	260	ASN	CB-CG-ND2	5.54	124.71	116.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	255	HIS	CA-CB-CG	5.52	119.32	113.80
1	A	137	GLN	N-CA-C	-5.52	105.17	111.07
1	C	272	GLU	CA-CB-CG	5.51	125.12	114.10
1	C	134	HIS	CB-CG-CD2	-5.50	124.06	131.20
1	C	104	MET	N-CA-C	5.48	118.16	109.50
2	D	41	ARG	NE-CZ-NH2	5.47	124.12	119.20
1	C	17	ARG	NE-CZ-NH2	5.47	124.12	119.20
1	A	234	ARG	NE-CZ-NH2	5.45	124.11	119.20
1	A	106	HIS	CA-C-N	5.43	125.10	119.56
1	A	106	HIS	C-N-CA	5.43	125.10	119.56
1	C	167	ARG	CB-CG-CD	5.43	123.79	111.30
1	C	309	VAL	N-CA-C	-5.43	97.79	107.18
1	C	232	LYS	CA-CB-CG	5.41	124.93	114.10
1	A	81	LEU	N-CA-C	-5.39	105.30	111.07
1	C	269	ARG	NE-CZ-NH2	5.39	124.05	119.20
2	B	40	GLN	OE1-CD-NE2	-5.39	117.21	122.60
1	C	2	ASN	N-CA-C	-5.39	103.34	110.40
1	A	153	ASP	CA-CB-CG	5.38	117.98	112.60
1	C	139	LEU	CA-C-O	-5.37	115.36	121.00
2	B	10	GLU	N-CA-C	5.36	116.99	108.79
1	C	198	ILE	N-CA-C	-5.36	105.49	110.53
1	C	183	ARG	NE-CZ-NH2	5.36	124.02	119.20
1	A	260	ASN	OD1-CG-ND2	-5.34	117.26	122.60
1	C	162	ASP	CA-CB-CG	5.32	117.92	112.60
1	A	121	ASN	OD1-CG-ND2	-5.32	117.28	122.60
2	B	21	ILE	CB-CA-C	-5.31	106.00	110.94
1	C	54	ARG	N-CA-C	5.30	116.91	111.03
2	B	19	ASP	CA-CB-CG	5.29	117.89	112.60
2	D	106	VAL	CB-CA-C	-5.29	105.66	110.63
1	A	53	THR	OG1-CB-CG2	5.28	119.86	109.30
2	B	39	ASP	CA-C-O	5.28	125.90	119.78
1	A	58	SER	N-CA-C	-5.27	105.45	111.14
1	A	113	ARG	NE-CZ-NH2	5.25	123.93	119.20
2	D	147	HIS	CG-CD2-NE2	-5.25	101.95	107.20
2	D	147	HIS	CB-CG-CD2	-5.25	124.37	131.20
2	B	87	ASP	N-CA-C	-5.25	99.89	108.76
1	A	106	HIS	CB-CG-CD2	-5.23	124.40	131.20
1	A	265	HIS	ND1-CG-CD2	5.23	111.33	106.10
1	C	236	ASP	CA-CB-CG	-5.21	107.39	112.60
1	A	168	THR	N-CA-C	5.21	116.65	111.07
2	D	10	GLU	CB-CG-CD	5.21	121.45	112.60
1	C	1	ALA	N-CA-C	-5.20	96.43	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	91	THR	N-CA-C	-5.20	105.50	111.07
1	A	41	HIS	CG-CD2-NE2	-5.20	102.00	107.20
1	C	33	ASN	CA-C-N	5.19	125.13	119.78
1	C	33	ASN	C-N-CA	5.19	125.13	119.78
1	A	121	ASN	CB-CG-ND2	5.19	124.18	116.40
1	A	157	VAL	CA-C-O	5.18	125.39	120.31
1	C	255	HIS	N-CA-C	5.18	117.67	111.71
1	C	71	VAL	O-C-N	-5.18	117.56	123.10
1	A	284	TRP	CE2-CD2-CG	-5.16	101.00	107.20
1	A	291	ASN	OD1-CG-ND2	-5.16	117.44	122.60
1	C	90	ASP	CA-CB-CG	5.16	117.76	112.60
2	B	8	GLY	N-CA-C	-5.16	98.34	113.30
1	C	288	GLN	OE1-CD-NE2	-5.15	117.45	122.60
2	D	105	ASN	CA-CB-CG	5.14	117.74	112.60
1	A	84	LYS	CG-CD-CE	5.13	123.11	111.30
1	A	78	ASN	OD1-CG-ND2	-5.13	117.47	122.60
1	A	271	ASP	CA-CB-CG	5.13	117.73	112.60
1	A	53	THR	CA-CB-OG1	-5.12	101.93	109.60
1	A	232	LYS	CA-CB-CG	5.09	124.28	114.10
1	A	80	SER	N-CA-C	-5.09	106.58	112.89
1	A	65	ARG	NE-CZ-NH2	5.09	123.78	119.20
1	C	106	HIS	CA-C-N	5.09	124.70	119.56
1	C	106	HIS	C-N-CA	5.09	124.70	119.56
1	A	266	PRO	O-C-N	5.08	128.77	122.22
1	A	104	MET	N-CA-CB	-5.06	102.53	110.77
1	A	157	VAL	N-CA-C	5.05	115.72	107.24
1	A	219	MET	N-CA-C	5.04	117.16	111.11
2	D	40	GLN	N-CA-C	-5.02	103.09	110.52
1	A	204	GLU	O-C-N	-5.01	116.87	122.09
1	C	10	ILE	CB-CG1-CD1	-5.01	103.27	113.80

