



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2017 – 10:53 AM EST

PDB ID : 8AT1
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 : unknown
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

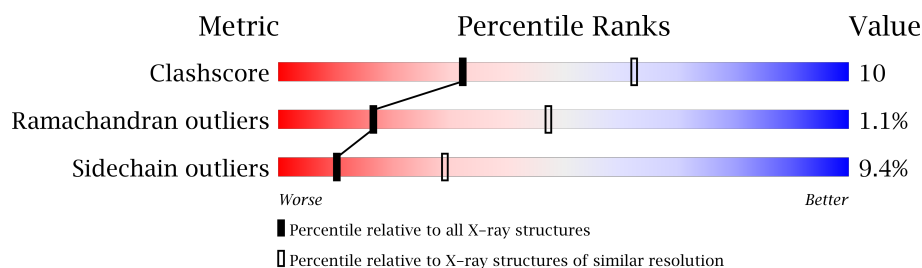
1 Overall quality at a glance

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	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	310	
1	C	310	
2	B	153	
2	D	153	

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
3	MAL	A	312	X	X	-	-
3	MAL	C	312	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

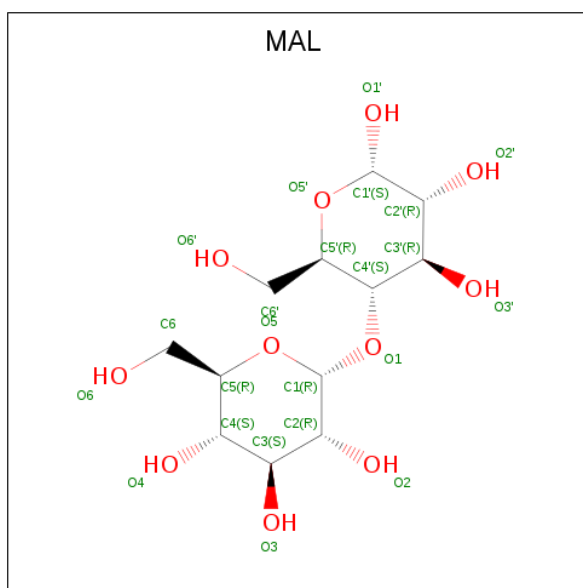
- Molecule 2 is a protein called ASPARTATE CARBAMOYLTRANSFERASE REGULATORY CHAIN.

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:

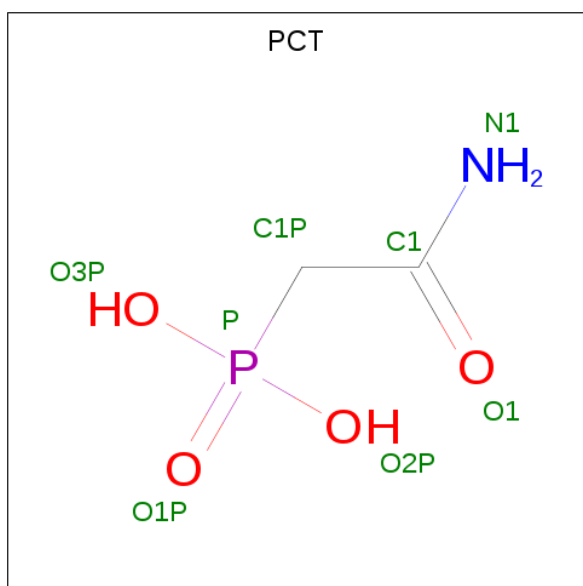
Chain	Residue	Modelled	Actual	Comment	Reference
B	8	GLY	GLN	CONFLICT	UNP P0A7F3
D	8	GLY	GLN	CONFLICT	UNP P0A7F3

- Molecule 3 is MALTOSE (three-letter code: MAL) (formula: $C_{12}H_{22}O_{11}$).



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

- Molecule 4 is PHOSPHONOACETAMIDE (three-letter code: PCT) (formula: $C_2H_6NO_4P$).



