



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:03 PM UTC

PDB ID : 1MIY / pdb_00001miy
Title : Crystal structure of Bacillus stearothermophilus CCA-adding enzyme in complex with CTP
Authors : Li, F.; Xiong, Y.; Wang, J.; Cho, H.D.; Weiner, A.M.; Steitz, T.A.
Deposited on : 2002-08-23
Resolution : 3.52 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

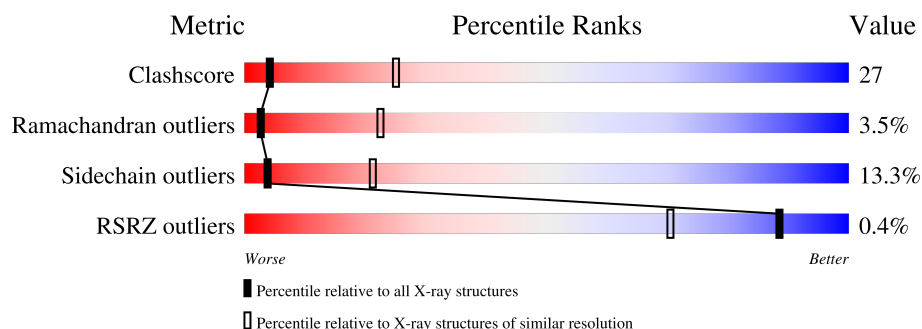
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.52 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	1079 (3.56-3.48)
Ramachandran outliers	187476	1052 (3.56-3.48)
Sidechain outliers	187428	1053 (3.56-3.48)
RSRZ outliers	180081	1024 (3.56-3.48)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	404	
1	B	404	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6320 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.

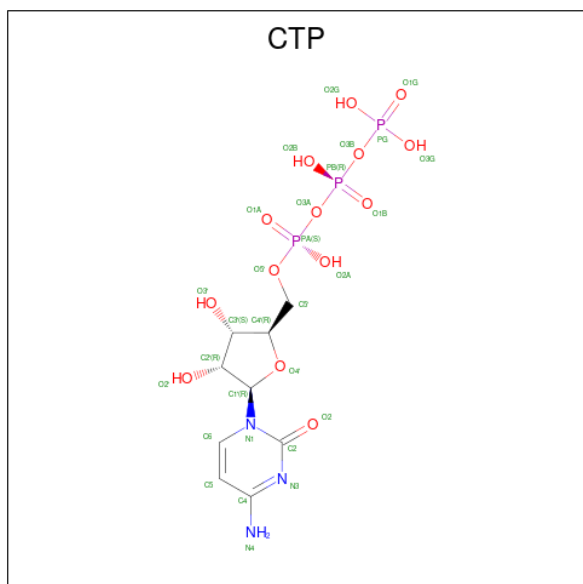
- Molecule 1 is a protein called tRNA CCA-adding enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	0	0	0
			3112	1986	560	554	12			
1	B	395	Total	C	N	O	S	0	0	0
			3112	1986	560	554	12			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mg	0	0
			2	2		
2	B	2	Total	Mg	0	0
			2	2		

- Molecule 3 is CYTIDINE-5'-TRIPHOSPHATE (CCD ID: CTP) (formula: C₉H₁₆N₃O₁₄P₃).



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

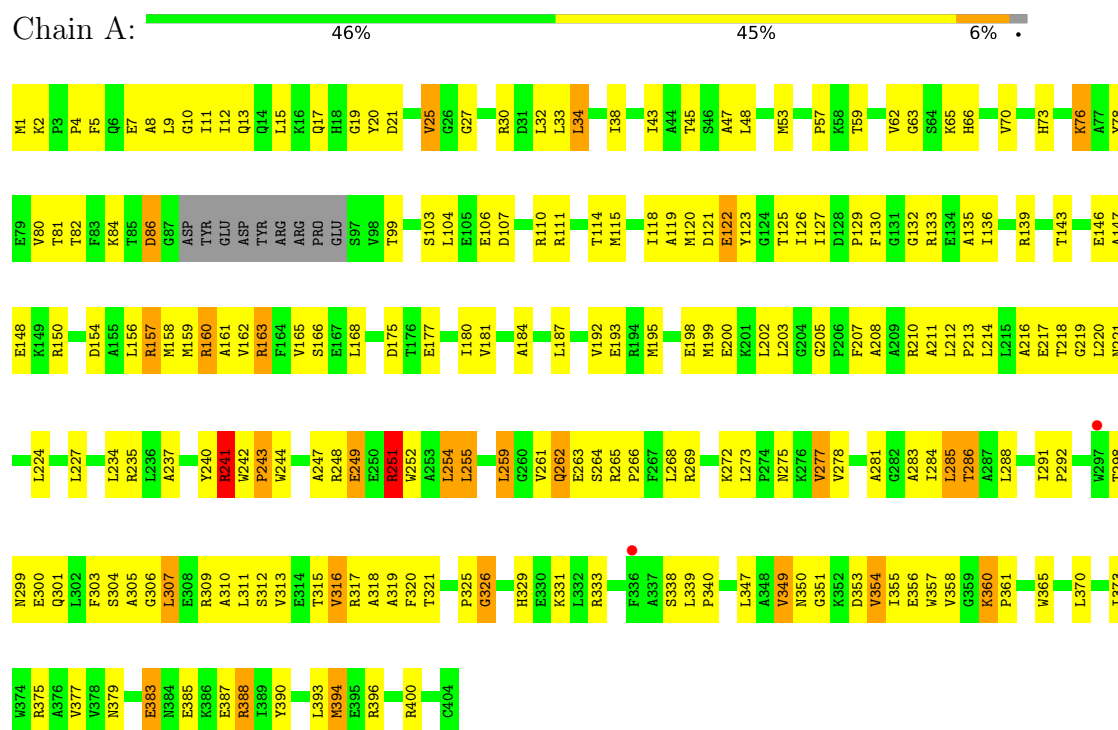
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	17	Total	O	0	0
			17	17		
4	B	17	Total	O	0	0
			17	17		

3 Residue-property plots

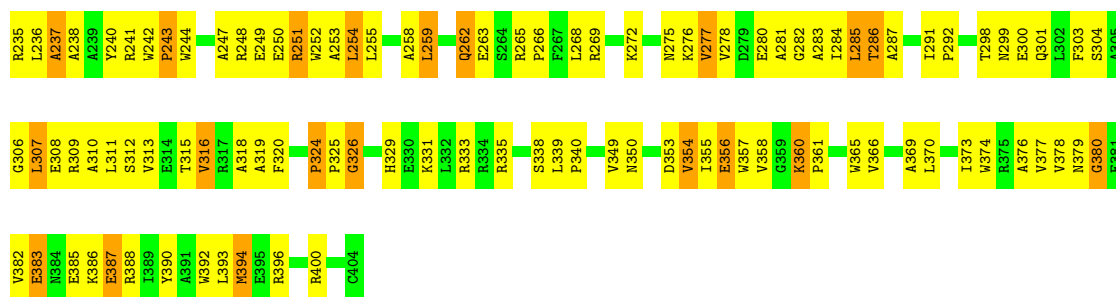
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

