



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 01:11 pm BST

PDB ID : 1ZA2
Title : Structure of wild-type E. coli Aspartate Transcarbamoylase in the presence of CTP, carbamoyl phosphate at 2.50 Å resolution
Authors : Wang, J.; Stieglitz, K.A.; Cardia, J.P.; Kantrowitz, E.R.
Deposited on : 2005-04-05
Resolution : 2.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

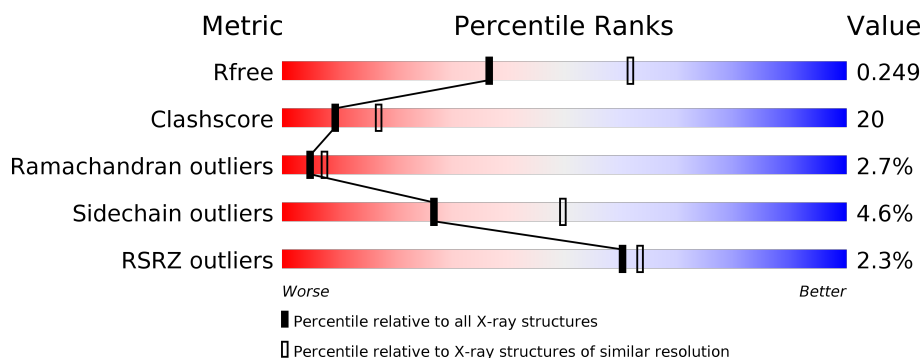
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 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	310	<div> <div>%</div> <div> <div></div> <div>71%</div> <div>25%</div> <div>•</div> </div> </div>
1	C	310	<div> <div>%</div> <div> <div></div> <div>73%</div> <div>23%</div> <div>•</div> </div> </div>
2	B	153	<div> <div>5%</div> <div> <div></div> <div>57%</div> <div>33%</div> <div>•</div> <div>6%</div> </div> </div>
2	D	153	<div> <div>5%</div> <div> <div></div> <div>48%</div> <div>41%</div> <div>6%</div> <div>6%</div> </div> </div>

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
4	CP	A	313	-	-	X	-
4	CP	C	312	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7558 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 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			

- Molecule 2 is a protein called Aspartate carbamoyltransferase regulatory chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	144	Total	C	N	O	S	0	0	0
			1127	707	199	216	5			
2	D	144	Total	C	N	O	S	0	0	0
			1127	707	199	216	5			

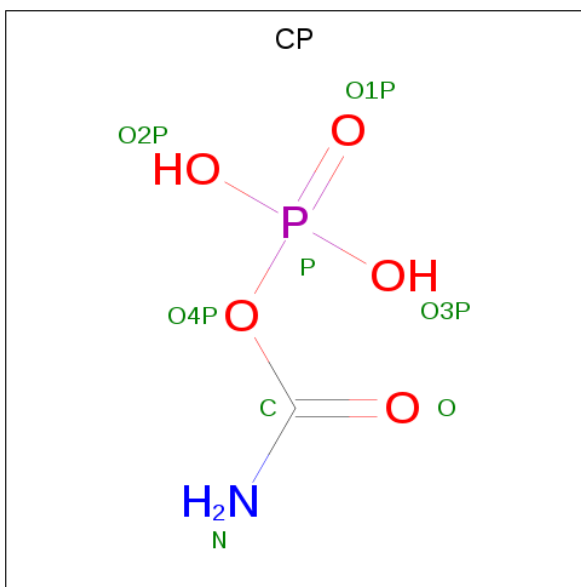
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	INITIATING METHIONINE	UNP P00478
D	1	MET	-	INITIATING METHIONINE	UNP P00478

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		

- Molecule 4 is PHOSPHORIC ACID MONO(FORMAMIDE)ESTER (three-letter code: CP) (formula: CH₄NO₅P).

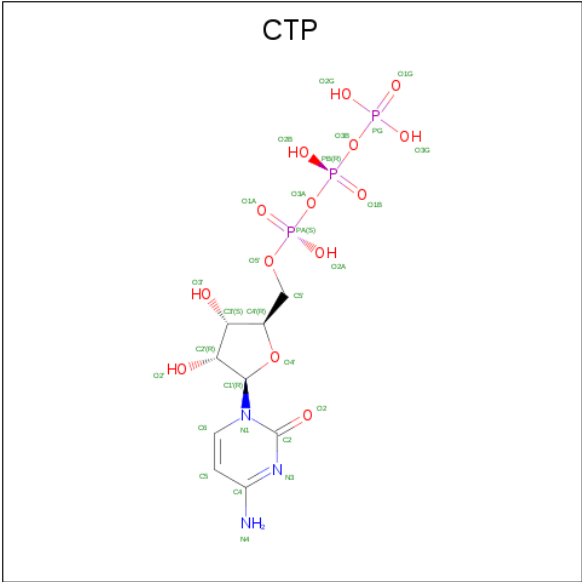


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			8	1	1	5	1		
4	A	1	Total	C	N	O	P	0	0
			8	1	1	5	1		
4	C	1	Total	C	N	O	P	0	0
			8	1	1	5	1		
4	C	1	Total	C	N	O	P	0	0
			8	1	1	5	1		

- 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: 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		

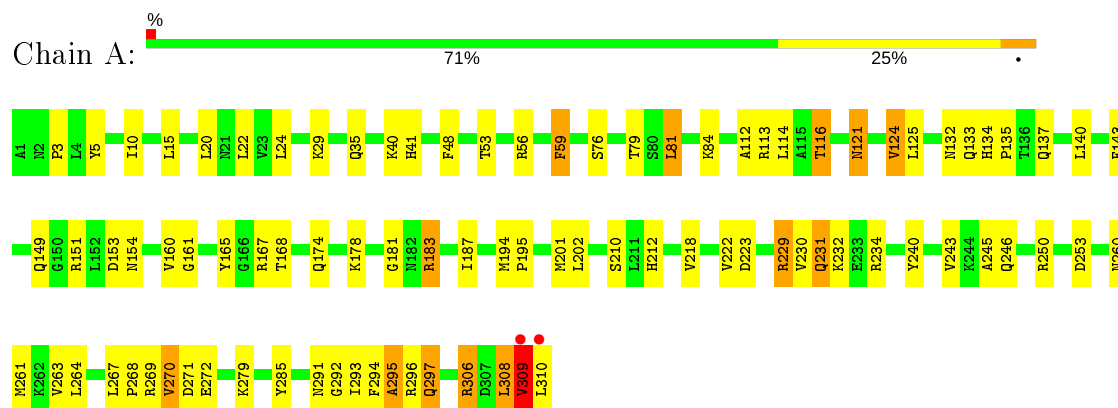
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	150	Total	O	0	0
			150	150		
7	B	57	Total	O	0	0
			57	57		
7	C	133	Total	O	0	0
			133	133		
7	D	41	Total	O	0	0
			41	41		