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

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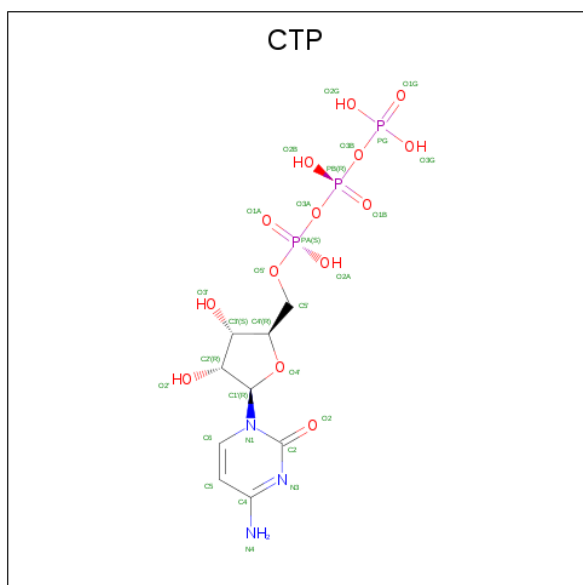
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total	C	N	O	P	0	0
			8	2	1	4	1		

- Molecule 5 is ZINC ION (three-letter code: 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 (three-letter code: CTP) (formula: $C_9H_{16}N_3O_{14}P_3$).



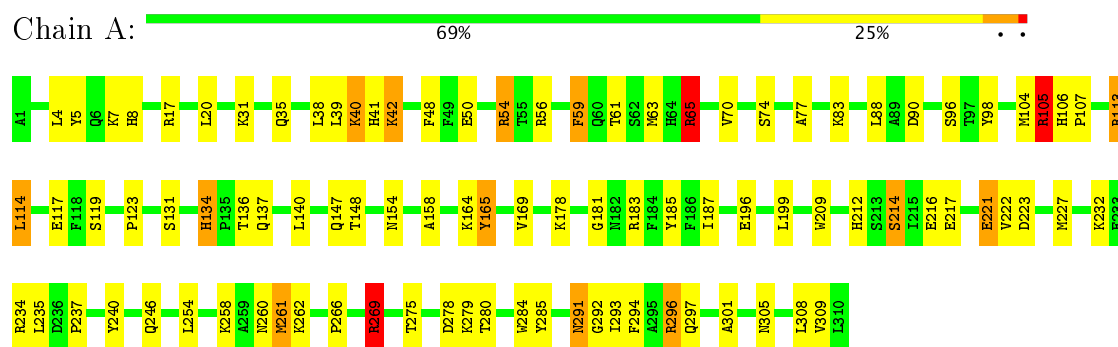
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	P	0	0
			29	9	3	14	3		
6	D	1	Total	C	N	O	P	0	0
			29	9	3	14	3		

3 Residue-property plots [i](#)

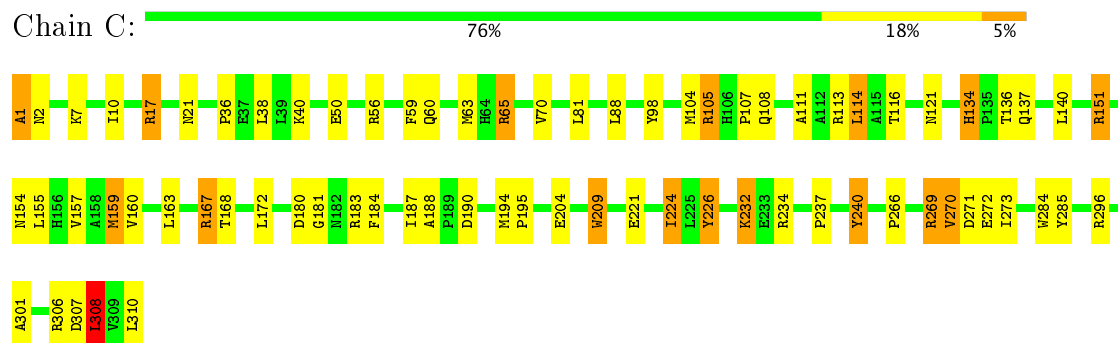
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

Note EDS was not executed.