• Molecule 1: tRNA CCA-adding enzyme



• Molecule 1: tRNA CCA-adding enzyme





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	105.55Å 105.55Å 182.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	81.65 – 3.52 81.65 – 3.52	Depositor EDS
% Data completeness (in resolution range)	99.5 (81.65-3.52) 93.6 (81.65-3.52)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 3.49Å)	Xtriage
Refinement program	REFMAC 5.1.17	Depositor
R, R_{free}	0.286 , 0.331 0.282 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	157.1	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 137.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.045 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6320	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 31.19 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1528e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.86	0/3180	1.13	6/4305 (0.1%)
1	B	0.79	1/3180 (0.0%)	1.09	9/4305 (0.2%)
All	All	0.83	1/6360 (0.0%)	1.11	15/8610 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	394	MET	SD-CE	5.03	1.92	1.79

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	132	GLY	N-CA-C	-6.92	104.47	112.50
1	B	259	LEU	N-CA-C	-6.00	105.79	113.23
1	A	316	VAL	N-CA-C	-5.65	105.22	110.53
1	A	211	ALA	N-CA-C	-5.45	104.20	112.04
1	B	324	PRO	CA-C-N	5.42	124.94	119.19
1	B	324	PRO	C-N-CA	5.42	124.94	119.19
1	B	132	GLY	N-CA-C	-5.41	106.23	112.50
1	B	380	GLY	CA-C-N	5.28	127.30	120.44
1	B	380	GLY	C-N-CA	5.28	127.30	120.44
1	B	201	LYS	N-CA-C	-5.18	105.80	112.68
1	B	247	ALA	N-CA-C	5.18	117.68	109.50
1	A	247	ALA	CA-C-N	5.15	127.95	120.79
1	A	247	ALA	C-N-CA	5.15	127.95	120.79
1	B	29	VAL	N-CA-C	-5.08	105.90	110.82
1	A	259	LEU	N-CA-C	-5.01	107.16	113.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3112	0	3161	174	0
1	B	3112	0	3161	171	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	29	0	12	2	0
3	B	29	0	12	2	0
4	A	17	0	0	0	0
4	B	17	0	0	0	0
All	All	6320	0	6346	345	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 27.

