



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:16 AM UTC

PDB ID : 7PZA / pdb_00007pza
Title : Structure of the Clr-cAMP-DNA complex
Authors : Werel, L.; Essen, L.-O.
Deposited on : 2021-10-11
Resolution : 2.72 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:

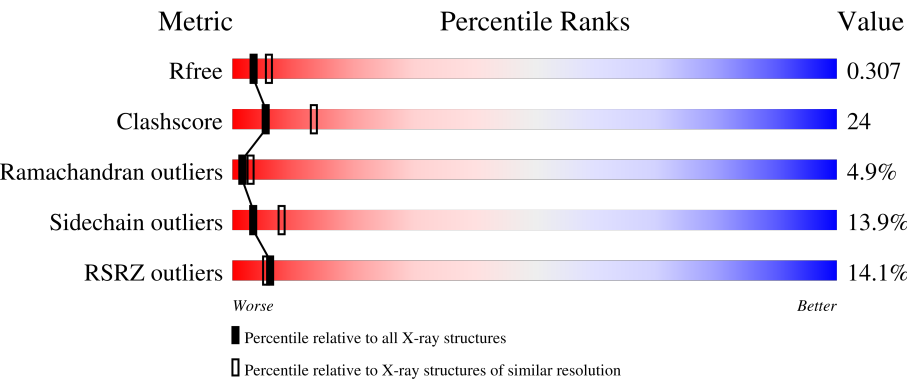
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal 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 i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.72 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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

Mol	Chain	Length	Quality of chain
1	A	244	7% 50% 30% 9% • 8%
1	B	244	21% 47% 32% 9% 5% 7%
2	C	18	6% 33% 61% 6%
2	E	18	44% 28% 28%
3	D	14	7% 71% 29%

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Mol	Chain	Length	Quality of chain
3	F	14	7%21%43%36%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4588 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 Putative cAMP-binding protein-catabolite gene activator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	225	Total	C	N	O	S	0	0	0
			1731	1093	308	322	8			
1	B	226	Total	C	N	O	S	0	0	0
			1727	1091	306	322	8			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	235	LEU	-	expression tag	UNP Q92SD2
A	236	GLU	-	expression tag	UNP Q92SD2
A	237	HIS	-	expression tag	UNP Q92SD2
A	238	HIS	-	expression tag	UNP Q92SD2
A	239	HIS	-	expression tag	UNP Q92SD2
A	240	HIS	-	expression tag	UNP Q92SD2
A	241	HIS	-	expression tag	UNP Q92SD2
A	242	HIS	-	expression tag	UNP Q92SD2
A	243	HIS	-	expression tag	UNP Q92SD2
A	244	HIS	-	expression tag	UNP Q92SD2
B	235	LEU	-	expression tag	UNP Q92SD2
B	236	GLU	-	expression tag	UNP Q92SD2
B	237	HIS	-	expression tag	UNP Q92SD2
B	238	HIS	-	expression tag	UNP Q92SD2
B	239	HIS	-	expression tag	UNP Q92SD2
B	240	HIS	-	expression tag	UNP Q92SD2
B	241	HIS	-	expression tag	UNP Q92SD2
B	242	HIS	-	expression tag	UNP Q92SD2
B	243	HIS	-	expression tag	UNP Q92SD2
B	244	HIS	-	expression tag	UNP Q92SD2