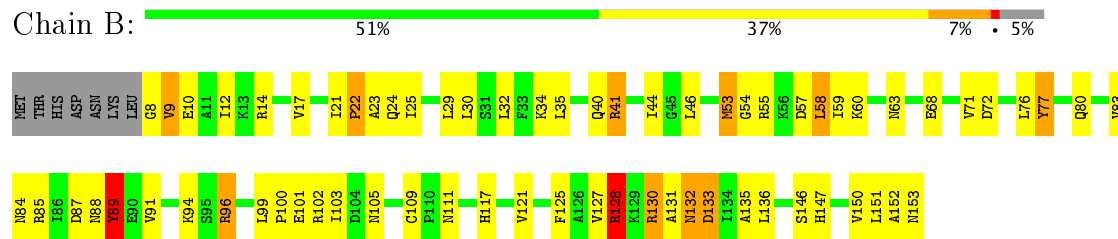
• Molecule 1: ASPARTATE CARBAMOYLTRANSFERASE (R STATE), CATALYTIC CHAIN



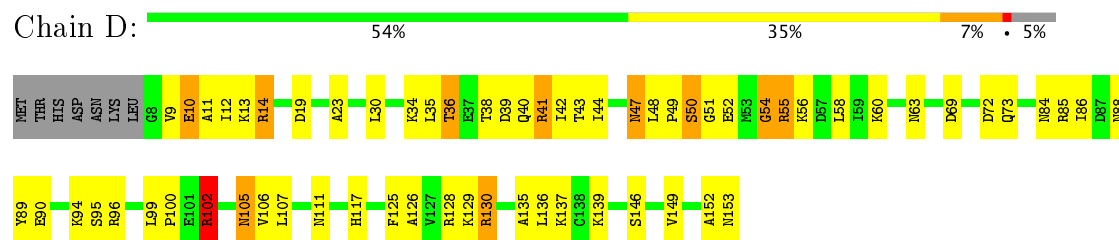
• Molecule 1: ASPARTATE CARBAMOYLTRANSFERASE (R STATE), CATALYTIC CHAIN



• Molecule 2: ASPARTATE CARBAMOYLTRANSFERASE REGULATORY CHAIN



• Molecule 2: ASPARTATE CARBAMOYLTRANSFERASE REGULATORY CHAIN



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	122.20 Å 122.20 Å 155.20 Å 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

5.1 Standard geometry

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

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	0.81	0/2461	1.58	34/3339 (1.0%)
1	C	0.83	0/2461	1.58	29/3339 (0.9%)
2	B	0.77	0/1155	1.51	10/1561 (0.6%)
2	D	0.74	0/1155	1.49	12/1561 (0.8%)
All	All	0.80	0/7232	1.55	85/9800 (0.9%)

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

There are no bond length outliers.