All (345) 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:B:360:LYS:HG3	1:B:361:PRO:HD2	1.17	1.09
1:A:360:LYS:HG3	1:A:361:PRO:HD2	1.29	1.08
1:A:242:TRP:HB2	1:A:243:PRO:HD3	1.48	0.95
1:B:242:TRP:HB2	1:B:243:PRO:HD3	1.51	0.90
1:A:370:LEU:HA	1:A:373:ILE:HD12	1.52	0.89
1:B:360:LYS:HG3	1:B:361:PRO:CD	2.00	0.89
1:A:298:THR:HB	1:A:301:GLN:HG3	1.54	0.89
1:B:143:THR:HG21	1:B:147:ALA:HA	1.55	0.88
1:B:118:ILE:HG22	1:B:119:ALA:H	1.37	0.88
1:A:312:SER:O	1:A:316:VAL:HG23	1.74	0.87
1:B:249:GLU:H	1:B:249:GLU:CD	1.84	0.86
1:A:118:ILE:HG22	1:A:119:ALA:H	1.41	0.85
1:A:300:GLU:HG2	1:A:379:ASN:HD21	1.39	0.85
1:A:143:THR:HG21	1:A:147:ALA:HA	1.58	0.85
1:A:118:ILE:HG22	1:A:119:ALA:N	1.91	0.85
1:A:249:GLU:CD	1:A:249:GLU:H	1.86	0.84
1:A:34:LEU:HD13	1:A:136:ILE:HG23	1.58	0.84
1:B:318:ALA:HB2	1:B:325:PRO:HG3	1.60	0.83
1:A:318:ALA:HB2	1:A:325:PRO:HG3	1.61	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:370:LEU:HA	1:B:373:ILE:HD12	1.61	0.82
1:A:360:LYS:HG3	1:A:361:PRO:CD	2.10	0.81
1:B:11:ILE:HD11	1:B:73:HIS:HB2	1.62	0.81
1:B:298:THR:HB	1:B:301:GLN:HG3	1.63	0.80
1:B:118:ILE:HG22	1:B:119:ALA:N	1.97	0.79
1:B:15:LEU:O	1:B:20:TYR:HB2	1.84	0.78
1:A:53:MET:HG2	1:A:59:THR:HG21	1.68	0.76
1:A:199:MET:CE	1:A:202:LEU:HD22	2.16	0.76
1:A:241:ARG:HB2	1:A:319:ALA:HA	1.68	0.76
1:B:120:MET:HE3	1:B:126:ILE:HD11	1.68	0.74
1:A:11:ILE:HD11	1:A:73:HIS:HB2	1.68	0.74
1:A:43:ILE:O	1:A:80:VAL:HA	1.86	0.74
1:A:255:LEU:HD12	1:A:255:LEU:O	1.87	0.74
1:A:11:ILE:HD11	1:A:73:HIS:CB	2.18	0.73
1:B:203:LEU:HB3	1:B:251:ARG:HG2	1.70	0.73
1:A:203:LEU:HB3	1:A:251:ARG:HG2	1.71	0.73
1:A:318:ALA:CB	1:A:325:PRO:HG3	2.19	0.72
1:A:242:TRP:HB2	1:A:243:PRO:CD	2.20	0.72
1:A:15:LEU:O	1:A:20:TYR:HB2	1.89	0.71
1:A:298:THR:CB	1:A:301:GLN:HG3	2.19	0.71
1:A:192:VAL:HA	1:A:195:MET:HE3	1.73	0.71
1:B:199:MET:CE	1:B:202:LEU:HD22	2.21	0.70
1:B:11:ILE:HD11	1:B:73:HIS:CB	2.21	0.70
1:A:309:ARG:O	1:A:313:VAL:HG23	1.92	0.69
1:A:118:ILE:CG2	1:A:119:ALA:H	2.06	0.69
1:B:118:ILE:CG2	1:B:119:ALA:H	2.05	0.68
1:A:111:ARG:HB3	3:A:5501:CTP:O2'	1.94	0.67
1:B:300:GLU:HG2	1:B:379:ASN:HD21	1.59	0.67
1:B:318:ALA:CB	1:B:325:PRO:HG3	2.24	0.67
1:A:310:ALA:O	1:A:313:VAL:HB	1.95	0.67
1:A:358:VAL:HG12	1:A:358:VAL:O	1.95	0.67
1:B:242:TRP:HB2	1:B:243:PRO:CD	2.23	0.66
1:A:217:GLU:HG3	1:A:235:ARG:HH21	1.61	0.66
1:A:311:LEU:HD23	1:A:329:HIS:HA	1.78	0.66
1:A:120:MET:HE3	1:A:126:ILE:HD11	1.78	0.65
1:A:154:ASP:HB3	1:A:157:ARG:HD2	1.78	0.65
1:A:118:ILE:CG2	1:A:119:ALA:N	2.60	0.65
1:B:53:MET:HG2	1:B:59:THR:HG21	1.79	0.65
1:B:298:THR:O	1:B:301:GLN:N	2.30	0.65
1:B:34:LEU:HD13	1:B:136:ILE:HG23	1.79	0.64
1:A:315:THR:OG1	1:A:329:HIS:CE1	2.50	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:ASP:O	1:A:123:TYR:N	2.31	0.64
1:B:212:LEU:HB2	1:B:213:PRO:HD3	1.79	0.64
1:A:354:VAL:HG13	1:A:365:TRP:CH2	2.33	0.64
1:B:53:MET:HA	1:B:59:THR:HG21	1.80	0.64
1:B:358:VAL:HG12	1:B:358:VAL:O	1.98	0.64
1:A:217:GLU:CG	1:A:235:ARG:HH21	2.11	0.63
1:A:212:LEU:HB2	1:A:213:PRO:HD3	1.81	0.63
1:A:53:MET:HA	1:A:59:THR:HG21	1.81	0.62
1:B:217:GLU:HG3	1:B:235:ARG:HH21	1.64	0.62
1:A:34:LEU:CD1	1:A:136:ILE:HG23	2.30	0.62
1:B:241:ARG:HB2	1:B:319:ALA:HA	1.82	0.62
1:A:177:GLU:HA	1:A:180:ILE:HD12	1.81	0.62
1:B:248:ARG:HB3	1:B:249:GLU:OE2	1.99	0.62
1:B:309:ARG:O	1:B:313:VAL:HG23	1.99	0.62
1:B:350:ASN:HB3	1:B:353:ASP:OD2	1.99	0.62
1:A:48:LEU:HD23	1:A:82:THR:HG21	1.82	0.62
1:B:154:ASP:HB3	1:B:157:ARG:HD2	1.80	0.62
1:B:356:GLU:HG3	1:B:356:GLU:O	2.00	0.61
1:B:315:THR:OG1	1:B:329:HIS:CE1	2.53	0.61
1:B:43:ILE:O	1:B:80:VAL:HA	2.01	0.61
1:A:121:ASP:C	1:A:123:TYR:H	2.08	0.61
1:B:1:MET:HE3	1:B:9:LEU:HD22	1.82	0.60
1:A:187:LEU:HD11	1:A:220:LEU:HD22	1.83	0.60
1:A:355:ILE:C	1:A:357:TRP:H	2.08	0.60
1:A:161:ALA:O	1:A:165:VAL:HG23	2.01	0.60
1:B:177:GLU:HA	1:B:180:ILE:HD12	1.83	0.60
1:B:234:LEU:O	1:B:237:ALA:HB3	2.02	0.59
1:B:312:SER:O	1:B:316:VAL:HG23	2.03	0.59
1:A:350:ASN:HB3	1:A:353:ASP:OD2	2.02	0.59
1:A:1:MET:HE3	1:A:9:LEU:HD22	1.85	0.58
1:A:199:MET:HE1	1:A:202:LEU:HD22	1.86	0.58
1:B:192:VAL:HA	1:B:195:MET:HE3	1.85	0.58
1:B:112:ASP:OD2	1:B:160:ARG:NH2	2.29	0.57
1:B:277:VAL:HG12	1:B:278:VAL:N	2.18	0.57
1:A:252:TRP:CE3	1:A:255:LEU:HD23	2.40	0.57
1:A:298:THR:O	1:A:301:GLN:N	2.37	0.57
1:B:158:MET:O	1:B:159:MET:C	2.48	0.56
1:B:298:THR:CB	1:B:301:GLN:HG3	2.35	0.56
1:B:354:VAL:HG22	1:B:393:LEU:CD2	2.34	0.56
1:A:300:GLU:CG	1:A:379:ASN:HD21	2.15	0.56
1:A:354:VAL:HG22	1:A:393:LEU:CD2	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:ILE:CG2	1:B:119:ALA:N	2.66	0.56
1:B:48:LEU:HD23	1:B:82:THR:HG21	1.88	0.56
1:B:315:THR:OG1	1:B:329:HIS:NE2	2.38	0.56
1:B:254:LEU:HD13	1:B:319:ALA:HB3	1.86	0.56
1:B:217:GLU:CG	1:B:235:ARG:HH21	2.19	0.56
1:B:390:TYR:CE1	1:B:394:MET:HG2	2.41	0.55
1:B:25:VAL:HG21	1:B:111:ARG:HH22	1.71	0.55
1:A:240:TYR:CE2	1:A:320:PHE:HD1	2.24	0.55
1:A:248:ARG:HB3	1:A:249:GLU:OE2	2.06	0.55
1:A:242:TRP:CB	1:A:243:PRO:HD3	2.30	0.55
1:B:224:LEU:HB2	1:B:227:LEU:HD12	1.88	0.54
1:A:103:SER:HB3	1:A:106:GLU:HB2	1.88	0.54
1:B:286:THR:HG22	1:B:287:ALA:N	2.21	0.54