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	165	TYR	Sidechain
1	A	185	TYR	Sidechain
1	A	5	TYR	Sidechain
2	B	77	TYR	Sidechain
2	B	89	TYR	Sidechain
1	C	240	TYR	Sidechain
2	D	89	TYR	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	2415	0	2422	43	0
1	C	2415	0	2422	34	0
2	B	1138	0	1154	37	0
2	D	1138	0	1154	39	0
3	E	7	0	0	0	0
3	F	7	0	0	0	0
4	A	8	0	4	1	0
4	C	8	0	4	0	0
5	B	1	0	0	0	0
5	D	1	0	0	0	0
6	B	29	0	12	3	0
6	D	29	0	12	2	0
All	All	7196	0	7184	149	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 10.

All (149) 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:C:10:ILE:HD11	1:C:116:THR:HG21	1.43	0.98
1:C:137:GLN:HG2	1:C:168:THR:HG22	1.71	0.72
2:B:14:ARG:HG3	2:B:87:ASP:HA	1.70	0.72
2:D:30:LEU:HD12	2:D:36:THR:HG23	1.70	0.71
1:C:301:ALA:HB1	1:C:308:LEU:HD11	1.73	0.71
2:B:71:VAL:HG13	2:B:83:VAL:HG21	1.71	0.71
2:B:80:GLN:HA	2:B:96:ARG:HH22	1.55	0.70
1:A:292:GLY:O	1:A:296:ARG:HB2	1.91	0.69
2:D:130:ARG:HD2	2:D:135:ALA:HB2	1.74	0.67
1:A:154:ASN:HA	1:A:181:GLY:O	1.94	0.67
1:A:235:LEU:HB2	1:A:240:TYR:HE2	1.61	0.65
2:B:40:GLN:NE2	2:B:63:ASN:HB2	2.12	0.65
2:D:94:LYS:HE2	6:D:155:CTP:O2A	1.97	0.65
2:D:72:ASP:HB3	2:D:100:PRO:HG3	1.77	0.64
2:D:47:ASN:HD21	2:D:55:ARG:HE	1.44	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1:ALA:HA	1:C:306:ARG:HG2	1.80	0.63
2:D:11:ALA:HA	6:D:155:CTP:N3	2.15	0.62
2:B:133:ASP:HB2	2:B:147:HIS:CE1	2.36	0.61
2:D:146:SER:HB3	2:D:149:VAL:HG23	1.83	0.59
1:A:4:LEU:HD23	1:A:7:LYS:HD3	1.85	0.59
2:B:103:ILE:HD11	2:B:127:VAL:HG22	1.83	0.59
1:C:134:HIS:HB2	1:C:167:ARG:HD3	1.84	0.58
2:D:49:PRO:O	2:D:56:LYS:HG2	2.03	0.58
2:B:22:PRO:HB2	2:B:25:ILE:HD12	1.86	0.58
1:A:258:LYS:HG2	1:A:260:ASN:OD1	2.04	0.57
1:A:196:GLU:HA	1:A:199:LEU:HD12	1.85	0.57
1:C:157:VAL:HG13	1:C:224:ILE:HG22	1.88	0.56
1:A:136:THR:HB	1:A:296:ARG:HH21	1.71	0.56
1:C:154:ASN:HA	1:C:181:GLY:O	2.07	0.55
2:D:102:ARG:HA	2:D:125:PHE:O	2.07	0.55
1:A:20:LEU:HD11	1:A:178:LYS:HD2	1.89	0.55
1:C:136:THR:HB	1:C:296:ARG:NH2	2.21	0.55
2:B:91:VAL:HG11	2:B:94:LYS:HZ2	1.72	0.55
1:C:10:ILE:HD11	1:C:116:THR:CG2	2.29	0.55
1:A:137:GLN:O	1:A:140:LEU:HG	2.06	0.54
1:C:159:MET:HE3	1:C:172:LEU:HD23	1.88	0.54
1:C:159:MET:CE	1:C:172:LEU:HD23	2.38	0.54
1:A:254:LEU:HD12	1:A:280:THR:HG21	1.90	0.54
1:C:184:PHE:O	1:C:209:TRP:HA	2.08	0.54
1:C:88:LEU:HD23	1:C:114:LEU:HD12	1.90	0.53
2:B:136:LEU:HD12	2:B:150:VAL:HG21	1.89	0.53
2:D:23:ALA:HB2	2:D:55:ARG:HG3	1.89	0.53