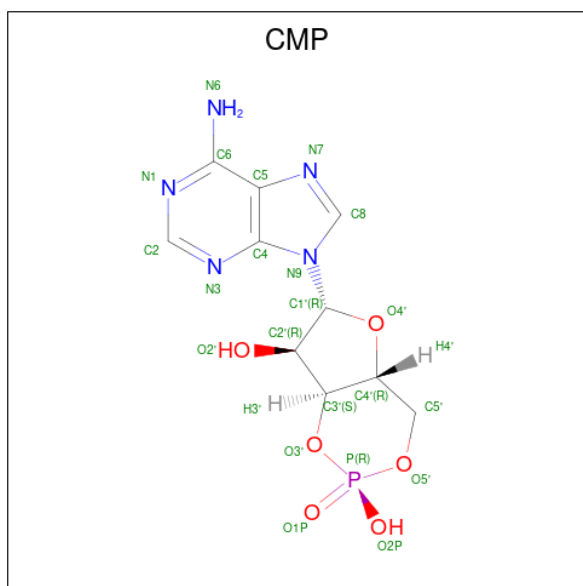
- Molecule 2 is a DNA chain called DNA (5'-D(*CP*TP*AP*GP*GP*TP*AP*AP*CP*AP*TP*TP*AP*CP*TP*CP*GP)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	17	Total	C	N	O	P	0	0	0
			344	166	62	100	16			
2	E	13	Total	C	N	O	P	0	0	0
			264	128	49	75	12			

- Molecule 3 is a DNA chain called DNA (5'-D(*GP*CP*GP*AP*GP*TP*AP*AP*TP*GP*TP*TP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	14	Total	C	N	O	P	0	0	0
			287	138	54	82	13			
3	F	9	Total	C	N	O	P	0	0	0
			184	89	31	55	9			

- Molecule 4 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (CCD ID: CMP) (formula: $C_{10}H_{12}N_5O_6P$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
4	B	1	Total	C	N	O	P	0	0
			22	10	5	6	1		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Mg 1	0	0
5	B	1	Total 1	Mg 1	0	0

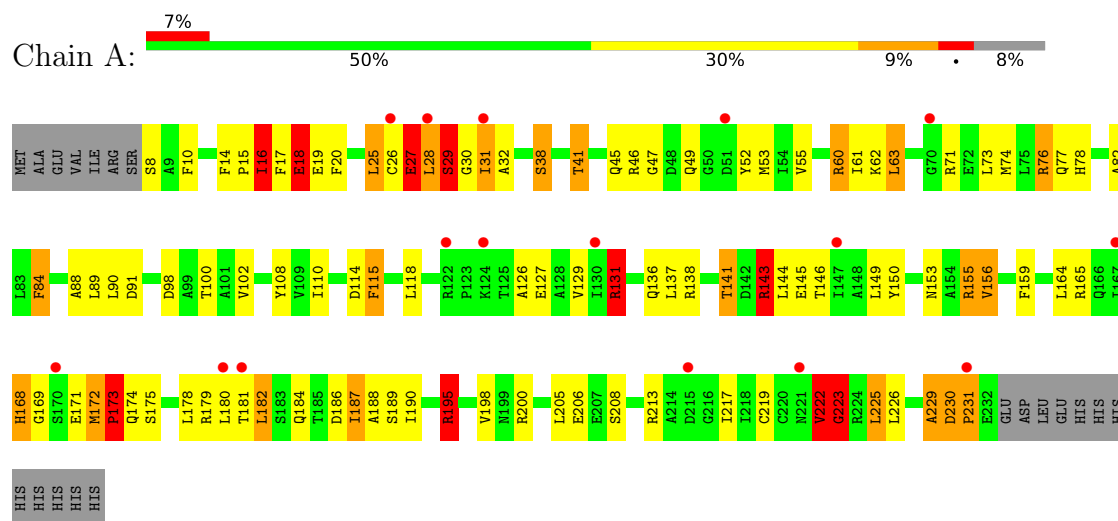
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	3	Total 3	O 3	0	0
6	B	1	Total 1	O 1	0	0
6	F	1	Total 1	O 1	0	0