1:B:252:TRP:CE3	1:B:255:LEU:HD23	2.43	0.54
1:A:298:THR:O	1:A:299:ASN:C	2.51	0.53
1:B:283:ALA:HB1	1:B:309:ARG:CZ	2.39	0.53
1:A:216:ALA:O	1:A:218:THR:O	2.26	0.53
1:B:355:ILE:C	1:B:357:TRP:H	2.15	0.53
1:B:383:GLU:HG3	1:B:388:ARG:HD2	1.89	0.53
1:B:283:ALA:HB1	1:B:309:ARG:NE	2.21	0.53
1:A:254:LEU:HD13	1:A:319:ALA:HB3	1.91	0.53
1:B:199:MET:HE1	1:B:202:LEU:HD22	1.90	0.53
1:B:268:LEU:HB2	1:B:278:VAL:HG13	1.91	0.53
1:A:277:VAL:O	1:A:281:ALA:N	2.38	0.53
1:A:303:PHE:C	1:A:303:PHE:CD2	2.87	0.53
1:B:221:ASN:ND2	1:B:227:LEU:HB3	2.23	0.53
1:B:298:THR:O	1:B:299:ASN:C	2.52	0.53
1:A:25:VAL:HG21	1:A:111:ARG:HH22	1.74	0.53
1:B:11:ILE:HB	1:B:43:ILE:HD13	1.90	0.52
1:B:265:ARG:O	1:B:269:ARG:HB2	2.09	0.52
1:A:115:MET:HE1	1:A:168:LEU:CD1	2.39	0.52
1:B:111:ARG:HB3	3:B:6501:CTP:O2'	2.10	0.52
1:B:307:LEU:HD21	1:B:333:ARG:HG2	1.92	0.52
1:A:306:GLY:O	1:A:307:LEU:C	2.52	0.52
1:B:299:ASN:N	1:B:335:ARG:NH2	2.57	0.52
1:A:162:VAL:HG13	1:A:214:LEU:HD23	1.91	0.52
1:A:163:ARG:NH2	3:A:5501:CTP:O3G	2.42	0.52
1:A:275:ASN:O	1:A:277:VAL:N	2.43	0.52
1:B:299:ASN:N	1:B:335:ARG:HH22	2.08	0.52
1:B:62:VAL:HG21	1:B:70:VAL:HG23	1.91	0.52
1:B:121:ASP:O	1:B:123:TYR:N	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:ASN:O	1:A:353:ASP:HB2	2.10	0.52
1:B:163:ARG:HD3	1:B:198:GLU:OE2	2.10	0.52
1:B:184:ALA:HB2	1:B:219:GLY:HA3	1.92	0.52
1:A:221:ASN:ND2	1:A:227:LEU:HB3	2.25	0.51
1:A:263:GLU:O	1:A:266:PRO:HD2	2.10	0.51
1:B:240:TYR:CE2	1:B:320:PHE:HD1	2.28	0.51
1:B:121:ASP:C	1:B:123:TYR:H	2.18	0.51
1:A:268:LEU:HB2	1:A:278:VAL:HG13	1.90	0.51
1:A:20:TYR:CD2	1:A:47:ALA:HB2	2.46	0.51
1:B:284:ILE:HG23	1:B:313:VAL:HG22	1.92	0.51
1:B:350:ASN:O	1:B:353:ASP:HB2	2.11	0.51
1:A:86:ASP:OD2	1:A:110:ARG:HD2	2.11	0.51
1:B:303:PHE:C	1:B:303:PHE:CD2	2.89	0.51
1:B:45:THR:O	1:B:82:THR:HA	2.10	0.50
1:B:236:LEU:C	1:B:238:ALA:H	2.19	0.50
1:A:217:GLU:HG3	1:A:235:ARG:NH2	2.25	0.50
1:B:236:LEU:O	1:B:238:ALA:N	2.45	0.50
1:A:259:LEU:HB2	1:A:261:VAL:HG23	1.93	0.50
1:A:224:LEU:HB2	1:A:227:LEU:HD12	1.93	0.50
1:A:255:LEU:HD12	1:A:255:LEU:C	2.35	0.50
1:B:236:LEU:C	1:B:238:ALA:N	2.68	0.50
1:B:306:GLY:O	1:B:307:LEU:C	2.53	0.50
1:B:146:GLU:HG3	1:B:148:GLU:HB2	1.93	0.50
1:B:161:ALA:O	1:B:165:VAL:HG23	2.12	0.49
1:B:212:LEU:CD1	1:B:242:TRP:CZ2	2.95	0.49
1:B:354:VAL:HG22	1:B:393:LEU:HD21	1.94	0.49
1:A:32:LEU:HD21	1:A:120:MET:HE1	1.94	0.49
1:A:248:ARG:HD3	1:A:277:VAL:HG21	1.94	0.49
1:A:33:LEU:O	1:A:34:LEU:C	2.54	0.49
1:B:282:GLY:O	1:B:283:ALA:C	2.55	0.49
1:B:187:LEU:HD11	1:B:220:LEU:HD22	1.94	0.49
1:B:181:VAL:HG22	1:B:218:THR:HA	1.93	0.49
1:A:383:GLU:HG3	1:A:388:ARG:HD2	1.93	0.49
1:A:207:PHE:O	1:A:208:ALA:C	2.56	0.49
1:A:355:ILE:C	1:A:357:TRP:N	2.70	0.49
1:B:103:SER:HB3	1:B:106:GLU:HB2	1.93	0.49
1:B:107:ASP:O	1:B:110:ARG:HG2	2.12	0.48
1:A:10:GLY:O	1:A:11:ILE:C	2.56	0.48
1:A:163:ARG:HD3	1:A:198:GLU:OE2	2.13	0.48
1:A:158:MET:O	1:A:159:MET:C	2.55	0.48
1:A:373:ILE:O	1:A:377:VAL:HG23	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:TYR:HD2	1:A:254:LEU:HD11	1.78	0.48
1:A:20:TYR:CG	1:A:47:ALA:HB2	2.49	0.48
1:B:354:VAL:HG13	1:B:365:TRP:CH2	2.49	0.48
1:A:283:ALA:HB1	1:A:309:ARG:NE	2.28	0.48
1:A:315:THR:OG1	1:A:329:HIS:NE2	2.45	0.48
1:A:242:TRP:CB	1:A:243:PRO:CD	2.87	0.48
1:B:284:ILE:O	1:B:285:LEU:C	2.56	0.48
1:A:284:ILE:O	1:A:285:LEU:C	2.57	0.48
1:B:280:GLU:O	1:B:283:ALA:HB3	2.14	0.48
1:A:192:VAL:O	1:A:193:GLU:C	2.57	0.47
1:A:268:LEU:HD11	1:A:281:ALA:HB1	1.97	0.47
1:B:143:THR:HG21	1:B:147:ALA:CA	2.37	0.47
1:B:373:ILE:O	1:B:377:VAL:HG23	2.13	0.47
1:A:350:ASN:OD1	1:A:351:GLY:N	2.47	0.47
1:B:263:GLU:O	1:B:266:PRO:HD2	2.14	0.47
1:A:76:LYS:HG3	1:A:78:TYR:OH	2.14	0.47
1:A:158:MET:SD	1:A:180:ILE:HA	2.55	0.47
1:B:27:GLY:HA2	1:B:30:ARG:HB3	1.96	0.47
1:B:277:VAL:O	1:B:281:ALA:N	2.40	0.47
1:B:311:LEU:HD23	1:B:329:HIS:HA	1.97	0.47
1:A:268:LEU:CD1	1:A:281:ALA:CB	2.92	0.47
1:B:8:ALA:HB2	1:B:78:TYR:CE2	2.49	0.47
1:B:76:LYS:HG3	1:B:78:TYR:OH	2.15	0.47
1:B:242:TRP:CB	1:B:243:PRO:HD3	2.34	0.47
1:A:146:GLU:HG3	1:A:148:GLU:HB2	1.96	0.47
1:B:242:TRP:CB	1:B:243:PRO:CD	2.92	0.46
1:A:298:THR:O	1:A:300:GLU:N	2.48	0.46
1:A:390:TYR:CE1	1:A:394:MET:HG2	2.49	0.46
1:B:62:VAL:HG21	1:B:70:VAL:CG2	2.46	0.46
1:A:184:ALA:HB2	1:A:219:GLY:HA3	1.98	0.46
1:B:244:TRP:CG	1:B:244:TRP:O	2.68	0.46
1:B:248:ARG:HD3	1:B:277:VAL:HG21	1.98	0.46
1:A:4:PRO:HG2	1:A:38:ILE:HD12	1.98	0.46
1:A:53:MET:HG2	1:A:59:THR:CG2	2.43	0.46
1:A:265:ARG:O	1:A:269:ARG:HB2	2.15	0.46
1:B:240:TYR:CZ	1:B:258:ALA:HB2	2.50	0.46
1:B:382:VAL:HG21	1:B:392:TRP:HB2	1.97	0.45
1:A:120:MET:HG3	1:A:126:ILE:HD13	1.98	0.45
1:A:214:LEU:C	1:A:216:ALA:N	2.72	0.45
1:B:5:PHE:HE1	1:B:38:ILE:HD13	1.82	0.45
1:A:107:ASP:O	1:A:110:ARG:HG2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:GLN:C	1:B:19:GLY:H	2.24	0.45
1:B:157:ARG:HA	1:B:160:ARG:HD3	1.97	0.45
1:A:187:LEU:CD1	1:A:220:LEU:HD22	2.46	0.45
1:B:275:ASN:O	1:B:277:VAL:N	2.50	0.45
1:B:298:THR:O	1:B:300:GLU:N	2.50	0.45
1:A:2:LYS:C	1:A:4:PRO:HD2	2.42	0.45
1:A:121:ASP:HB3	1:A:127:ILE:HD11	1.99	0.45