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	105	ARG	NE-CZ-NH2	15.21	127.91	120.30
1	C	105	ARG	NE-CZ-NH1	-14.38	113.11	120.30
1	A	105	ARG	NE-CZ-NH2	13.44	127.02	120.30
1	A	269	ARG	NE-CZ-NH2	13.42	127.01	120.30
1	A	269	ARG	NE-CZ-NH1	-11.28	114.66	120.30
1	C	151	ARG	NE-CZ-NH2	10.11	125.35	120.30
1	A	56	ARG	NE-CZ-NH2	9.17	124.88	120.30
1	A	284	TRP	CD1-CG-CD2	9.10	113.58	106.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	309	VAL	O-C-N	9.00	137.10	122.70
1	A	296	ARG	NE-CZ-NH2	8.75	124.67	120.30
1	A	309	VAL	CA-C-N	-8.70	98.06	117.20
1	C	209	TRP	CD1-CG-CD2	8.46	113.06	106.30
2	D	102	ARG	NE-CZ-NH2	8.43	124.52	120.30
2	D	14	ARG	NE-CZ-NH2	8.39	124.50	120.30
1	C	284	TRP	CD1-CG-CD2	8.28	112.93	106.30
1	C	56	ARG	NE-CZ-NH2	8.12	124.36	120.30
1	C	269	ARG	CA-CB-CG	7.94	130.87	113.40
1	A	284	TRP	CE2-CD2-CG	-7.87	101.00	107.30
2	D	41	ARG	NE-CZ-NH2	7.65	124.12	120.30
1	C	17	ARG	NE-CZ-NH2	7.64	124.12	120.30
1	A	234	ARG	NE-CZ-NH2	7.61	124.11	120.30
2	B	96	ARG	NE-CZ-NH1	-7.53	116.53	120.30
1	A	56	ARG	NE-CZ-NH1	-7.50	116.55	120.30
1	C	269	ARG	NE-CZ-NH2	7.50	124.05	120.30
1	C	183	ARG	NE-CZ-NH2	7.44	124.02	120.30
1	C	284	TRP	CE2-CD2-CG	-7.38	101.39	107.30
1	A	113	ARG	NE-CZ-NH2	7.26	123.93	120.30
1	C	209	TRP	CE2-CD2-CG	-7.24	101.51	107.30
1	C	308	LEU	CA-CB-CG	7.19	131.85	115.30
1	A	104	MET	CG-SD-CE	-7.18	88.70	100.20
1	C	113	ARG	NE-CZ-NH1	-7.14	116.73	120.30
1	A	169	VAL	CA-CB-CG1	-7.08	100.28	110.90
2	D	152	ALA	CA-C-N	-7.00	101.80	117.20
1	A	65	ARG	NE-CZ-NH2	6.96	123.78	120.30
1	A	209	TRP	CE2-CD2-CG	-6.92	101.76	107.30
1	A	209	TRP	CD1-CG-CD2	6.90	111.82	106.30
1	C	113	ARG	NE-CZ-NH2	6.76	123.68	120.30
1	A	98	TYR	CB-CG-CD2	-6.75	116.95	121.00
1	A	183	ARG	NE-CZ-NH1	-6.75	116.92	120.30
2	B	89	TYR	CB-CG-CD2	-6.72	116.97	121.00
2	B	101	GLU	CA-CB-CG	6.70	128.13	113.40
2	B	96	ARG	CA-CB-CG	-6.67	98.72	113.40
1	A	17	ARG	NE-CZ-NH2	6.62	123.61	120.30
1	C	234	ARG	NE-CZ-NH2	6.59	123.60	120.30
1	C	234	ARG	NE-CZ-NH1	-6.52	117.04	120.30
1	C	159	MET	CA-CB-CG	-6.39	102.44	113.30
2	B	85	ARG	NE-CZ-NH2	6.37	123.49	120.30
1	A	234	ARG	NE-CZ-NH1	-6.31	117.15	120.30
1	C	114	LEU	CA-CB-CG	6.25	129.69	115.30
2	B	130	ARG	NE-CZ-NH2	6.23	123.42	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	54	ARG	NE-CZ-NH1	-6.23	117.19	120.30
2	D	41	ARG	NE-CZ-NH1	-6.15	117.23	120.30
1	C	60	GLN	CA-CB-CG	-6.12	99.94	113.40
1	A	232	LYS	N-CA-CB	-6.02	99.76	110.60
2	D	130	ARG	NE-CZ-NH1	-5.97	117.32	120.30
1	A	296	ARG	NE-CZ-NH1	-5.96	117.32	120.30
1	A	183	ARG	NE-CZ-NH2	5.94	123.27	120.30
2	D	14	ARG	NE-CZ-NH1	-5.92	117.34	120.30
2	B	8	GLY	N-CA-C	-5.90	98.34	113.10
2	B	41	ARG	NE-CZ-NH2	5.89	123.24	120.30
2	B	152	ALA	N-CA-C	5.84	126.76	111.00
1	A	284	TRP	CG-CD1-NE1	-5.70	104.40	110.10
1	C	65	ARG	NE-CZ-NH2	5.70	123.15	120.30
1	A	209	TRP	CB-CG-CD1	-5.65	119.65	127.00
1	C	209	TRP	CG-CD1-NE1	-5.59	104.51	110.10
1	A	65	ARG	NE-CZ-NH1	-5.50	117.55	120.30
2	B	128	ARG	NE-CZ-NH2	5.49	123.05	120.30
2	D	90	GLU	CA-CB-CG	5.45	125.39	113.40
2	D	85	ARG	NE-CZ-NH2	5.42	123.01	120.30
1	C	1	ALA	N-CA-C	-5.39	96.43	111.00
1	C	226	TYR	CB-CG-CD2	-5.38	117.77	121.00
1	C	98	TYR	CB-CG-CD2	-5.38	117.77	121.00
1	C	272	GLU	CA-CB-CG	5.33	125.12	113.40
2	D	96	ARG	NE-CZ-NH2	5.24	122.92	120.30
1	C	232	LYS	CA-CB-CG	5.24	124.93	113.40
1	A	90	ASP	CB-CG-OD2	5.22	123.00	118.30
1	C	104	MET	CA-CB-CG	5.20	122.14	113.30
1	A	169	VAL	CA-CB-CG2	5.19	118.69	110.90
1	A	114	LEU	CA-CB-CG	5.19	127.23	115.30
1	A	261	MET	CA-CB-CG	5.16	122.07	113.30
2	D	139	LYS	CB-CG-CD	-5.05	98.46	111.60
1	A	221	GLU	N-CA-CB	-5.04	101.53	110.60
1	A	309	VAL	C-N-CA	5.02	134.24	121.70
1	C	183	ARG	NE-CZ-NH1	-5.01	117.80	120.30
2	D	9	VAL	CA-C-N	-5.01	106.18	117.20