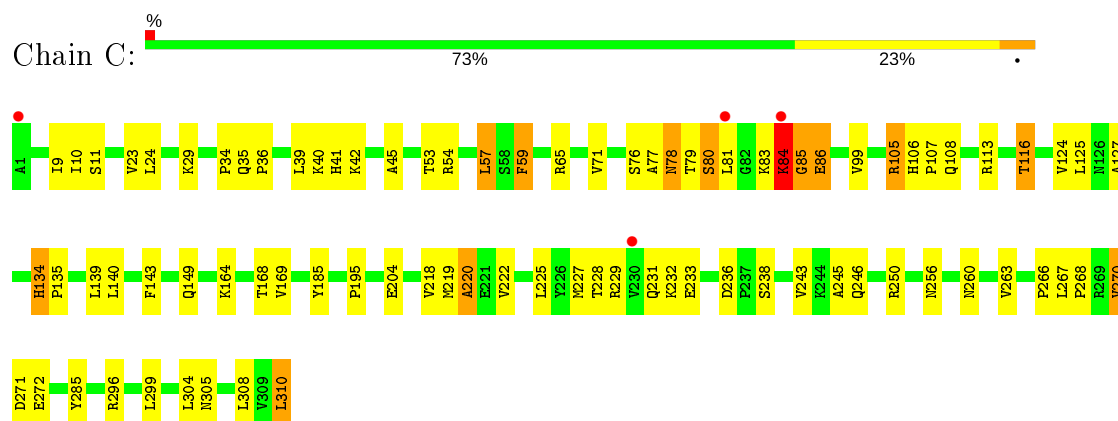
3 Residue-property plots

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.

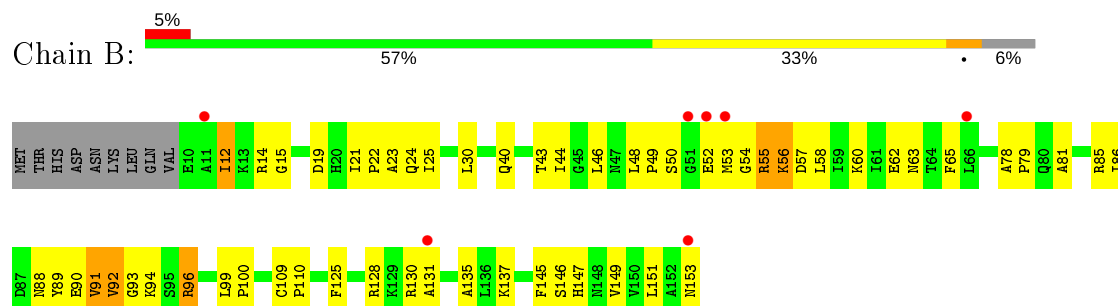
- Molecule 1: Aspartate carbamoyltransferase catalytic chain



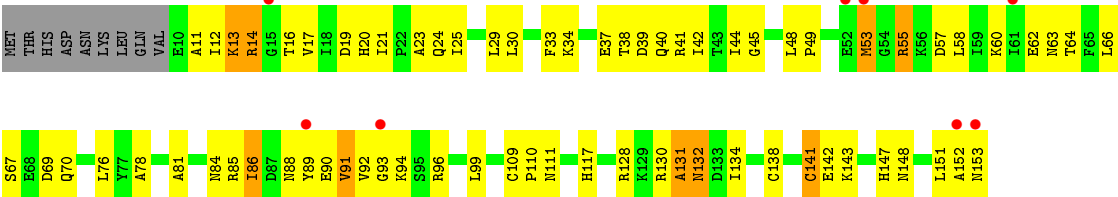
- Molecule 1: Aspartate carbamoyltransferase catalytic chain



- Molecule 2: Aspartate carbamoyltransferase regulatory chain



● Molecule 2: Aspartate carbamoyltransferase regulatory chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	121.44Å 121.44Å 142.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.50 49.34 – 2.49	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.50) 94.2 (49.34-2.49)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.48Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.197 , 0.251 0.194 , 0.249	Depositor DCC
R_{free} test set	4055 reflections (10.04%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7558	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.38	0/2461	0.65	1/3339 (0.0%)
1	C	0.35	0/2461	0.64	0/3339
2	B	0.32	0/1144	0.61	0/1546
2	D	0.34	0/1144	0.62	0/1546
All	All	0.35	0/7210	0.63	1/9770 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	81	LEU	N-CA-C	-6.25	94.12	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	85	0
1	C	2415	0	2422	77	0
2	B	1127	0	1142	51	0
2	D	1127	0	1142	72	0
3	A	1	0	0	0	1
4	A	16	0	4	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	16	0	4	8	0
5	B	1	0	0	0	0
5	D	1	0	0	0	0
6	B	29	0	12	0	0
6	D	29	0	12	5	0
7	A	150	0	0	5	1
7	B	57	0	0	0	0
7	C	133	0	0	2	1
7	D	41	0	0	2	0
All	All	7558	0	7160	283	3

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 20.