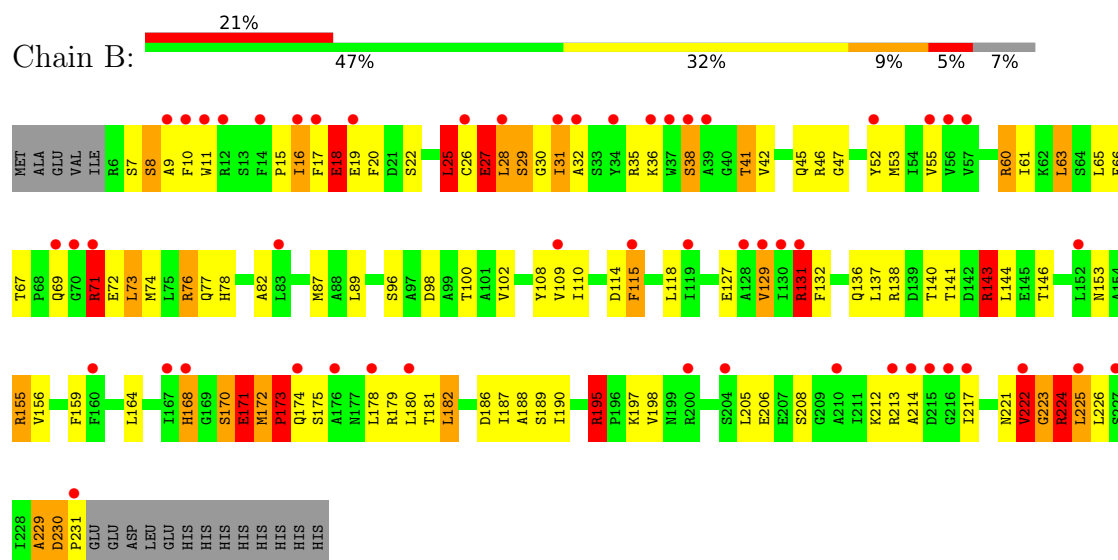
3 Residue-property plots [i](#)

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.

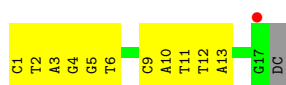
- Molecule 1: Putative cAMP-binding protein-catabolite gene activator



- Molecule 1: Putative cAMP-binding protein-catabolite gene activator



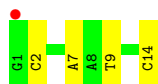
- Molecule 2: DNA (5'-D(*CP*TP*AP*GP*GP*TP*AP*AP*CP*AP*TP*TP*AP*CP*TP*CP*GP)-3')



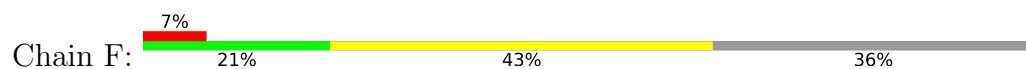
- Molecule 2: DNA (5'-D(*CP*TP*AP*GP*GP*TP*AP*AP*CP*AP*TP*TP*AP*CP*TP*CP*GP)-3')



- Molecule 3: DNA (5'-D(*GP*CP*GP*AP*GP*TP*AP*AP*TP*GP*TP*TP*AP*C)-3')



- Molecule 3: DNA (5'-D(*GP*CP*GP*AP*GP*TP*AP*AP*TP*GP*TP*TP*AP*C)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	50.10Å 69.93Å 200.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.03 – 2.72 66.03 – 2.72	Depositor EDS
% Data completeness (in resolution range)	74.9 (66.03-2.72) 74.9 (66.03-2.72)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.19.1_4122	Depositor
R, R_{free}	0.271 , 0.304 0.270 , 0.307	Depositor DCC
R_{free} test set	747 reflections (3.80%)	wwPDB-VP
Wilson B-factor (Å ²)	64.7	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 51.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4588	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.66% 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: MG, CMP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	1.00	9/1758 (0.5%)	1.42	22/2375 (0.9%)
1	B	0.99	7/1754 (0.4%)	1.44	22/2371 (0.9%)
2	C	1.00	0/385	0.87	0/592
2	E	0.94	0/296	0.94	0/455
3	D	0.89	0/322	0.79	0/496
3	F	0.87	0/205	0.80	0/314
All	All	0.98	16/4720 (0.3%)	1.29	44/6603 (0.7%)