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

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Mol	Chain	Res	Type	Group
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	42	0
1	C	2415	0	2422	29	0
2	B	1138	0	1154	37	0
2	D	1138	0	1154	40	0
3	A	7	0	0	0	0
3	C	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	144	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 (144) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:B:91:VAL:HG11	2:B:94:LYS:HZ2	1.70	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
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:106:VAL:HG23	2:D:107:LEU:HD13	1.91	0.52
2:D:49:PRO:HA	2:D:54:GLY:O	2.09	0.52
2:D:13:LYS:CG	2:D:88:ASN:HA	2.40	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
2:D:102:ARG:HB3	2:D:102:ARG:HH21	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:159:MET:HE1	1:C:172:LEU:HD23	1.91	0.50
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: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
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
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:D:99:LEU:HB3	2:D:129:LYS:HZ2	1.82	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'	2.33	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:99:LEU:HB3	2:D:129:LYS:NZ	2.32	0.43
2:D:42:ILE:HA	2:D:60:LYS:O	2.18	0.43
2:B:41:ARG:HD2	2:D:48:LEU:HD22	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:114:LEU:CD2	2:B:121:VAL:HG11	2.49	0.43
1:A:31:LYS:HD2	1:A:294:PHE:CD2	2.54	0.43
6:B:155:CTP:O3A	6:B:155:CTP:H3'	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
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
2:D:19:ASP:OD2	2:D:58:LEU:HD13	2.19	0.42
1:A:216:GLU:H	1:A:216:GLU:CD	2.23	0.42
1:C:273:ILE:HG21	1:C:273:ILE:HD13	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:147:HIS:O	2:B:151:LEU:HD12	2.20	0.41
2:B:89:TYR:HA	6:B:155:CTP:N4	2.35	0.41
1:A:275:THR:O	1:A:278:ASP:HB2	2.21	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:291:ASN:HA	1:A:291:ASN:HD22	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:HD3	1:C:7:LYS:HA	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%)	44	77
2	B	144/153 (94%)	125 (87%)	15 (10%)	4 (3%)	6	19
2	D	144/153 (94%)	126 (88%)	13 (9%)	5 (4%)	4	14
All	All	904/926 (98%)	842 (93%)	52 (6%)	10 (1%)	17	47