1:B:251:ARG:O	1:B:252:TRP:C	2.59	0.45
1:A:192:VAL:HA	1:A:195:MET:CE	2.45	0.45
1:A:307:LEU:HD21	1:A:333:ARG:HG2	1.99	0.45
1:B:163:ARG:HG2	1:B:164:PHE:N	2.30	0.45
1:B:324:PRO:HA	1:B:325:PRO:HD3	1.68	0.45
1:B:27:GLY:O	1:B:31:ASP:N	2.41	0.44
1:A:234:LEU:O	1:A:237:ALA:HB3	2.17	0.44
1:A:244:TRP:CD1	1:A:315:THR:HG23	2.53	0.44
1:A:298:THR:OG1	1:A:301:GLN:NE2	2.49	0.44
1:B:12:ILE:HG23	1:B:22:ALA:HB3	1.99	0.44
1:A:354:VAL:HG22	1:A:393:LEU:HD21	1.99	0.44
1:A:156:LEU:C	1:A:158:MET:N	2.75	0.44
1:B:193:GLU:OE2	1:B:272:LYS:HD3	2.16	0.44
1:A:17:GLN:C	1:A:19:GLY:H	2.25	0.44
1:A:251:ARG:O	1:A:252:TRP:C	2.59	0.44
1:A:393:LEU:O	1:A:394:MET:C	2.60	0.44
1:A:2:LYS:HB3	1:A:4:PRO:HD2	2.00	0.44
1:A:193:GLU:OE2	1:A:272:LYS:HD3	2.18	0.44
1:A:252:TRP:CE2	1:A:273:LEU:HD11	2.53	0.44
1:B:192:VAL:O	1:B:193:GLU:C	2.61	0.43
1:A:156:LEU:O	1:A:158:MET:N	2.51	0.43
1:A:181:VAL:HG22	1:A:218:THR:HA	1.99	0.43
1:B:291:ILE:HG12	1:B:301:GLN:HE22	1.83	0.43
1:A:143:THR:HG21	1:A:147:ALA:CA	2.39	0.43
1:A:163:ARG:HH11	1:A:198:GLU:CD	2.25	0.43
1:A:284:ILE:HG23	1:A:313:VAL:HG22	1.99	0.43
1:B:240:TYR:HD2	1:B:254:LEU:HD11	1.82	0.43
1:B:121:ASP:C	1:B:123:TYR:N	2.76	0.43
1:B:252:TRP:CZ3	1:B:255:LEU:HD23	2.54	0.43
1:B:355:ILE:C	1:B:357:TRP:N	2.76	0.43
1:A:159:MET:HE3	1:A:198:GLU:HB2	2.00	0.43
1:A:166:SER:OG	1:A:205:GLY:HA3	2.18	0.43
1:A:249:GLU:CD	1:A:249:GLU:N	2.65	0.43
1:B:291:ILE:HA	1:B:292:PRO:HD3	1.85	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:315:THR:HG1	1:B:329:HIS:CE1	2.35	0.43
1:A:157:ARG:HA	1:A:160:ARG:HD3	2.01	0.42
1:B:20:TYR:CD2	1:B:47:ALA:HB2	2.54	0.42
1:B:233:GLN:NE2	1:B:259:LEU:O	2.52	0.42
1:B:240:TYR:HB3	1:B:254:LEU:HD21	2.00	0.42
1:B:325:PRO:O	1:B:326:GLY:C	2.62	0.42
1:A:27:GLY:O	1:A:30:ARG:HB3	2.19	0.42
1:A:240:TYR:CE2	1:A:320:PHE:CD1	3.07	0.42
1:A:317:ARG:O	1:A:321:THR:HG23	2.20	0.42
1:A:390:TYR:CZ	1:A:394:MET:HG2	2.54	0.42
1:B:158:MET:SD	1:B:180:ILE:HA	2.60	0.42
1:B:163:ARG:NH1	1:B:198:GLU:OE1	2.51	0.42
1:B:307:LEU:O	1:B:308:GLU:C	2.60	0.42
1:B:360:LYS:CG	1:B:361:PRO:HD2	2.12	0.42
1:A:2:LYS:HG3	1:A:32:LEU:HD13	2.01	0.42
1:A:62:VAL:HG21	1:A:70:VAL:HG23	2.01	0.42
1:A:286:THR:C	1:A:288:LEU:N	2.76	0.42
1:B:255:LEU:HD12	1:B:255:LEU:O	2.18	0.42
1:A:351:GLY:HA2	1:A:370:LEU:HD11	2.00	0.42
1:B:63:GLY:HA2	1:B:65:LYS:HE2	2.02	0.42
1:A:5:PHE:O	1:A:9:LEU:N	2.53	0.42
1:B:268:LEU:CD1	1:B:281:ALA:CB	2.98	0.42
1:B:366:VAL:O	1:B:369:ALA:HB3	2.19	0.42
1:A:305:ALA:HA	1:A:309:ARG:NH1	2.35	0.42
1:B:163:ARG:HH11	1:B:198:GLU:CD	2.28	0.42
1:B:374:TRP:C	1:B:376:ALA:H	2.26	0.42
1:A:254:LEU:HD13	1:A:319:ALA:CB	2.50	0.42
1:A:347:LEU:C	1:A:349:VAL:H	2.27	0.42
1:B:115:MET:HE1	1:B:168:LEU:CD1	2.49	0.42
1:B:339:LEU:HA	1:B:340:PRO:HD3	1.81	0.42
1:A:121:ASP:C	1:A:123:TYR:N	2.68	0.42
1:A:162:VAL:HA	1:A:165:VAL:HG23	2.02	0.42
1:A:240:TYR:HB3	1:A:254:LEU:HD21	2.01	0.42
1:B:20:TYR:CG	1:B:47:ALA:HB2	2.55	0.42
1:B:217:GLU:HG3	1:B:235:ARG:NH2	2.32	0.42
1:A:339:LEU:HA	1:A:340:PRO:HD3	1.85	0.42
1:B:113:PHE:HB3	1:B:142:ARG:O	2.20	0.42
1:B:156:LEU:HB2	1:B:190:ILE:HG21	2.01	0.42
1:B:310:ALA:O	1:B:313:VAL:HB	2.20	0.42
1:A:199:MET:HE3	1:A:202:LEU:HD22	2.00	0.41
1:A:277:VAL:HG12	1:A:278:VAL:N	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:ARG:NH2	3:B:6501:CTP:O3G	2.53	0.41
1:B:387:GLU:O	1:B:390:TYR:HB3	2.20	0.41
1:A:244:TRP:O	1:A:244:TRP:CG	2.72	0.41
1:A:360:LYS:CG	1:A:361:PRO:HD2	2.22	0.41
1:B:216:ALA:O	1:B:218:THR:O	2.38	0.41
1:A:11:ILE:CD1	1:A:73:HIS:HB2	2.44	0.41
1:A:45:THR:HG23	1:A:47:ALA:H	1.86	0.41
1:A:45:THR:O	1:A:82:THR:HA	2.19	0.41
1:B:214:LEU:C	1:B:216:ALA:N	2.76	0.41
1:A:135:ALA:O	1:A:139:ARG:N	2.54	0.41
1:A:160:ARG:NH1	1:A:163:ARG:NH1	2.68	0.41
1:A:325:PRO:O	1:A:326:GLY:C	2.62	0.41
1:B:86:ASP:OD2	1:B:110:ARG:HD2	2.20	0.41
1:B:285:LEU:HD23	1:B:285:LEU:HA	1.94	0.41
1:A:291:ILE:HA	1:A:292:PRO:HD3	1.87	0.41
1:B:2:LYS:C	1:B:4:PRO:HD2	2.46	0.41
1:B:233:GLN:H	1:B:233:GLN:HG2	1.73	0.41
1:A:8:ALA:HB2	1:A:78:TYR:CE2	2.55	0.41
1:B:386:LYS:O	1:B:387:GLU:C	2.61	0.41
1:A:216:ALA:HA	1:A:221:ASN:HB2	2.03	0.41
1:A:129:PRO:C	1:A:130:PHE:CD1	2.98	0.41
1:A:220:LEU:HD23	1:A:220:LEU:HA	1.86	0.41
1:A:268:LEU:CD1	1:A:281:ALA:HB1	2.51	0.41
1:A:305:ALA:HB1	1:A:309:ARG:HD2	2.03	0.41
1:B:5:PHE:O	1:B:9:LEU:N	2.54	0.41
1:B:120:MET:HE3	1:B:126:ILE:CD1	2.44	0.41
1:B:184:ALA:N	1:B:185:PRO:CD	2.84	0.41
1:B:268:LEU:HD11	1:B:281:ALA:HB1	2.02	0.41
1:A:156:LEU:C	1:A:158:MET:H	2.29	0.40
1:A:355:ILE:O	1:A:357:TRP:N	2.53	0.40
1:B:378:VAL:C	1:B:380:GLY:H	2.29	0.40
1:A:275:ASN:C	1:A:277:VAL:N	2.77	0.40
1:B:265:ARG:HB3	1:B:266:PRO:HD3	2.03	0.40
1:B:212:LEU:O	1:B:213:PRO:C	2.62	0.40
1:B:250:GLU:O	1:B:253:ALA:HB3	2.21	0.40
1:B:207:PHE:O	1:B:208:ALA:C	2.64	0.40
1:B:216:ALA:HA	1:B:221:ASN:HB2	2.04	0.40
1:B:244:TRP:CD1	1:B:315:THR:HG23	2.55	0.40
1:B:374:TRP:C	1:B:376:ALA:N	2.80	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	391/404 (97%)	316 (81%)	61 (16%)	14 (4%)	2	21
1	B	391/404 (97%)	319 (82%)	59 (15%)	13 (3%)	3	23
All	All	782/808 (97%)	635 (81%)	120 (15%)	27 (4%)	3	22