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	13
1	B	0	15
All	All	0	28

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	28	LEU	CA-CB	8.56	1.64	1.53
1	A	28	LEU	CB-CG	8.00	1.69	1.53
1	A	229	ALA	CA-C	7.14	1.62	1.52
1	B	229	ALA	CA-C	7.10	1.62	1.52
1	B	195	ARG	C-O	-7.03	1.18	1.24
1	A	195	ARG	C-O	-7.00	1.18	1.24
1	B	27	GLU	CB-CG	6.95	1.73	1.52
1	A	29	SER	CA-C	6.48	1.61	1.52
1	B	224	ARG	CB-CG	5.88	1.70	1.52
1	B	27	GLU	CD-OE1	5.70	1.36	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	195	ARG	CA-C	5.61	1.58	1.52
1	B	195	ARG	CA-C	5.54	1.58	1.52
1	A	28	LEU	C-O	5.46	1.32	1.24
1	A	173	PRO	C-N	-5.36	1.26	1.33
1	B	173	PRO	C-N	-5.36	1.26	1.33
1	A	28	LEU	CG-CD1	5.11	1.69	1.52

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	28	LEU	N-CA-CB	10.99	126.42	110.16
1	B	28	LEU	CB-CA-C	-10.23	93.46	110.85
1	A	29	SER	N-CA-CB	-8.64	99.57	110.98
1	B	28	LEU	N-CA-C	8.09	120.18	111.36
1	A	29	SER	CB-CA-C	7.72	121.94	109.27
1	A	28	LEU	CA-C-N	7.49	133.44	121.98
1	A	28	LEU	C-N-CA	7.49	133.44	121.98
1	A	29	SER	N-CA-C	-7.39	103.67	114.39
1	A	29	SER	CA-CB-OG	7.36	125.83	111.10
1	B	18	GLU	CA-C-N	-7.32	112.08	124.31
1	B	18	GLU	C-N-CA	-7.32	112.08	124.31
1	A	18	GLU	CA-C-N	-7.32	112.09	124.31
1	A	18	GLU	C-N-CA	-7.32	112.09	124.31
1	A	28	LEU	CB-CG-CD1	7.21	132.33	110.70
1	B	28	LEU	CA-CB-CG	6.95	140.63	116.30
1	A	28	LEU	CB-CA-C	6.86	121.66	111.17
1	A	229	ALA	CA-C-O	-6.50	111.22	120.51
1	B	229	ALA	CA-C-O	-6.44	111.30	120.51
1	A	25	LEU	CB-CG-CD2	-6.02	92.65	110.70
1	B	25	LEU	CB-CG-CD2	-6.01	92.66	110.70
1	A	195	ARG	O-C-N	6.01	125.12	120.38
1	B	195	ARG	O-C-N	5.91	125.05	120.38
1	B	224	ARG	N-CA-CB	-5.84	100.62	110.49
1	A	28	LEU	N-CA-C	-5.78	106.84	112.97
1	B	115	PHE	CB-CA-C	-5.77	101.75	110.92
1	A	115	PHE	CB-CA-C	-5.76	101.76	110.92
1	A	179	ARG	CA-C-N	-5.64	111.14	121.52
1	A	179	ARG	C-N-CA	-5.64	111.14	121.52
1	B	179	ARG	CA-C-N	-5.63	111.16	121.52
1	B	179	ARG	C-N-CA	-5.63	111.16	121.52
1	A	223	GLY	CA-C-N	-5.62	112.68	120.38
1	A	223	GLY	C-N-CA	-5.62	112.68	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	131	ARG	CA-CB-CG	-5.51	103.08	114.10
1	B	171	GLU	N-CA-C	5.49	122.49	110.80
1	B	131	ARG	CA-CB-CG	-5.49	103.13	114.10
1	B	170	SER	CA-C-N	5.43	131.90	121.54
1	B	170	SER	C-N-CA	5.43	131.90	121.54
1	B	195	ARG	CG-CD-NE	5.30	123.67	112.00
1	A	195	ARG	CG-CD-NE	5.29	123.64	112.00
1	B	28	LEU	CA-C-N	-5.27	111.47	121.54
1	B	28	LEU	C-N-CA	-5.27	111.47	121.54
1	B	28	LEU	CB-CG-CD1	-5.11	95.38	110.70
1	B	16	ILE	CG1-CB-CG2	-5.04	95.58	110.70
1	A	16	ILE	CG1-CB-CG2	-5.03	95.61	110.70

There are no chirality outliers.