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
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%)	241 (92%)	20 (8%)	15	39
1	C	261/261 (100%)	239 (92%)	22 (8%)	13	35
2	B	129/136 (95%)	112 (87%)	17 (13%)	5	14
2	D	129/136 (95%)	115 (89%)	14 (11%)	7	22
All	All	780/794 (98%)	707 (91%)	73 (9%)	10	29

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	59	PHE
1	A	65	ARG
1	A	74	SER
1	A	105	ARG
1	A	113	ARG
1	A	117	GLU
1	A	131	SER
1	A	134	HIS
1	A	147	GLN

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Mol	Chain	Res	Type
1	A	214	SER
1	A	221	GLU
1	A	246	GLN
1	A	269	ARG
1	A	279	LYS
1	A	285	TYR
1	A	291	ASN
1	A	308	LEU
2	B	10	GLU
2	B	12	ILE
2	B	22	PRO
2	B	32	LEU
2	B	53	MET
2	B	58	LEU
2	B	68	GLU
2	B	72	ASP
2	B	76	LEU
2	B	88	ASN
2	B	102	ARG
2	B	105	ASN
2	B	128	ARG
2	B	132	ASN
2	B	133	ASP
2	B	146	SER
2	B	153	ASN
1	C	17	ARG
1	C	21	ASN
1	C	38	LEU
1	C	40	LYS
1	C	59	PHE
1	C	81	LEU
1	C	105	ARG
1	C	108	GLN
1	C	121	ASN
1	C	134	HIS
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

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Mol	Chain	Res	Type
1	C	269	ARG
1	C	285	TYR
1	C	307	ASP
1	C	308	LEU
1	C	310	LEU
2	D	34	LYS
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	55	ARG
2	D	95	SER
2	D	102	ARG
2	D	105	ASN
2	D	128	ARG
2	D	137	LYS
2	D	153	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	GLN
1	A	60	GLN
1	A	108	GLN
1	A	146	GLN
1	A	154	ASN
1	A	305	ASN
2	B	40	GLN
2	B	47	ASN
2	B	63	ASN
2	B	84	ASN
2	B	113	ASN
2	B	117	HIS
2	B	153	ASN
1	C	13	ASN
1	C	154	ASN
2	D	117	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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$
4	PCT	A	311	-	7,7,7	1.70	2 (28%)	8,10,10	3.35	3 (37%)
3	MAL	A	312	-	6,6,24	5.21	4 (66%)	5,7,35	4.48	5 (100%)
6	CTP	B	155	-	24,30,30	1.59	5 (20%)	24,47,47	1.49	5 (20%)
4	PCT	C	311	-	7,7,7	1.54	2 (28%)	8,10,10	2.12	3 (37%)
3	MAL	C	312	-	6,6,24	4.94	4 (66%)	5,7,35	4.06	5 (100%)
6	CTP	D	155	-	24,30,30	1.58	5 (20%)	24,47,47	1.38	3 (12%)

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
4	PCT	A	311	-	-	0/4/5/5	0/0/0/0
3	MAL	A	312	-	2/2/2/10	0/4/4/48	0/0/0/2
6	CTP	B	155	-	-	0/18/38/38	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PCT	C	311	-	-	0/4/5/5	0/0/0/0
3	MAL	C	312	-	2/2/2/10	0/4/4/48	0/0/0/2
6	CTP	D	155	-	-	0/18/38/38	0/2/2/2