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	307	LEU
1	B	383	GLU
1	A	326	GLY
1	A	383	GLU
1	B	122	GLU
1	B	307	LEU
1	B	326	GLY
1	A	122	GLU
1	A	262	GLN
1	B	34	LEU
1	B	57	PRO
1	B	262	GLN
1	B	356	GLU
1	A	57	PRO
1	A	86	ASP
1	A	241	ARG
1	A	243	PRO
1	A	251	ARG
1	A	356	GLU
1	B	86	ASP
1	B	237	ALA
1	A	21	ASP
1	A	34	LEU
1	B	276	LYS
1	B	243	PRO

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Mol	Chain	Res	Type
1	A	63	GLY
1	B	63	GLY

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	316/325 (97%)	271 (86%)	45 (14%)	3	19
1	B	316/325 (97%)	277 (88%)	39 (12%)	4	23
All	All	632/650 (97%)	548 (87%)	84 (13%)	4	21

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	GLU
1	A	12	ILE
1	A	13	GLN
1	A	25	VAL
1	A	65	LYS
1	A	66	HIS
1	A	76	LYS
1	A	81	THR
1	A	84	LYS
1	A	99	THR
1	A	104	LEU
1	A	114	THR
1	A	122	GLU
1	A	125	THR
1	A	133	ARG
1	A	150	ARG
1	A	157	ARG
1	A	160	ARG
1	A	163	ARG
1	A	175	ASP
1	A	200	GLU

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Mol	Chain	Res	Type
1	A	210	ARG
1	A	241	ARG
1	A	249	GLU
1	A	251	ARG
1	A	254	LEU
1	A	255	LEU
1	A	262	GLN
1	A	264	SER
1	A	277	VAL
1	A	285	LEU
1	A	286	THR
1	A	304	SER
1	A	331	LYS
1	A	338	SER
1	A	349	VAL
1	A	354	VAL
1	A	360	LYS
1	A	375	ARG
1	A	385	GLU
1	A	387	GLU
1	A	388	ARG
1	A	394	MET
1	A	396	ARG
1	A	400	ARG
1	B	7	GLU
1	B	12	ILE
1	B	13	GLN
1	B	25	VAL
1	B	65	LYS
1	B	66	HIS
1	B	76	LYS
1	B	79	GLU
1	B	84	LYS
1	B	105	GLU
1	B	114	THR
1	B	122	GLU
1	B	126	ILE
1	B	133	ARG
1	B	150	ARG
1	B	157	ARG
1	B	160	ARG
1	B	163	ARG

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Mol	Chain	Res	Type
1	B	175	ASP
1	B	187	LEU
1	B	200	GLU
1	B	210	ARG
1	B	251	ARG
1	B	254	LEU
1	B	262	GLN
1	B	277	VAL
1	B	285	LEU
1	B	286	THR
1	B	304	SER
1	B	316	VAL
1	B	331	LYS
1	B	338	SER
1	B	349	VAL
1	B	354	VAL
1	B	360	LYS
1	B	385	GLU
1	B	387	GLU
1	B	396	ARG
1	B	400	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	73	HIS
1	A	182	GLN
1	A	221	ASN
1	A	233	GLN
1	A	379	ASN
1	B	6	GLN
1	B	182	GLN
1	B	221	ASN
1	B	233	GLN
1	B	257	HIS
1	B	301	GLN
1	B	379	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	CTP	A	5501	2	29,30,30	1.80	5 (17%)	43,47,47	1.25	4 (9%)
3	CTP	B	6501	2	29,30,30	1.97	4 (13%)	43,47,47	1.30	6 (13%)

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	CTP	A	5501	2	-	8/22/38/38	0/2/2/2
3	CTP	B	6501	2	-	7/22/38/38	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	6501	CTP	PB-O3A	6.23	1.66	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	6501	CTP	PA-O3A	5.54	1.65	1.59
3	A	5501	CTP	PB-O3B	4.86	1.64	1.59
3	A	5501	CTP	PA-O3A	4.55	1.64	1.59
3	A	5501	CTP	PB-O3A	4.01	1.63	1.59
3	B	6501	CTP	PB-O3B	3.40	1.63	1.59
3	A	5501	CTP	PG-O1G	3.38	1.61	1.50
3	B	6501	CTP	PG-O1G	3.28	1.60	1.50
3	A	5501	CTP	C6-C5	2.47	1.40	1.35

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	6501	CTP	O2-C2-N3	-3.27	117.17	122.33
3	A	5501	CTP	O3G-PG-O3B	3.18	115.30	104.64
3	A	5501	CTP	O2B-PB-O3B	3.07	115.58	107.27
3	B	6501	CTP	O3G-PG-O3B	3.04	114.83	104.64
3	B	6501	CTP	O4'-C1'-N1	2.66	114.38	108.36
3	B	6501	CTP	O4'-C4'-C3'	2.33	109.79	105.15
3	A	5501	CTP	O4'-C1'-N1	2.13	113.19	108.36
3	B	6501	CTP	O2B-PB-O3A	2.13	113.04	107.27
3	A	5501	CTP	O2A-PA-O3A	2.13	113.02	107.27
3	B	6501	CTP	O2B-PB-O3B	2.05	112.82	107.27

There are no chirality outliers.