All (283) 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:THR:HG23	1:A:81:LEU:O	1.53	1.07
1:A:168:THR:HG23	4:A:313:CP:HN1	1.20	1.05
1:A:168:THR:HG23	4:A:313:CP:N	1.72	1.04
1:C:168:THR:HG23	4:C:312:CP:HN1	1.28	0.99
1:C:229:ARG:HH11	1:C:268:PRO:HG2	1.23	0.99
1:C:79:THR:O	1:C:86:GLU:HG2	1.65	0.96
1:A:149:GLN:HE22	1:A:260:ASN:HD22	1.13	0.95
1:C:77:ALA:HB1	7:C:426:HOH:O	1.71	0.89
2:D:128:ARG:HH21	2:D:130:ARG:HD3	1.38	0.87
2:B:55:ARG:HB2	2:B:55:ARG:NH1	1.90	0.86
1:C:229:ARG:NH1	1:C:268:PRO:HG2	1.95	0.82
2:D:138:CYS:HB3	2:D:141:CYS:HB2	1.62	0.81
1:C:35:GLN:HE22	1:C:310:LEU:HB3	1.45	0.80
1:A:149:GLN:HE22	1:A:260:ASN:ND2	1.81	0.79
1:A:149:GLN:NE2	1:A:260:ASN:HD22	1.82	0.78
1:C:78:ASN:HD22	1:C:78:ASN:C	1.88	0.77
2:D:141:CYS:HB3	2:D:143:LYS:H	1.49	0.76
1:C:24:LEU:HD22	1:C:143:PHE:HB2	1.68	0.76
1:C:168:THR:HG23	4:C:312:CP:N	2.02	0.75
2:D:86:ILE:O	2:D:86:ILE:HD12	1.86	0.74
1:A:308:LEU:HD21	1:A:310:LEU:HD13	1.69	0.74
1:A:294:PHE:O	1:A:295:ALA:HB3	1.88	0.74
1:A:154:ASN:HD22	1:A:181:GLY:HA3	1.53	0.73
1:C:168:THR:CG2	4:C:312:CP:HN1	2.00	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:GLY:HA3	7:A:463:HOH:O	1.88	0.72
1:A:29:LYS:HD3	1:A:310:LEU:HB3	1.71	0.72
1:A:294:PHE:O	1:A:295:ALA:CB	2.36	0.72
1:A:168:THR:H	4:A:313:CP:HN2	1.36	0.72
1:C:83:LYS:C	1:C:84:LYS:HD3	2.11	0.71
1:A:270:VAL:HG22	1:A:271:ASP:N	2.05	0.71
2:D:148:ASN:O	2:D:152:ALA:HB3	1.90	0.70
2:B:48:LEU:O	2:B:55:ARG:HA	1.91	0.70
2:B:96:ARG:HB2	2:B:96:ARG:NH1	2.07	0.70
1:A:292:GLY:O	1:A:296:ARG:HB2	1.92	0.70
1:A:168:THR:CG2	4:A:313:CP:HN1	2.01	0.68
1:C:243:VAL:O	1:C:246:GLN:HG2	1.93	0.68
2:D:12:ILE:HD11	2:D:62:GLU:HG3	1.75	0.68
1:A:174:GLN:HE22	1:A:201:MET:HG3	1.61	0.66
2:D:16:THR:HG22	2:D:64:THR:O	1.95	0.66
2:B:92:VAL:HG13	2:B:93:GLY:H	1.60	0.66
2:B:96:ARG:HB2	2:B:96:ARG:HH11	1.59	0.66
1:A:231:GLN:HG3	1:A:234:ARG:HE	1.61	0.65
1:C:149:GLN:HE22	1:C:260:ASN:HD22	1.44	0.65
2:B:55:ARG:CZ	2:B:55:ARG:HB2	2.24	0.65
2:B:40:GLN:NE2	2:B:63:ASN:HD22	1.94	0.65
1:C:85:GLY:O	1:C:86:GLU:C	2.34	0.65
1:A:183:ARG:NH2	1:A:210:SER:OG	2.29	0.65
2:D:11:ALA:HB1	6:D:155:CTP:N3	2.12	0.64
2:D:78:ALA:HB1	2:D:81:ALA:HB2	1.78	0.64
2:D:17:VAL:HG13	2:D:84:ASN:HB2	1.77	0.64
2:B:78:ALA:HB1	2:B:81:ALA:HB2	1.78	0.64
2:B:128:ARG:HG3	2:B:128:ARG:O	1.98	0.64
1:A:231:GLN:HB2	1:A:234:ARG:HD2	1.78	0.63
1:A:232:LYS:HE2	1:A:240:TYR:OH	1.97	0.63
1:A:308:LEU:HD23	1:A:309:VAL:N	2.13	0.63
1:A:76:SER:O	1:A:79:THR:HG22	1.96	0.63
2:D:67:SER:OG	2:D:70:GLN:HG3	1.97	0.63
1:A:308:LEU:CD2	1:A:310:LEU:HD13	2.28	0.62
2:B:30:LEU:HD11	2:B:44:ILE:HD13	1.82	0.62
1:A:15:LEU:O	1:A:178:LYS:HE3	1.99	0.62
1:C:76:SER:HB2	1:C:81:LEU:HD13	1.81	0.62
2:D:12:ILE:HG23	2:D:86:ILE:HD13	1.82	0.62
1:A:308:LEU:HD23	1:A:309:VAL:H	1.66	0.61
1:C:79:THR:HG22	1:C:86:GLU:HB2	1.81	0.61
1:C:76:SER:HB2	1:C:81:LEU:CD1	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:LYS:HG2	1:C:85:GLY:N	2.16	0.60
1:C:229:ARG:HD2	1:C:231:GLN:HG3	1.84	0.60
1:C:113:ARG:O	1:C:116:THR:HB	2.01	0.60
1:C:270:VAL:HG13	1:C:271:ASP:H	1.66	0.60
1:A:154:ASN:ND2	1:A:181:GLY:HA3	2.17	0.59
2:D:138:CYS:CB	2:D:141:CYS:HB2	2.28	0.59
2:D:12:ILE:HD11	2:D:62:GLU:HA	1.85	0.59
2:D:81:ALA:O	2:D:96:ARG:HD3	2.03	0.59
1:A:308:LEU:HD23	1:A:310:LEU:H	1.67	0.58
1:C:225:LEU:HD23	1:C:263:VAL:HG22	1.86	0.58
1:A:151:ARG:HD2	1:A:153:ASP:OD2	2.04	0.58
1:C:84:LYS:N	1:C:84:LYS:HD3	2.19	0.58
2:D:66:LEU:HA	2:D:70:GLN:OE1	2.03	0.57
1:A:291:ASN:O	1:A:294:PHE:O	2.22	0.57
1:A:271:ASP:HB3	7:A:455:HOH:O	2.03	0.57
1:C:76:SER:HB3	1:C:81:LEU:HB2	1.86	0.57
2:D:131:ALA:O	2:D:132:ASN:HB2	2.03	0.57
2:D:128:ARG:HH21	2:D:130:ARG:CD	2.14	0.57
1:A:309:VAL:HG13	1:A:309:VAL:O	2.04	0.57
2:B:55:ARG:HH11	2:B:55:ARG:HB2	1.70	0.56
1:C:164:LYS:HA	1:C:195:PRO:HD3	1.86	0.56
1:A:10:ILE:HD11	1:A:124:VAL:HG22	1.86	0.56
2:B:14:ARG:HH12	2:B:65:PHE:HE1	1.53	0.56
1:C:39:LEU:HD13	1:C:304:LEU:HD12	1.86	0.56
2:B:53:MET:HE2	2:B:55:ARG:O	2.05	0.56
2:B:81:ALA:O	2:B:96:ARG:HD3	2.05	0.56
2:D:138:CYS:HB3	2:D:141:CYS:CB	2.35	0.56
1:A:250:ARG:HG2	1:A:253:ASP:OD2	2.05	0.56
2:D:99:LEU:HD11	2:D:134:ILE:HD13	1.88	0.56