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	312	MAL	O5-C1	-7.24	1.22	1.40
3	A	312	MAL	O1-C1	-7.06	1.23	1.40
3	C	312	MAL	O1-C1	-7.01	1.23	1.40
3	C	312	MAL	O5-C1	-6.82	1.23	1.40
3	A	312	MAL	C4-C3	-6.03	1.23	1.51
3	C	312	MAL	C4-C3	-5.85	1.24	1.51
3	A	312	MAL	O3-C3	-4.37	1.22	1.43
3	C	312	MAL	O3-C3	-4.07	1.24	1.43
6	B	155	CTP	C2'-C1'	-3.47	1.48	1.53
6	B	155	CTP	C6-N1	-3.14	1.31	1.35
6	D	155	CTP	C2-N3	-2.89	1.32	1.38
6	B	155	CTP	C2-N3	-2.46	1.33	1.38
6	D	155	CTP	C6-N1	-2.44	1.32	1.35
6	D	155	CTP	PB-O1B	2.22	1.59	1.50
6	B	155	CTP	PG-O1G	2.25	1.58	1.50
4	A	311	PCT	P-C1P	2.54	1.84	1.79
4	C	311	PCT	P-C1P	2.65	1.84	1.79
6	D	155	CTP	PG-O1G	2.66	1.59	1.50
6	B	155	CTP	PG-O3B	2.92	1.64	1.60
4	C	311	PCT	P-O1P	2.97	1.56	1.50
4	A	311	PCT	P-O1P	3.30	1.57	1.50
6	D	155	CTP	PG-O3B	4.01	1.66	1.60

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	311	PCT	C1P-C1-N1	-4.95	109.55	115.37
6	B	155	CTP	C6-N1-C2	-4.32	114.28	121.28
6	D	155	CTP	C6-N1-C2	-4.31	114.30	121.28
4	C	311	PCT	O1P-P-C1P	-2.34	105.39	110.97
6	B	155	CTP	O2'-C2'-C1'	-2.15	104.88	111.61
6	B	155	CTP	C4'-O4'-C1'	-2.04	107.59	109.77
6	D	155	CTP	C4'-O4'-C1'	2.21	112.12	109.77
6	D	155	CTP	N4-C4-N3	2.28	120.48	116.64
6	B	155	CTP	O4'-C1'-N1	2.42	112.92	108.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	312	MAL	O3-C3-C4	2.47	120.77	109.46
3	C	312	MAL	C4-C3-C2	2.63	118.21	112.56
6	B	155	CTP	N4-C4-N3	2.71	121.20	116.64
4	C	311	PCT	P-C1P-C1	2.85	120.23	114.70
4	A	311	PCT	P-C1P-C1	3.13	120.78	114.70
3	C	312	MAL	O3-C3-C4	3.20	124.13	109.46
3	A	312	MAL	C4-C3-C2	3.82	120.75	112.56
3	C	312	MAL	O1-C1-C2	4.02	117.02	109.13
4	C	311	PCT	O1-C1-C1P	4.32	124.37	119.77
3	C	312	MAL	O3-C3-C2	4.63	117.67	109.26
3	A	312	MAL	O3-C3-C2	5.07	118.45	109.26
3	A	312	MAL	O1-C1-C2	5.16	119.25	109.13
3	A	312	MAL	O5-C1-C2	5.25	119.44	109.13
3	C	312	MAL	O5-C1-C2	5.27	119.47	109.13
4	A	311	PCT	O1-C1-C1P	7.18	127.41	119.77

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	312	MAL	C3
3	A	312	MAL	C1
3	C	312	MAL	C3
3	C	312	MAL	C1

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 6 short contacts:

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

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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.