1:A:237:PRO:HA	1:A:240:TYR:CD1	2.44	0.53
2:B:109:CYS:HA	2:B:125:PHE:HZ	1.73	0.53
2:B:128:ARG:HB3	2:B:135:ALA:HB3	1.91	0.53
1:A:164:LYS:HD3	1:A:165:TYR:CE2	2.44	0.53
2:B:23:ALA:HB3	2:B:55:ARG:HH21	1.73	0.52
2:D:14:ARG:HG2	2:D:63:ASN:HA	1.92	0.52
1:A:96:SER:OG	1:A:119:SER:HA	2.10	0.52
2:B:53:MET:SD	2:B:54:GLY:N	2.83	0.52
2:D:13:LYS:CG	2:D:88:ASN:HA	2.40	0.52
2:D:49:PRO:HA	2:D:54:GLY:O	2.09	0.52
2:D:106:VAL:HG23	2:D:107:LEU:HD13	1.91	0.52
2:B:133:ASP:HB2	2:B:147:HIS:HE1	1.75	0.52
2:D:50:SER:HB3	2:D:56:LYS:HD3	1.92	0.52
2:B:30:LEU:HA	2:B:35:LEU:HD12	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:102:ARG:HH21	2:D:102:ARG:HB3	1.75	0.51
1:A:88:LEU:HD23	1:A:114:LEU:HD12	1.91	0.51
2:B:30:LEU:HD13	2:B:59:ILE:HG12	1.92	0.51
1:A:223:ASP:O	1:A:261:MET:HA	2.10	0.51
2:D:43:THR:HB	2:D:60:LYS:HB2	1.91	0.51
1:C:226:TYR:OH	1:C:266:PRO:HG3	2.11	0.51
2:D:111:ASN:O	2:D:117:HIS:HE1	1.92	0.51
1:C:36:PRO:HA	1:C:65:ARG:O	2.11	0.51
1:A:54:ARG:HE	4:A:311:PCT:H1P2	1.75	0.51
1:C:163:LEU:HG	1:C:188:ALA:HB2	1.94	0.50
2:B:23:ALA:O	2:B:24:GLN:HB2	2.11	0.50
1:A:63:MET:SD	1:A:70:VAL:HG22	2.53	0.49
2:D:13:LYS:HG2	2:D:88:ASN:HA	1.94	0.49
2:B:17:VAL:HG13	2:B:60:LYS:HG2	1.93	0.49
1:A:254:LEU:HD22	1:A:261:MET:HE1	1.94	0.49
1:A:293:ILE:O	1:A:297:GLN:HG2	2.12	0.49
2:D:102:ARG:HG3	2:D:125:PHE:O	2.13	0.49
2:B:46:LEU:HB2	2:D:42:ILE:HD13	1.93	0.49
2:D:12:ILE:HG22	2:D:41:ARG:NH1	2.28	0.48
1:C:50:GLU:HB2	1:C:107:PRO:HD3	1.95	0.48
2:D:14:ARG:HA	2:D:86:ILE:O	2.12	0.48
1:A:8:HIS:ND1	1:A:123:PRO:HA	2.29	0.48
2:B:99:LEU:HD22	2:B:127:VAL:HG11	1.94	0.48
2:D:69:ASP:O	2:D:73:GLN:HG3	2.13	0.48
1:C:254:LEU:CD2	1:C:261:MET:HE1	2.43	0.48
2:D:84:ASN:HD21	2:D:94:LYS:HD3	1.78	0.48
2:B:111:ASN:O	2:B:117:HIS:HE1	1.96	0.47
1:A:158:ALA:HB2	1:A:222:VAL:HG11	1.96	0.47
2:B:111:ASN:O	2:B:117:HIS:CE1	2.68	0.47
1:C:254:LEU:HD22	1:C:261:MET:HE1	1.97	0.47
2:D:48:LEU:O	2:D:55:ARG:HA	2.15	0.47
1:A:77:ALA:O	1:A:83:LYS:HG2	2.14	0.47
1:A:216:GLU:H	1:A:216:GLU:CD	2.23	0.47
2:B:100:PRO:HG2	2:B:127:VAL:HG21	1.97	0.47
1:C:137:GLN:O	1:C:140:LEU:HG	2.15	0.47
1:C:111:ALA:O	1:C:114:LEU:HB3	2.15	0.46
1:A:40:LYS:O	1:A:41:HIS:HB2	2.15	0.46
2:D:130:ARG:CD	2:D:135:ALA:HB2	2.44	0.46
1:A:301:ALA:O	1:A:305:ASN:HB2	2.16	0.45
2:B:9:VAL:HG11	2:B:58:LEU:HD21	1.97	0.45
1:A:136:THR:HB	1:A:296:ARG:NH2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:LEU:O	1:A:42:LYS:HB2	2.17	0.45
2:B:29:LEU:HD21	2:B:77:TYR:HB2	1.97	0.45