All (28) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	131	ARG	Sidechain
1	A	143	ARG	Sidechain
1	A	155	ARG	Sidechain
1	A	195	ARG	Sidechain
1	A	223	GLY	Peptide
1	A	27	GLU	Mainchain,Peptide
1	A	29	SER	Peptide
1	A	46	ARG	Sidechain
1	A	47	GLY	Peptide
1	A	60	ARG	Sidechain
1	A	71	ARG	Sidechain
1	A	76	ARG	Sidechain
1	B	131	ARG	Sidechain
1	B	143	ARG	Sidechain
1	B	155	ARG	Sidechain
1	B	170	SER	Peptide
1	B	171	GLU	Peptide
1	B	195	ARG	Sidechain
1	B	223	GLY	Peptide
1	B	27	GLU	Mainchain,Sidechain
1	B	46	ARG	Sidechain
1	B	47	GLY	Peptide
1	B	60	ARG	Sidechain
1	B	71	ARG	Sidechain
1	B	76	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	B	8	SER	Peptide

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	1731	0	1754	96	0
1	B	1727	0	1744	92	0
2	C	344	0	194	12	0
2	E	264	0	149	6	0
3	D	287	0	160	7	0
3	F	184	0	104	11	0
4	A	22	0	11	6	0
4	B	22	0	11	5	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	3	0	0	2	0
6	B	1	0	0	0	0
6	F	1	0	0	1	0
All	All	4588	0	4127	206	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 24.