2:D:12:ILE:HG13	2:D:62:GLU:OE2	2.06	0.55
2:D:17:VAL:CG1	2:D:84:ASN:HB2	2.37	0.55
2:B:56:LYS:HD2	2:B:56:LYS:O	2.06	0.55
2:D:12:ILE:HG12	2:D:13:LYS:N	2.22	0.55
2:D:11:ALA:HB1	6:D:155:CTP:C4	2.42	0.54
1:A:3:PRO:HG2	1:A:22:LEU:HD22	1.88	0.54
1:A:267:LEU:HB3	1:A:268:PRO:HA	1.90	0.54
1:C:85:GLY:O	1:C:86:GLU:O	2.26	0.54
2:D:30:LEU:HD11	2:D:44:ILE:HD13	1.88	0.54
1:A:5:TYR:CD2	1:A:306:ARG:HA	2.42	0.54
2:B:22:PRO:HG2	2:B:25:ILE:HG13	1.89	0.54
2:D:84:ASN:OD1	2:D:94:LYS:HG2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:147:HIS:O	2:B:151:LEU:HD23	2.08	0.54
1:C:40:LYS:O	1:C:41:HIS:HB2	2.08	0.53
2:D:34:LYS:HE2	2:D:37:GLU:OE1	2.08	0.53
2:B:79:PRO:O	2:B:96:ARG:HD2	2.07	0.53
1:C:84:LYS:HG2	1:C:85:GLY:H	1.73	0.53
2:D:128:ARG:NH2	2:D:130:ARG:HD3	2.17	0.53
2:B:44:ILE:HB	2:D:44:ILE:HB	1.91	0.53
2:D:60:LYS:HE3	6:D:155:CTP:O2	2.08	0.53
2:D:25:ILE:O	2:D:29:LEU:HG	2.09	0.53
1:C:78:ASN:C	1:C:78:ASN:ND2	2.60	0.53
1:A:270:VAL:HG22	1:A:271:ASP:HB2	1.90	0.52
2:B:85:ARG:C	2:B:86:ILE:HD12	2.30	0.52
2:B:40:GLN:NE2	2:B:40:GLN:HA	2.25	0.52
2:D:19:ASP:HA	2:D:58:LEU:HD12	1.92	0.52
1:A:218:VAL:HG22	1:A:222:VAL:HG13	1.90	0.52
1:A:125:LEU:HD12	1:A:125:LEU:N	2.25	0.52
1:A:309:VAL:CG1	1:A:309:VAL:O	2.58	0.52
1:C:185:TYR:CG	1:C:218:VAL:HG21	2.44	0.52
2:B:146:SER:O	2:B:149:VAL:HG22	2.10	0.51
1:C:53:THR:O	1:C:57:LEU:HB2	2.09	0.51
1:A:194:MET:SD	1:A:195:PRO:HD2	2.50	0.51
2:D:38:THR:HG22	2:D:39:ASP:N	2.25	0.51
1:A:113:ARG:O	1:A:116:THR:HB	2.10	0.51
1:A:230:VAL:N	7:A:463:HOH:O	2.43	0.51
1:A:3:PRO:HG2	1:A:22:LEU:CD2	2.40	0.51
1:C:84:LYS:CG	1:C:85:GLY:N	2.72	0.51
2:D:111:ASN:O	2:D:117:HIS:HE1	1.94	0.51
2:D:13:LYS:CE	2:D:88:ASN:HA	2.41	0.51
2:D:13:LYS:HG3	2:D:14:ARG:N	2.26	0.51
1:C:35:GLN:NE2	1:C:310:LEU:HB3	2.20	0.51
1:C:79:THR:HG22	1:C:86:GLU:CB	2.41	0.51
2:D:130:ARG:O	2:D:131:ALA:HB3	2.11	0.51
2:B:52:GLU:HG2	2:B:53:MET:N	2.26	0.50
2:D:85:ARG:HB2	2:D:93:GLY:H	1.75	0.50
1:C:125:LEU:HD12	1:C:125:LEU:N	2.25	0.50
2:D:13:LYS:HE2	2:D:88:ASN:HA	1.93	0.50
2:B:99:LEU:HD12	2:B:100:PRO:HD2	1.93	0.50
2:D:23:ALA:O	2:D:24:GLN:HB2	2.12	0.50
1:A:230:VAL:O	1:A:232:LYS:N	2.45	0.50
1:C:29:LYS:HD3	1:C:310:LEU:HD13	1.94	0.50
1:C:9:ILE:HG21	1:C:299:LEU:HD21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:LYS:O	1:C:85:GLY:C	2.51	0.49
1:C:229:ARG:HD2	1:C:231:GLN:CG	2.43	0.49
2:B:43:THR:HB	2:B:60:LYS:HB2	1.94	0.49
1:C:105:ARG:HD3	1:C:127:ALA:O	2.12	0.49
1:C:78:ASN:HD21	1:C:80:SER:HB3	1.76	0.49
2:D:131:ALA:O	2:D:132:ASN:CB	2.61	0.49
1:C:83:LYS:O	1:C:84:LYS:O	2.31	0.49
1:A:309:VAL:O	1:A:310:LEU:C	2.51	0.49
2:D:20:HIS:HB3	2:D:53:MET:HE1	1.95	0.49
1:A:59:PHE:CE2	1:A:296:ARG:HG3	2.48	0.48
2:B:49:PRO:HA	2:B:54:GLY:O	2.12	0.48
1:C:11:SER:HB2	2:D:142:GLU:OE1	2.14	0.48
1:A:293:ILE:O	1:A:297:GLN:HB2	2.12	0.48
2:D:45:GLY:HA3	2:D:48:LEU:HD11	1.95	0.48
2:B:23:ALA:O	2:B:24:GLN:HB2	2.13	0.48
2:B:145:PHE:HB3	2:B:149:VAL:CG2	2.43	0.48
1:C:76:SER:CB	1:C:81:LEU:HB2	2.44	0.48
1:A:270:VAL:HG13	1:A:272:GLU:OE2	2.12	0.48
1:C:80:SER:HA	1:C:86:GLU:HB3	1.96	0.48
1:A:149:GLN:NE2	1:A:223:ASP:HB3	2.29	0.48
1:C:250:ARG:HG2	1:C:250:ARG:HH21	1.77	0.48
1:A:308:LEU:CG	1:A:310:LEU:HD13	2.43	0.48
1:A:308:LEU:HD21	1:A:310:LEU:HD22	1.95	0.47
1:A:59:PHE:CE1	1:A:296:ARG:HD2	2.49	0.47
1:C:168:THR:H	4:C:312:CP:HN1	1.61	0.47
1:C:218:VAL:O	1:C:218:VAL:HG22	2.14	0.47
1:A:5:TYR:CE2	1:A:306:ARG:HA	2.49	0.47
1:C:10:ILE:HD11	1:C:124:VAL:HG22	1.97	0.47
1:C:59:PHE:CD1	1:C:296:ARG:HD3	2.50	0.47
2:B:48:LEU:HD23	2:D:41:ARG:HG3	1.96	0.47
2:D:13:LYS:HD3	2:D:89:TYR:CE2	2.50	0.46
1:A:187:ILE:HD12	1:A:187:ILE:N	2.30	0.46
1:C:227:MET:O	1:C:266:PRO:HD2	2.15	0.46
1:A:229:ARG:HG2	1:A:229:ARG:O	2.14	0.46
1:A:270:VAL:CG2	1:A:271:ASP:N	2.75	0.46
1:C:54:ARG:CG	4:C:311:CP:O4P	2.64	0.46
1:C:79:THR:O	1:C:80:SER:HB2	2.16	0.46
2:B:90:GLU:HA	2:B:90:GLU:OE1	2.15	0.46
2:B:91:VAL:O	2:B:91:VAL:HG12	2.17	0.46
1:C:218:VAL:HG22	1:C:222:VAL:HG13	1.98	0.46
2:D:49:PRO:HA	2:D:55:ARG:HD2	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:86:ILE:HG22	2:D:91:VAL:HA	1.97	0.45
1:A:114:LEU:HD23	1:A:114:LEU:C	2.37	0.45
2:B:145:PHE:HB3	2:B:149:VAL:HG21	1.97	0.45