2:B:25:ILE:O	2:B:29:LEU:HG	2.17	0.45
2:B:41:ARG:HD2	2:D:48:LEU:CD2	2.47	0.45
2:D:12:ILE:HG22	2:D:41:ARG:HH12	1.82	0.44
1:A:48:PHE:CD2	1:A:105:ARG:HB2	2.52	0.44
1:C:151:ARG:HG3	1:C:155:LEU:HD21	1.99	0.44
1:A:134:HIS:CD2	1:A:137:GLN:HB2	2.53	0.44
2:B:84:ASN:ND2	6:B:155:CTP:H5'2	2.33	0.44
1:C:237:PRO:HA	1:C:240:TYR:CD2	2.53	0.44
1:C:160:VAL:HA	1:C:187:ILE:O	2.18	0.44
2:D:30:LEU:HD13	2:D:35:LEU:HB2	2.00	0.44
2:D:42:ILE:HA	2:D:60:LYS:O	2.18	0.43
2:D:99:LEU:HB3	2:D:129:LYS:NZ	2.32	0.43
2:B:41:ARG:HD2	2:D:48:LEU:HD22	2.00	0.43
1:C:163:LEU:HD13	1:C:194:MET:HE2	2.00	0.43
2:D:10:GLU:HB3	2:D:60:LYS:HZ1	1.83	0.43
1:A:227:MET:O	1:A:266:PRO:HD2	2.18	0.43
1:A:61:THR:O	1:A:65:ARG:HG3	2.18	0.43
1:C:2:ASN:N	1:C:306:ARG:O	2.52	0.43
2:D:126:ALA:O	2:D:136:LEU:HA	2.19	0.43
2:D:10:GLU:HB3	2:D:60:LYS:NZ	2.34	0.43
1:A:31:LYS:HD2	1:A:294:PHE:CD2	2.54	0.43
1:A:114:LEU:CD2	2:B:121:VAL:HG11	2.49	0.43
6:B:155:CTP:H3'	6:B:155:CTP:O3A	2.19	0.43
1:C:194:MET:HA	1:C:195:PRO:HD2	1.91	0.43
1:A:59:PHE:CE1	1:A:296:ARG:HD2	2.54	0.42
2:B:44:ILE:HB	2:D:44:ILE:HD13	2.01	0.42
1:C:219:MET:SD	1:C:261:MET:HE2	2.59	0.42
2:B:21:ILE:HB	2:B:57:ASP:HB2	2.02	0.42
1:A:106:HIS:HA	1:A:107:PRO:HD2	1.79	0.42
1:C:201:MET:HE3	1:C:201:MET:HB3	1.83	0.42
2:D:19:ASP:OD2	2:D:58:LEU:HD13	2.19	0.42
1:C:273:ILE:HD13	1:C:273:ILE:HG21	1.80	0.42
1:A:187:ILE:HG12	1:A:212:HIS:HB2	2.02	0.41
1:C:237:PRO:HA	1:C:240:TYR:CE2	2.55	0.41
1:C:270:VAL:HG12	1:C:271:ASP:N	2.35	0.41
2:B:9:VAL:HG11	2:B:58:LEU:CD2	2.50	0.41
1:A:35:GLN:HB3	1:A:38:LEU:HD13	2.03	0.41
1:A:148:THR:HG21	1:A:262:LYS:HG3	2.03	0.41
2:B:89:TYR:HA	6:B:155:CTP:N4	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:147:HIS:O	2:B:151:LEU:HD12	2.20	0.41
1:A:214:SER:CB	1:A:217:GLU:HG3	2.51	0.41
1:A:269:ARG:NH1	1:A:278:ASP:OD2	2.54	0.41
1:A:275:THR:O	1:A:278:ASP:HB2	2.21	0.41
1:A:291:ASN:HD22	1:A:291:ASN:HA	1.67	0.41
1:C:63:MET:SD	1:C:70:VAL:HG22	2.61	0.40
2:D:84:ASN:ND2	2:D:94:LYS:HD3	2.36	0.40
2:B:128:ARG:NH2	2:B:130:ARG:HG3	2.36	0.40
1:C:7:LYS:HA	1:C:7:LYS:HD3	1.98	0.40
1:A:40:LYS:HA	1:A:40:LYS:HD3	1.95	0.40
2:B:84:ASN:HD21	2:B:91:VAL:CG1	2.33	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	308/310 (99%)	295 (96%)	13 (4%)	0	100	100
1	C	308/310 (99%)	296 (96%)	11 (4%)	1 (0%)	36	66
2	B	144/153 (94%)	125 (87%)	15 (10%)	4 (3%)	4	14
2	D	144/153 (94%)	126 (88%)	13 (9%)	5 (4%)	3	10
All	All	904/926 (98%)	842 (93%)	52 (6%)	10 (1%)	11	36