All (15) torsion outliers are listed below:

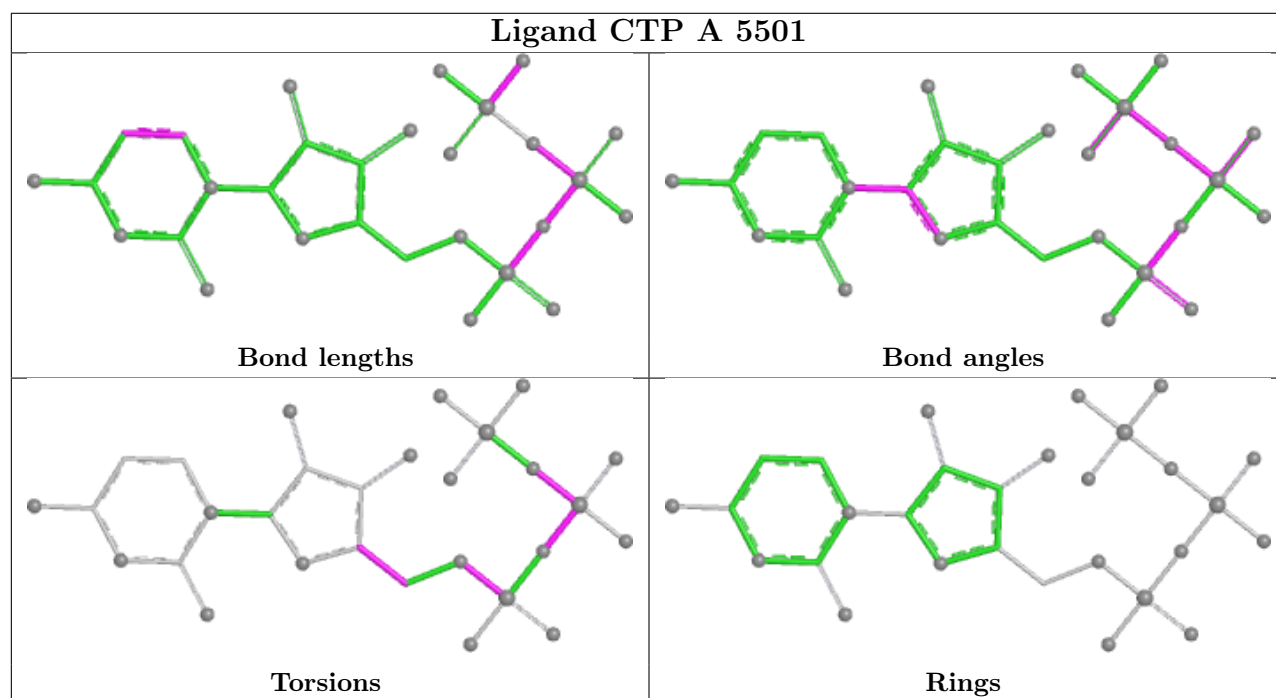
Mol	Chain	Res	Type	Atoms
3	B	6501	CTP	C5'-O5'-PA-O1A
3	B	6501	CTP	C5'-O5'-PA-O3A
3	A	5501	CTP	PG-O3B-PB-O3A
3	B	6501	CTP	PG-O3B-PB-O3A
3	A	5501	CTP	O4'-C4'-C5'-O5'
3	B	6501	CTP	O4'-C4'-C5'-O5'
3	A	5501	CTP	C5'-O5'-PA-O1A
3	A	5501	CTP	C5'-O5'-PA-O3A
3	B	6501	CTP	C3'-C4'-C5'-O5'
3	A	5501	CTP	PA-O3A-PB-O2B
3	A	5501	CTP	PG-O3B-PB-O1B
3	B	6501	CTP	PA-O3A-PB-O1B
3	B	6501	CTP	PG-O3B-PB-O1B
3	A	5501	CTP	C3'-C4'-C5'-O5'
3	A	5501	CTP	PG-O3B-PB-O2B

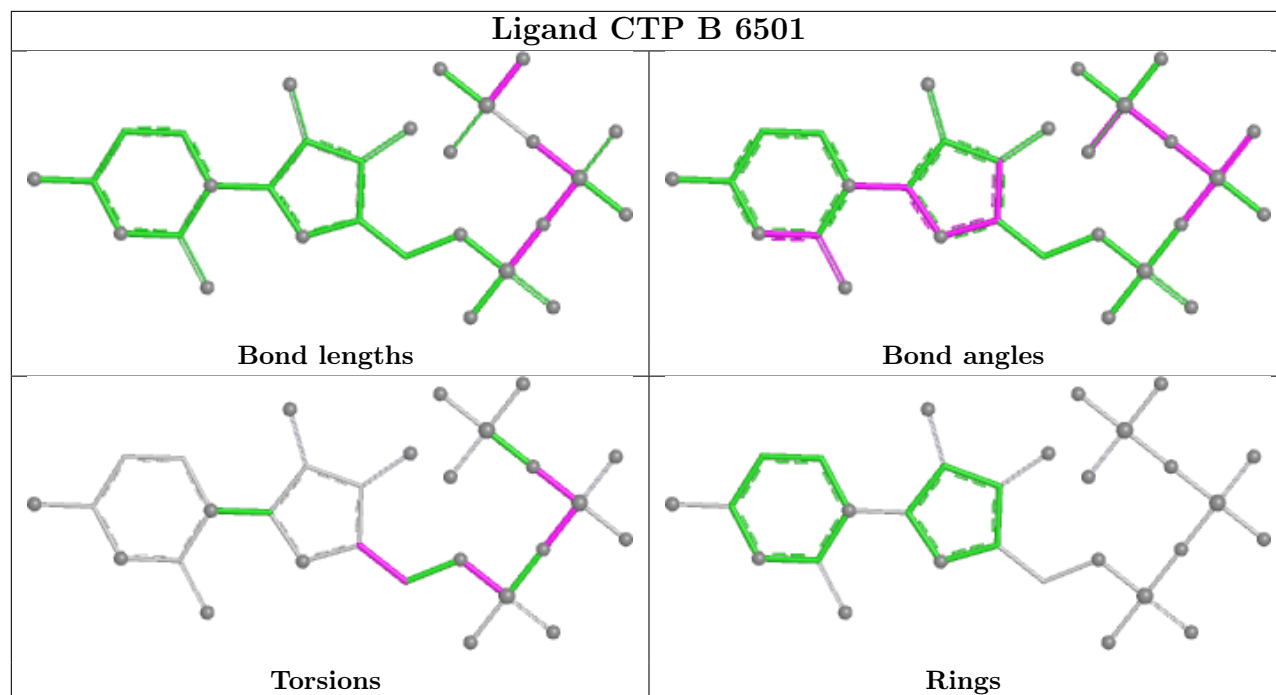
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	5501	CTP	2	0
3	B	6501	CTP	2	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	395/404 (97%)	-0.72	2 (0%) 87 67	6, 33, 73, 85	0
1	B	395/404 (97%)	-0.68	1 (0%) 90 75	6, 33, 73, 85	0
All	All	790/808 (97%)	-0.70	3 (0%) 88 70	6, 33, 73, 85	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	161	ALA	3.0
1	A	297	TRP	2.1
1	A	336	PHE	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	A	5611	1/1	0.51	0.16	9,9,9,9	0
2	MG	B	6611	1/1	0.65	0.18	9,9,9,9	0