1:C:106:HIS:CG	1:C:107:PRO:HD2	2.52	0.45
1:A:261:MET:SD	1:A:261:MET:C	2.95	0.45
2:D:40:GLN:O	2:D:42:ILE:HG13	2.16	0.45
2:B:52:GLU:C	2:B:54:GLY:H	2.18	0.45
2:D:91:VAL:HG21	6:D:155:CTP:O1A	2.16	0.45
2:D:76:LEU:HD13	2:D:151:LEU:HD21	1.98	0.45
1:C:270:VAL:HG13	1:C:272:GLU:OE2	2.17	0.44
2:D:21:ILE:HG22	2:D:25:ILE:HB	1.99	0.44
1:A:160:VAL:HG22	1:A:187:ILE:HB	1.98	0.44
2:B:30:LEU:CD1	2:B:44:ILE:HD13	2.48	0.44
2:B:56:LYS:C	2:B:56:LYS:HD2	2.37	0.44
1:C:39:LEU:O	1:C:42:LYS:HB2	2.18	0.44
1:A:218:VAL:HG22	1:A:218:VAL:O	2.17	0.44
1:A:229:ARG:C	7:A:463:HOH:O	2.55	0.44
2:B:130:ARG:CZ	2:B:135:ALA:HB2	2.48	0.44
2:B:14:ARG:NH1	2:B:65:PHE:HE1	2.14	0.44
2:D:94:LYS:NZ	6:D:155:CTP:O2G	2.49	0.44
2:B:125:PHE:HA	2:B:137:LYS:O	2.18	0.44
1:C:267:LEU:HB3	1:C:268:PRO:HA	2.00	0.44
4:C:312:CP:O3P	4:C:312:CP:O	2.32	0.43
1:C:45:ALA:HB2	1:C:99:VAL:HG11	2.00	0.43
2:B:46:LEU:HA	2:B:57:ASP:OD1	2.18	0.43
2:D:109:CYS:HA	2:D:110:PRO:HD3	1.84	0.43
1:A:201:MET:HE1	1:A:202:LEU:CD2	2.48	0.43
2:B:92:VAL:HG22	2:B:93:GLY:N	2.32	0.43
1:C:59:PHE:CG	1:C:296:ARG:HD3	2.53	0.43
2:D:147:HIS:O	2:D:151:LEU:HD23	2.18	0.43
1:A:201:MET:HE1	1:A:202:LEU:HD23	1.99	0.43
1:A:35:GLN:NE2	1:A:35:GLN:HA	2.34	0.43
2:D:33:PHE:O	2:D:34:LYS:C	2.57	0.43
2:D:69:ASP:HB3	7:D:187:HOH:O	2.18	0.43
2:B:52:GLU:OE1	2:B:52:GLU:N	2.52	0.43
1:C:169:VAL:HG13	1:C:228:THR:HG21	2.01	0.43
1:A:165:TYR:CD2	1:A:234:ARG:HD3	2.54	0.43
1:C:36:PRO:HA	1:C:65:ARG:O	2.18	0.43
2:D:76:LEU:CD2	2:D:134:ILE:HD12	2.48	0.43
2:B:19:ASP:HA	2:B:58:LEU:CD2	2.48	0.43
2:D:86:ILE:HA	2:D:90:GLU:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:141:CYS:O	2:D:142:GLU:HB2	2.19	0.42
2:D:153:ASN:ND2	7:D:194:HOH:O	2.49	0.42
1:A:20:LEU:O	1:A:24:LEU:HG	2.18	0.42
1:A:308:LEU:HG	1:A:310:LEU:HD13	2.01	0.42
1:A:132:ASN:ND2	1:A:133:GLN:HG2	2.35	0.42
1:C:270:VAL:HG22	1:C:271:ASP:N	2.33	0.42
2:B:130:ARG:NH2	2:B:135:ALA:HB2	2.35	0.42
1:A:167:ARG:HG2	4:A:313:CP:N	2.34	0.42
2:B:52:GLU:HG2	2:B:53:MET:H	1.84	0.42
2:B:15:GLY:HA3	2:B:62:GLU:HA	2.02	0.42
2:D:21:ILE:HB	2:D:57:ASP:HB2	2.01	0.42
2:D:42:ILE:HG23	2:D:60:LYS:O	2.19	0.42
1:A:40:LYS:O	1:A:41:HIS:HB2	2.20	0.42
1:A:48:PHE:CE1	1:A:56:ARG:HB2	2.55	0.42
1:A:263:VAL:C	1:A:264:LEU:HD23	2.40	0.42
1:C:23:VAL:HG11	1:C:139:LEU:HD13	2.00	0.42
2:D:12:ILE:CG1	2:D:13:LYS:N	2.83	0.42
2:D:85:ARG:O	2:D:86:ILE:C	2.57	0.42
1:A:149:GLN:NE2	1:A:260:ASN:ND2	2.53	0.41
1:C:219:MET:O	1:C:222:VAL:HG22	2.20	0.41
1:C:308:LEU:N	1:C:308:LEU:HD22	2.35	0.41
2:D:111:ASN:O	2:D:117:HIS:CE1	2.72	0.41
1:C:29:LYS:CD	1:C:310:LEU:HD13	2.50	0.41
1:A:269:ARG:HG2	7:A:331:HOH:O	2.20	0.41
2:D:14:ARG:HD2	2:D:63:ASN:OD1	2.19	0.41
1:C:236:ASP:OD2	1:C:238:SER:HB2	2.20	0.41
2:D:38:THR:HG22	2:D:40:GLN:H	1.85	0.41
1:A:210:SER:OG	1:A:212:HIS:NE2	2.47	0.41
2:D:58:LEU:HD21	2:D:60:LYS:NZ	2.36	0.41
2:D:66:LEU:HB3	2:D:70:GLN:HB2	2.01	0.41
1:C:232:LYS:HG3	1:C:233:GLU:N	2.36	0.41
1:C:220:ALA:HB2	1:C:256:ASN:HD22	1.86	0.41
1:A:121:ASN:C	1:A:121:ASN:HD22	2.24	0.41
1:A:133:GLN:HB3	1:A:135:PRO:HD3	2.02	0.41
1:A:134:HIS:CE1	1:A:137:GLN:HB2	2.56	0.41
1:A:140:LEU:C	1:A:140:LEU:HD12	2.41	0.41
1:A:24:LEU:HD22	1:A:143:PHE:HB2	2.02	0.41
2:B:130:ARG:HG3	2:B:130:ARG:HH11	1.86	0.41
2:B:12:ILE:HB	2:B:62:GLU:HG2	2.03	0.41
1:C:168:THR:CB	4:C:312:CP:HN1	2.34	0.41
1:C:45:ALA:HA	1:C:71:VAL:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:GLN:O	1:A:140:LEU:HG	2.21	0.41
2:B:50:SER:HB2	2:B:56:LYS:HG3	2.03	0.41
1:C:54:ARG:HG2	4:C:311:CP:P	2.61	0.40
2:D:33:PHE:C	2:D:34:LYS:HG2	2.41	0.40
2:B:109:CYS:HA	2:B:110:PRO:HD3	1.77	0.40
1:A:243:VAL:HA	1:A:246:GLN:NE2	2.36	0.40
2:B:21:ILE:HB	2:B:57:ASP:HB2	2.02	0.40
1:C:185:TYR:CD2	1:C:218:VAL:HG21	2.56	0.40
2:D:12:ILE:CD1	2:D:62:GLU:HG3	2.45	0.40
1:C:134:HIS:N	1:C:135:PRO:CD	2.84	0.40
1:A:10:ILE:HD12	1:A:112:ALA:HB1	2.03	0.40
2:B:22:PRO:HA	2:B:53:MET:CE	2.51	0.40
1:C:34:PRO:HB3	7:C:395:HOH:O	2.21	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:311:NA:NA	3:A:311:NA:NA[2_655]	0.80	1.40
7:C:361:HOH:O	7:C:361:HOH:O[2_655]	1.86	0.34
7:A:456:HOH:O	7:A:457:HOH:O[3_665]	2.14	0.06