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	34	LYS
2	B	132	ASN
2	D	51	GLY
2	D	54	GLY

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Mol	Chain	Res	Type
2	D	105	ASN
2	D	50	SER
2	D	10	GLU
2	B	9	VAL
1	C	270	VAL
2	B	131	ALA

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	261/261 (100%)	242 (93%)	19 (7%)	13	38
1	C	261/261 (100%)	241 (92%)	20 (8%)	12	36
2	B	129/136 (95%)	112 (87%)	17 (13%)	4	14
2	D	129/136 (95%)	112 (87%)	17 (13%)	4	14
All	All	780/794 (98%)	707 (91%)	73 (9%)	8	27

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	LYS
1	A	42	LYS
1	A	50	GLU
1	A	65	ARG
1	A	74	SER
1	A	105	ARG
1	A	117	GLU
1	A	131	SER
1	A	147	GLN
1	A	214	SER
1	A	221	GLU
1	A	233	GLU
1	A	246	GLN
1	A	269	ARG
1	A	279	LYS

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Mol	Chain	Res	Type
1	A	285	TYR
1	A	291	ASN
1	A	296	ARG
1	A	308	LEU
2	B	10	GLU
2	B	12	ILE
2	B	22	PRO
2	B	32	LEU
2	B	43	THR
2	B	58	LEU
2	B	68	GLU
2	B	72	ASP
2	B	76	LEU
2	B	86	ILE
2	B	88	ASN
2	B	92	VAL
2	B	103	ILE
2	B	105	ASN
2	B	128	ARG
2	B	132	ASN
2	B	146	SER
1	C	21	ASN
1	C	38	LEU
1	C	40	LYS
1	C	65	ARG
1	C	81	LEU
1	C	105	ARG
1	C	108	GLN
1	C	121	ASN
1	C	167	ARG
1	C	180	ASP
1	C	190	ASP
1	C	204	GLU
1	C	221	GLU
1	C	224	ILE
1	C	232	LYS
1	C	258	LYS
1	C	285	TYR
1	C	307	ASP
1	C	308	LEU
1	C	310	LEU
2	D	34	LYS

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Mol	Chain	Res	Type
2	D	36	THR
2	D	38	THR
2	D	39	ASP
2	D	40	GLN
2	D	47	ASN
2	D	52	GLU
2	D	61	ILE
2	D	71	VAL
2	D	92	VAL
2	D	95	SER
2	D	101	GLU
2	D	102	ARG
2	D	105	ASN
2	D	107	LEU
2	D	137	LYS
2	D	153	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	60	GLN
1	A	108	GLN
1	A	146	GLN
1	A	170	HIS
1	A	246	GLN
1	A	305	ASN
2	B	40	GLN
2	B	47	ASN
2	B	84	ASN
2	B	105	ASN
2	B	113	ASN
2	B	117	HIS
2	B	153	ASN
1	C	13	ASN
1	C	21	ASN
1	C	64	HIS
1	C	108	GLN
1	C	146	GLN
1	C	154	ASN
1	C	255	HIS
2	D	63	ASN

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Mol	Chain	Res	Type
2	D	70	GLN
2	D	117	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 ⓘ

Of 4 monosaccharides modelled in this entry, 2 were used 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	GLC	E	2	3	5,5,12	4.11	3 (60%)	5,5,17	2.02	3 (60%)
3	GLC	F	2	3	5,5,12	3.90	3 (60%)	5,5,17	2.12	2 (40%)