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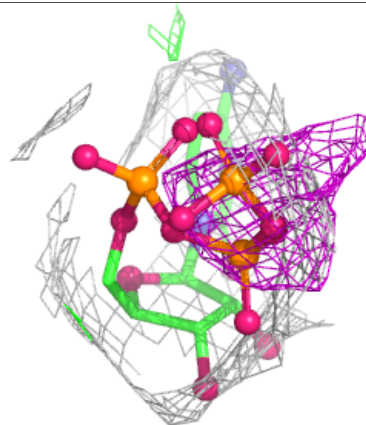
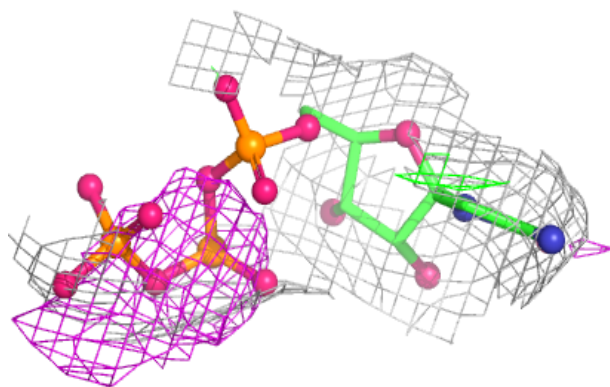
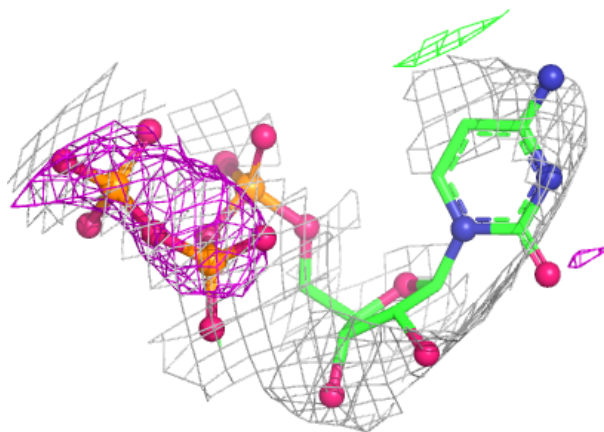
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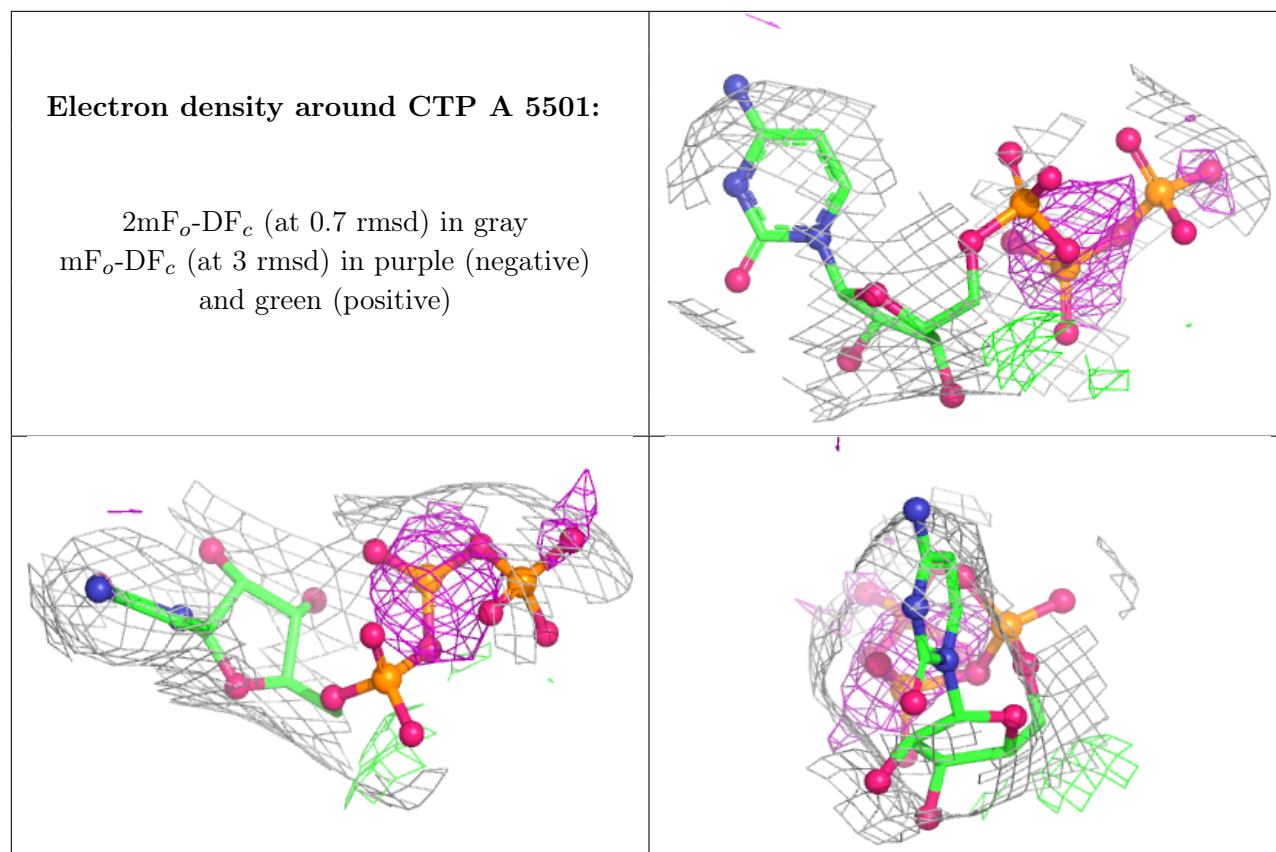
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	B	6601	1/1	0.77	0.13	9,9,9,9	0
2	MG	A	5601	1/1	0.87	0.08	9,9,9,9	0
3	CTP	B	6501	29/29	0.88	0.07	19,25,29,29	0
3	CTP	A	5501	29/29	0.92	0.06	19,25,29,29	0

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 B 6501:

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.