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%)	284 (92%)	18 (6%)	6 (2%)	8	13
1	C	308/310 (99%)	281 (91%)	21 (7%)	6 (2%)	8	13
2	B	142/153 (93%)	121 (85%)	16 (11%)	5 (4%)	3	4
2	D	142/153 (93%)	122 (86%)	13 (9%)	7 (5%)	2	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	900/926 (97%)	808 (90%)	68 (8%)	24 (3%)	5 7

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	295	ALA
1	A	306	ARG
1	A	309	VAL
2	B	12	ILE
2	B	131	ALA
1	C	80	SER
1	C	84	LYS
2	D	86	ILE
2	D	92	VAL
1	C	85	GLY
1	C	245	ALA
2	D	13	LYS
1	A	231	GLN
1	A	245	ALA
2	B	92	VAL
2	D	132	ASN
1	A	308	LEU
2	B	89	TYR
1	C	220	ALA
2	D	53	MET
2	D	131	ALA
1	C	86	GLU
2	D	91	VAL
2	B	91	VAL

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%)	248 (95%)	13 (5%)	24 46

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	261/261 (100%)	247 (95%)	14 (5%)	22	42
2	B	128/137 (93%)	122 (95%)	6 (5%)	26	49
2	D	128/137 (93%)	125 (98%)	3 (2%)	50	76
All	All	778/796 (98%)	742 (95%)	36 (5%)	27	50

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	THR
1	A	59	PHE
1	A	84	LYS
1	A	116	THR
1	A	121	ASN
1	A	124	VAL
1	A	183	ARG
1	A	229	ARG
1	A	270	VAL
1	A	279	LYS
1	A	285	TYR
1	A	297	GLN
1	A	309	VAL
2	B	55	ARG
2	B	56	LYS
2	B	88	ASN
2	B	94	LYS
2	B	96	ARG
2	B	153	ASN
1	C	57	LEU
1	C	59	PHE
1	C	78	ASN
1	C	84	LYS
1	C	105	ARG
1	C	108	GLN
1	C	116	THR
1	C	134	HIS
1	C	140	LEU
1	C	204	GLU
1	C	270	VAL
1	C	285	TYR
1	C	305	ASN
1	C	310	LEU

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Mol	Chain	Res	Type
2	D	14	ARG
2	D	55	ARG
2	D	141	CYS

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

Mol	Chain	Res	Type
1	A	21	ASN
1	A	35	GLN
1	A	121	ASN
1	A	149	GLN
1	A	154	ASN
1	A	174	GLN
1	A	196	GLN
1	A	231	GLN
1	A	242	ASN
1	A	246	GLN
1	A	291	ASN
2	B	40	GLN
2	B	73	GLN
2	B	88	ASN
2	B	132	ASN
2	B	153	ASN
1	C	21	ASN
1	C	35	GLN
1	C	78	ASN
1	C	149	GLN
1	C	196	GLN
1	C	291	ASN
2	D	47	ASN
2	D	88	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 9 ligands modelled in this entry, 3 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	CP	A	312	-	6,7,7	1.83	1 (16%)	7,10,10	1.56	1 (14%)
4	CP	A	313	-	6,7,7	1.88	1 (16%)	7,10,10	1.63	2 (28%)
6	CTP	B	155	-	23,30,30	1.77	6 (26%)	30,47,47	1.61	6 (20%)
6	CTP	D	155	-	23,30,30	1.78	6 (26%)	30,47,47	1.49	5 (16%)
4	CP	C	311	-	6,7,7	1.85	1 (16%)	7,10,10	1.70	2 (28%)
4	CP	C	312	-	6,7,7	1.89	1 (16%)	7,10,10	1.73	2 (28%)