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	GLC	E	2	3	2/2/1/5	1/3/3/22	-
3	GLC	F	2	3	2/2/1/5	1/3/3/22	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	2	GLC	C4-C3	-6.65	1.23	1.51
3	F	2	GLC	C4-C3	-6.45	1.24	1.51
3	E	2	GLC	O3-C3	-5.01	1.22	1.43
3	F	2	GLC	O3-C3	-4.67	1.24	1.43
3	E	2	GLC	O5-C1	-3.78	1.22	1.42
3	F	2	GLC	O5-C1	-3.58	1.23	1.42

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	2	GLC	O3-C3-C4	3.41	124.13	109.45
3	E	2	GLC	O3-C3-C4	2.63	120.77	109.45
3	F	2	GLC	O5-C1-C2	2.48	119.47	111.34
3	E	2	GLC	O5-C1-C2	2.47	119.44	111.34
3	E	2	GLC	O3-C3-C2	2.02	118.45	109.56

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	E	2	GLC	C1
3	E	2	GLC	C3
3	F	2	GLC	C1
3	F	2	GLC	C3

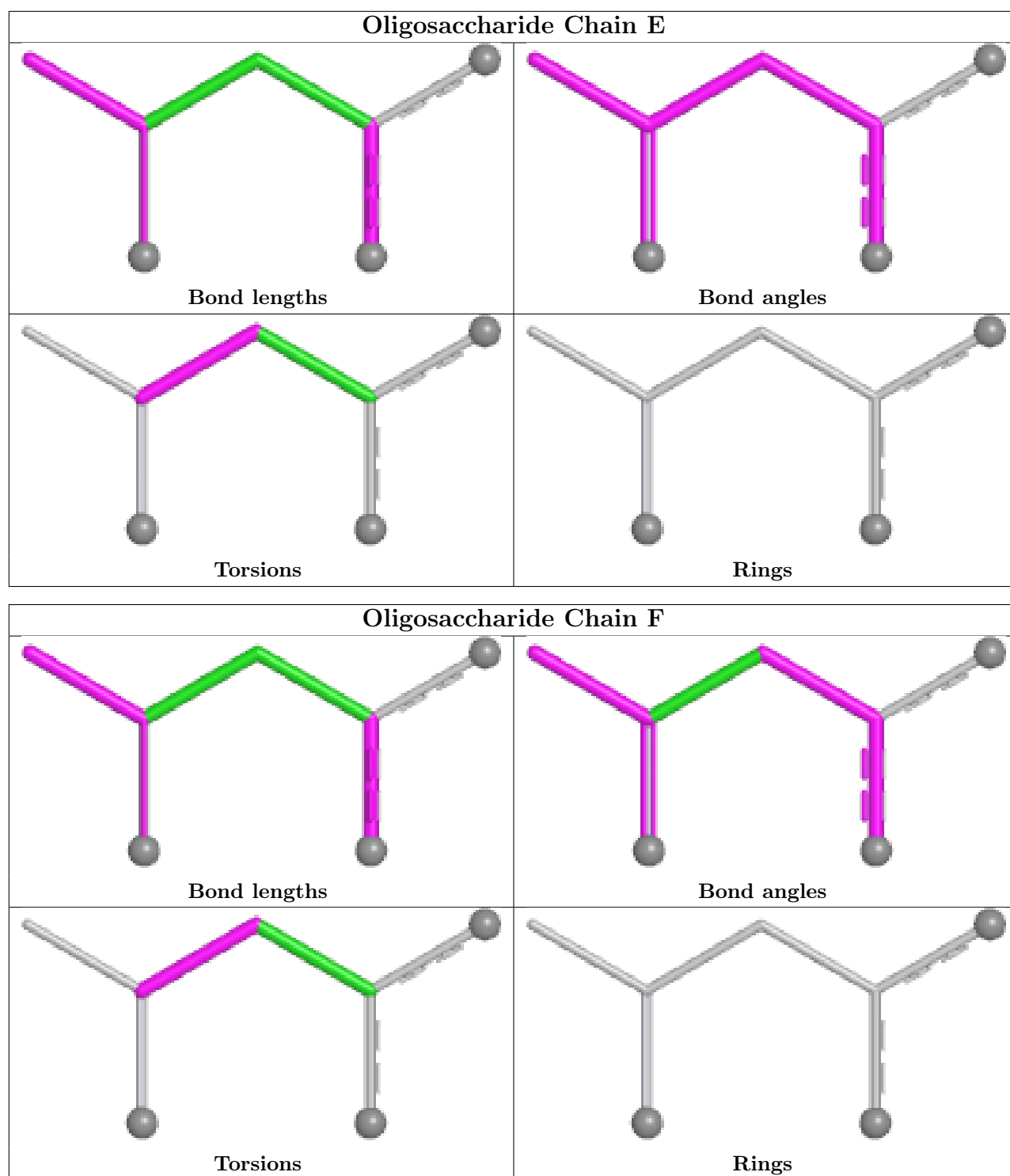
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	2	GLC	C1-C2-C3-C4
3	F	2	GLC	C1-C2-C3-C4

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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
6	CTP	D	155	-	29,30,30	2.47	9 (31%)	43,47,47	1.90	6 (13%)
6	CTP	B	155	-	29,30,30	2.20	8 (27%)	43,47,47	1.93	7 (16%)
4	PCT	C	311	-	7,7,7	1.67	2 (28%)	9,10,10	1.90	2 (22%)
4	PCT	A	311	-	7,7,7	1.83	2 (28%)	9,10,10	3.07	2 (22%)

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
6	CTP	D	155	-	-	4/22/38/38	0/2/2/2
6	CTP	B	155	-	-	7/22/38/38	0/2/2/2
4	PCT	C	311	-	-	3/4/5/5	-
4	PCT	A	311	-	-	3/4/5/5	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	155	CTP	PB-O3B	7.43	1.67	1.59
6	D	155	CTP	PA-O3A	6.29	1.66	1.59
6	B	155	CTP	PA-O3A	5.62	1.65	1.59
6	B	155	CTP	PB-O3B	5.54	1.65	1.59
6	D	155	CTP	PB-O3A	5.40	1.65	1.59
6	B	155	CTP	PB-O3A	4.93	1.64	1.59
4	A	311	PCT	P-O1P	3.54	1.57	1.50
4	C	311	PCT	P-O1P	3.19	1.56	1.50
6	B	155	CTP	C2-N1	-3.09	1.33	1.40
6	D	155	CTP	PG-O1G	3.00	1.59	1.50
6	D	155	CTP	C2-N1	-2.93	1.33	1.40
4	C	311	PCT	P-C1P	2.85	1.84	1.79
6	B	155	CTP	C6-N1	-2.75	1.31	1.38
4	A	311	PCT	P-C1P	2.72	1.84	1.79
6	B	155	CTP	PG-O1G	2.55	1.58	1.50
6	D	155	CTP	PB-O1B	2.43	1.59	1.50
6	D	155	CTP	C6-N1	-2.35	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	155	CTP	PB-O1B	2.18	1.58	1.50
6	D	155	CTP	PA-O1A	2.17	1.58	1.50
6	B	155	CTP	PA-O1A	2.17	1.58	1.50
6	D	155	CTP	C5-C4	-2.04	1.38	1.42

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	155	CTP	O2-C2-N3	-7.56	110.42	122.33
6	B	155	CTP	O2-C2-N3	-7.19	111.00	122.33
4	A	311	PCT	O1-C1-C1P	7.17	127.41	119.86
6	B	155	CTP	N1-C2-N3	6.18	129.53	118.80
6	D	155	CTP	N1-C2-N3	5.63	128.58	118.80
4	A	311	PCT	C1P-C1-N1	-4.90	109.55	115.41
4	C	311	PCT	O1-C1-C1P	4.28	124.37	119.86
6	B	155	CTP	C6-N1-C2	-3.66	114.28	120.46
6	D	155	CTP	C6-N1-C2	-3.65	114.30	120.46
6	B	155	CTP	C4-N3-C2	-3.47	114.80	120.26
6	D	155	CTP	C4-N3-C2	-3.07	115.44	120.26
6	D	155	CTP	C1'-N1-C2	2.99	125.04	118.44
4	C	311	PCT	O1P-P-C1P	-2.47	105.39	111.10
6	B	155	CTP	C3'-C2'-C1'	2.34	105.88	101.46
6	D	155	CTP	C3'-C2'-C1'	2.17	105.56	101.46
6	B	155	CTP	C2'-C1'-N1	-2.10	107.39	113.25
6	B	155	CTP	O4'-C1'-N1	2.01	112.92	108.36

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	311	PCT	C1-C1P-P-O1P
4	A	311	PCT	C1-C1P-P-O3P
4	C	311	PCT	C1-C1P-P-O1P
4	C	311	PCT	C1-C1P-P-O2P
4	C	311	PCT	C1-C1P-P-O3P
6	B	155	CTP	C5'-O5'-PA-O1A
6	B	155	CTP	C5'-O5'-PA-O2A
4	A	311	PCT	C1-C1P-P-O2P
6	D	155	CTP	C3'-C4'-C5'-O5'
6	D	155	CTP	O4'-C4'-C5'-O5'
6	B	155	CTP	O4'-C4'-C5'-O5'
6	B	155	CTP	C4'-C5'-O5'-PA