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	CP	A	312	-	-	0/3/5/5	-
4	CP	A	313	-	-	0/3/5/5	-
6	CTP	B	155	-	-	2/20/38/38	0/2/2/2
6	CTP	D	155	-	-	4/20/38/38	0/2/2/2
4	CP	C	311	-	-	0/3/5/5	-
4	CP	C	312	-	-	0/3/5/5	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	155	CTP	C6-N1	3.95	1.40	1.35
6	D	155	CTP	C6-N1	3.83	1.40	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	312	CP	P-O1P	3.41	1.61	1.50
4	A	313	CP	P-O1P	3.40	1.61	1.50
6	D	155	CTP	PG-O1G	3.40	1.61	1.50
6	B	155	CTP	PG-O1G	3.36	1.61	1.50
4	A	312	CP	P-O1P	3.34	1.61	1.50
4	C	311	CP	P-O1P	3.31	1.61	1.50
6	D	155	CTP	C4-N3	3.21	1.40	1.35
6	B	155	CTP	C4-N3	3.09	1.40	1.35
6	B	155	CTP	PA-O1A	2.98	1.61	1.50
6	D	155	CTP	PA-O1A	2.97	1.61	1.50
6	D	155	CTP	PB-O1B	2.97	1.61	1.50
6	B	155	CTP	PB-O1B	2.97	1.61	1.50
6	D	155	CTP	O4'-C1'	2.41	1.44	1.41
6	B	155	CTP	O4'-C1'	2.15	1.44	1.41

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	155	CTP	C2-N3-C4	4.02	120.42	116.34
6	B	155	CTP	C2-N3-C4	3.85	120.25	116.34
6	B	155	CTP	PB-O3B-PG	-3.79	119.84	132.83
6	D	155	CTP	PB-O3A-PA	-3.49	120.86	132.83
6	B	155	CTP	O4'-C1'-C2'	-3.41	101.94	106.93
6	D	155	CTP	PB-O3B-PG	-3.35	121.33	132.83
6	B	155	CTP	PB-O3A-PA	-3.33	121.38	132.83
4	A	313	CP	O-C-N	-2.93	120.67	125.51
4	C	312	CP	O3P-P-O4P	2.91	114.14	105.25
4	A	312	CP	O-C-N	-2.85	120.82	125.51
4	C	312	CP	O-C-N	-2.80	120.89	125.51
4	C	311	CP	O3P-P-O4P	2.79	113.75	105.25
4	C	311	CP	O-C-N	-2.52	121.34	125.51
6	B	155	CTP	O3G-PG-O3B	2.44	112.81	104.64
6	D	155	CTP	O3G-PG-O3B	2.42	112.74	104.64
4	A	313	CP	O3P-P-O4P	2.30	112.26	105.25
6	D	155	CTP	N4-C4-N3	2.15	119.88	116.49
6	B	155	CTP	N4-C4-N3	2.10	119.81	116.49

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	155	CTP	C5'-O5'-PA-O3A

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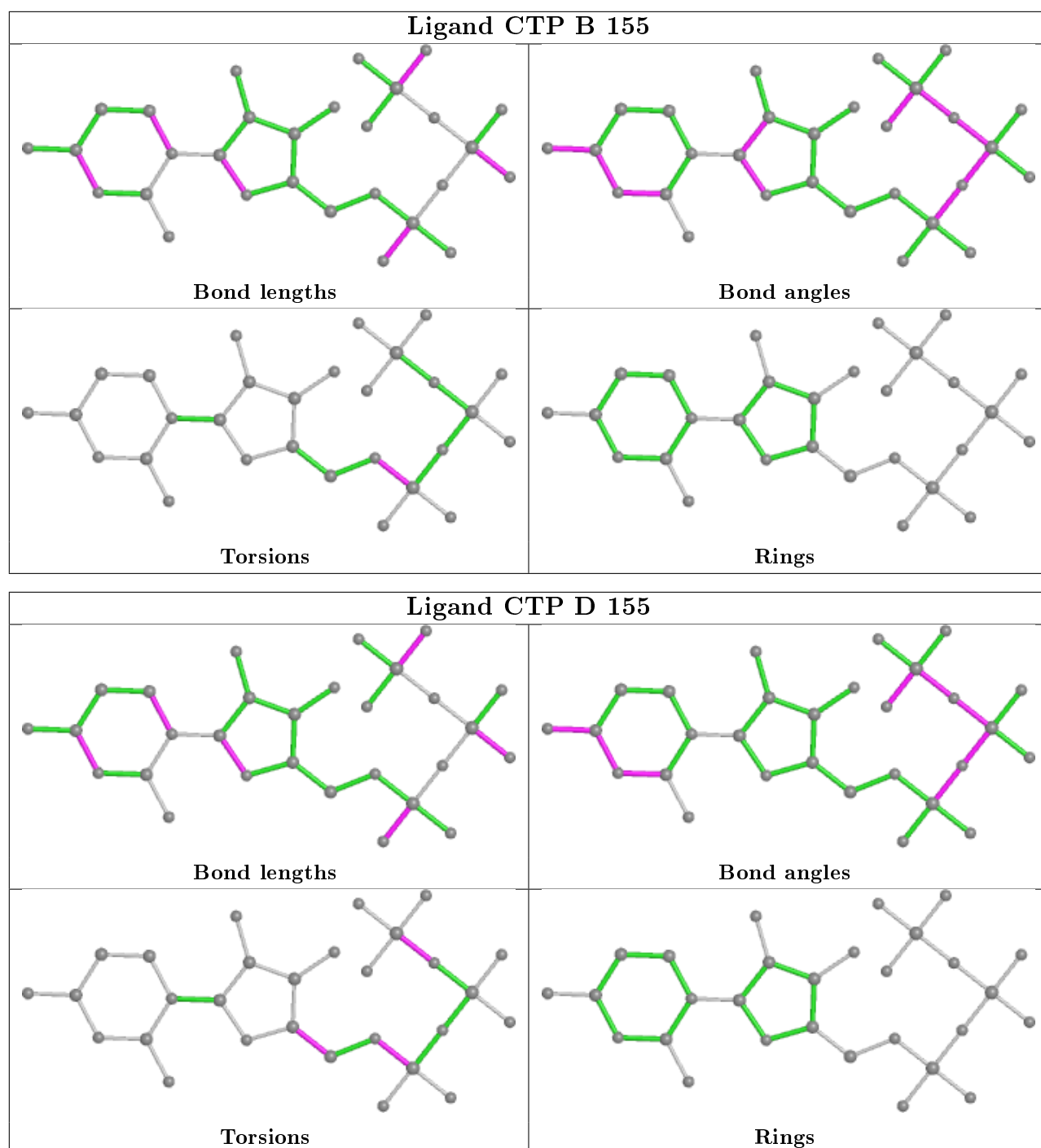
Mol	Chain	Res	Type	Atoms
6	D	155	CTP	PB-O3B-PG-O1G
6	B	155	CTP	C5'-O5'-PA-O3A
6	B	155	CTP	C5'-O5'-PA-O2A
6	D	155	CTP	C5'-O5'-PA-O2A
6	D	155	CTP	O4'-C4'-C5'-O5'

There are no ring outliers.