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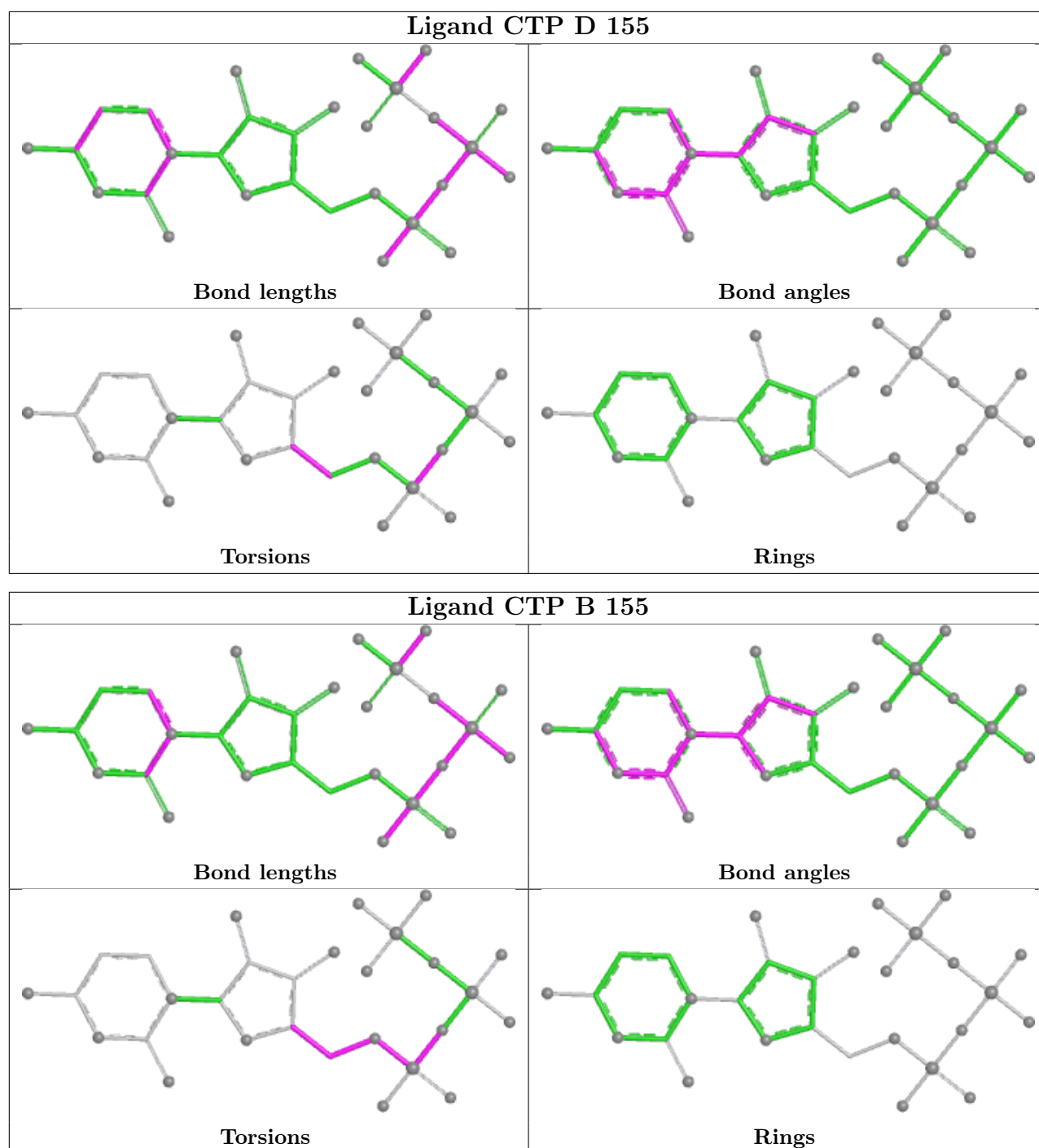
Mol	Chain	Res	Type	Atoms
6	B	155	CTP	PB-O3A-PA-O1A
6	D	155	CTP	PB-O3A-PA-O2A
6	B	155	CTP	C5'-O5'-PA-O3A
6	B	155	CTP	C3'-C4'-C5'-O5'
6	D	155	CTP	PB-O3A-PA-O1A

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	155	CTP	2	0
6	B	155	CTP	3	0
4	A	311	PCT	1	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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.