4 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	313	CP	5	0
6	D	155	CTP	5	0
4	C	311	CP	2	0
4	C	312	CP	6	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 ⓘ

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	310/310 (100%)	-0.56	2 (0%) 89 90	24, 39, 63, 112	0
1	C	310/310 (100%)	-0.48	4 (1%) 77 79	28, 43, 81, 91	0
2	B	144/153 (94%)	-0.08	7 (4%) 29 31	33, 62, 103, 119	0
2	D	144/153 (94%)	0.05	8 (5%) 24 25	35, 66, 107, 119	1 (0%)
All	All	908/926 (98%)	-0.36	21 (2%) 60 63	24, 47, 88, 119	1 (0%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	51	GLY	4.8
2	D	53	MET	4.1
2	B	153	ASN	4.0
2	D	52	GLU	3.9
1	A	310	LEU	3.8
1	C	84	LYS	3.8
2	D	89	TYR	3.4
1	A	309	VAL	3.3
2	B	53	MET	3.1
2	D	153	ASN	3.1
2	D	152	ALA	3.0
2	B	52	GLU	2.8
1	C	230	VAL	2.8
2	B	131	ALA	2.5
2	D	61	ILE	2.4
1	C	1	ALA	2.4
2	B	66	LEU	2.3
2	D	15	GLY	2.1
2	D	93	GLY	2.1
2	B	11	ALA	2.1
1	C	81	LEU	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

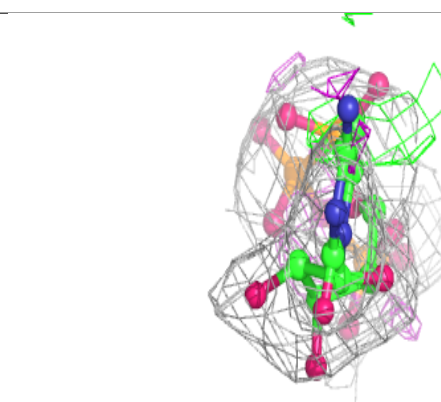
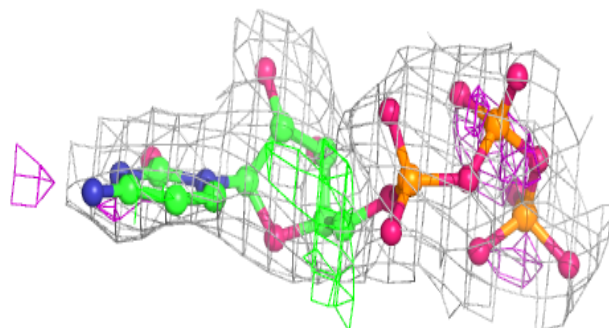
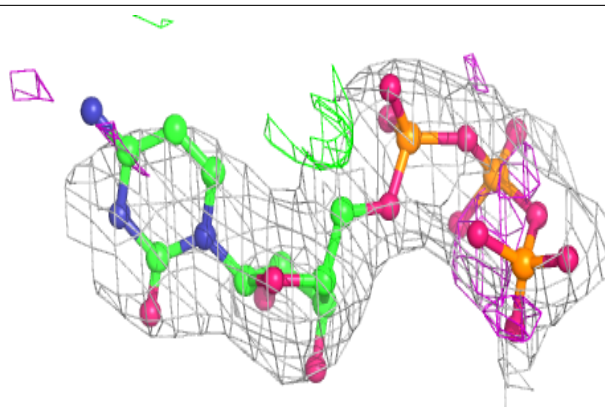
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	CP	C	312	8/8	0.86	0.17	74,82,87,87	0
6	CTP	D	155	29/29	0.89	0.16	82,85,94,95	0
4	CP	A	313	8/8	0.89	0.15	75,82,85,85	0
6	CTP	B	155	29/29	0.93	0.12	73,79,92,92	0
3	NA	A	311	1/1	0.98	0.16	40,40,40,40	0
4	CP	A	312	8/8	0.98	0.11	22,27,30,31	0
5	ZN	B	154	1/1	0.99	0.11	35,35,35,35	0
4	CP	C	311	8/8	0.99	0.12	24,29,31,31	0
5	ZN	D	154	1/1	1.00	0.10	37,37,37,37	0

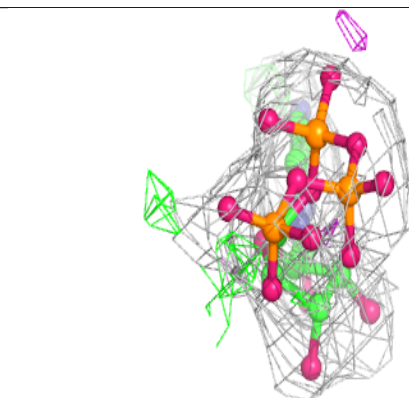
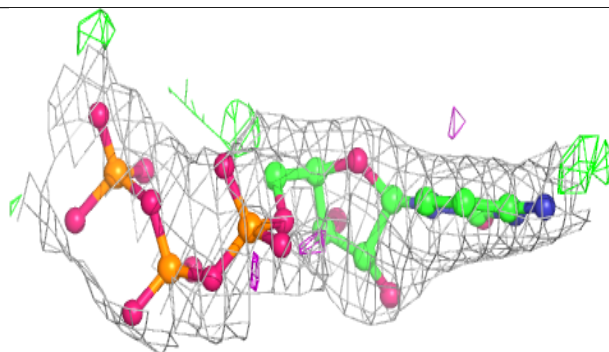
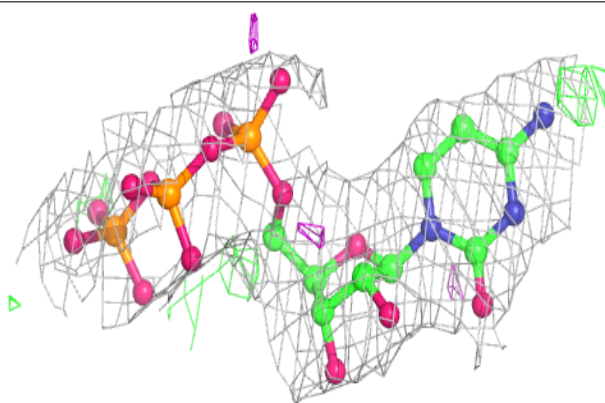
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CTP D 155:

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

**Electron density around CTP B 155:**

$2mF_o-DF_c$ (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.