All (206) 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)
4:A:301:CMP:C2	4:A:301:CMP:H2	0.97	1.49
4:B:301:CMP:H2	4:B:301:CMP:C2	0.97	1.48
1:B:11:TRP:HE1	1:B:29:SER:HB3	1.27	0.95
1:A:168:HIS:HB3	1:A:171:GLU:OE2	1.69	0.92
1:B:20:PHE:HB2	1:B:25:LEU:HD13	1.50	0.91
1:B:172:MET:HE2	1:B:222:VAL:HG21	1.55	0.88
1:A:172:MET:HE2	1:A:222:VAL:HG21	1.55	0.85
1:A:76:ARG:HE	4:A:301:CMP:H2	1.42	0.84
1:B:28:LEU:C	1:B:30:GLY:H	1.86	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:LEU:HD23	1:A:186:ASP:HB3	1.62	0.81
1:B:221:ASN:OD1	1:B:224:ARG:HG3	1.82	0.80
1:B:8:SER:C	1:B:10:PHE:H	1.90	0.79
1:B:223:GLY:H	1:B:225:LEU:H	1.31	0.79
1:A:180:LEU:O	1:A:181:THR:HG22	1.84	0.78
1:A:230:ASP:HA	1:B:69:GLN:NE2	1.99	0.78
1:B:11:TRP:NE1	1:B:29:SER:HB3	1.97	0.77
1:B:180:LEU:O	1:B:181:THR:HG22	1.84	0.77
1:A:126:ALA:O	1:A:129:VAL:HG12	1.83	0.77
1:B:146:THR:HG22	1:B:155:ARG:HG2	1.67	0.77
1:A:165:ARG:HG3	1:A:172:MET:SD	2.25	0.77
1:A:146:THR:HG22	1:A:155:ARG:HG2	1.67	0.76
1:A:171:GLU:HG3	1:A:172:MET:H	1.50	0.76
1:A:230:ASP:HA	1:B:69:GLN:HE22	1.52	0.75
1:B:53:MET:HE1	1:B:115:PHE:CE1	2.22	0.74
1:A:141:THR:HG21	4:B:301:CMP:HN61	1.51	0.74
1:A:29:SER:C	1:A:31:ILE:H	1.96	0.74
1:A:53:MET:HE1	1:A:115:PHE:CE1	2.22	0.73
1:B:182:LEU:HD23	1:B:186:ASP:HB3	1.72	0.72
1:B:197:LYS:HE2	2:E:5:DG:N7	2.05	0.72
1:A:141:THR:CG2	4:B:301:CMP:HN61	2.03	0.72
1:B:164:LEU:CD1	1:B:178:LEU:HD22	2.20	0.72
1:A:164:LEU:CD1	1:A:178:LEU:HD22	2.20	0.71
1:B:212:LYS:HD3	1:B:213:ARG:H	1.55	0.71
1:B:38:SER:O	1:B:41:THR:HG23	1.91	0.71
1:A:38:SER:O	1:A:41:THR:HG23	1.91	0.71
1:B:66:PHE:CE1	1:B:72:GLU:HG3	2.25	0.71
1:B:20:PHE:HB2	1:B:25:LEU:CD1	2.21	0.71
1:B:89:LEU:HD23	1:B:115:PHE:HD1	1.58	0.69
1:B:28:LEU:C	1:B:30:GLY:N	2.44	0.69
1:A:89:LEU:HD23	1:A:115:PHE:HD1	1.58	0.68
1:B:221:ASN:O	1:B:224:ARG:HB2	1.94	0.68
1:A:18:GLU:O	1:A:19:GLU:HB3	1.94	0.67
3:F:9:DT:C2	3:F:10:DG:C8	2.83	0.67
1:A:27:GLU:C	1:A:29:SER:H	2.01	0.67
1:A:20:PHE:HB2	1:A:25:LEU:HD13	1.77	0.66
2:C:12:DT:H2''	2:C:13:DA:H5'	1.77	0.66
1:B:18:GLU:O	1:B:19:GLU:HB3	1.94	0.66
1:A:19:GLU:OE2	1:A:131:ARG:CZ	2.43	0.66
1:B:19:GLU:OE2	1:B:131:ARG:CZ	2.43	0.66
1:A:180:LEU:HD21	1:A:187:ILE:HD11	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:GLY:H	1:A:225:LEU:H	1.45	0.65
1:B:8:SER:H	1:B:10:PHE:HB3	1.61	0.64
3:F:9:DT:C2	3:F:10:DG:N7	2.67	0.63
1:A:127:GLU:OE1	1:A:131:ARG:NH1	2.32	0.63
1:B:15:PRO:O	1:B:16:ILE:HB	1.98	0.63
1:B:127:GLU:OE1	1:B:131:ARG:NH1	2.32	0.62
1:B:223:GLY:HA2	1:B:226:LEU:HD12	1.81	0.62
1:A:76:ARG:NE	4:A:301:CMP:H2	2.15	0.61
1:A:223:GLY:HA2	1:A:226:LEU:HD12	1.81	0.61
1:A:195:ARG:HE	3:D:9:DT:C7	2.14	0.60
1:A:14:PHE:CD1	1:A:55:VAL:HG22	2.36	0.60
1:B:71:ARG:HG3	1:B:71:ARG:HH11	1.67	0.60
1:B:188:ALA:HA	1:B:198:VAL:HG21	1.84	0.60
1:A:188:ALA:HA	1:A:198:VAL:HG21	1.84	0.58
3:F:8:DA:H2"	6:F:101:HOH:O	2.03	0.58
1:A:16:ILE:HG22	1:A:17:PHE:N	2.17	0.58
1:A:27:GLU:CB	1:A:28:LEU:HD12	2.34	0.58
1:A:60:ARG:HB2	1:A:102:VAL:HG23	1.85	0.58
1:A:159:PHE:CE2	1:A:187:ILE:HD13	2.38	0.58
1:B:60:ARG:HB2	1:B:102:VAL:HG23	1.85	0.58
1:A:53:MET:HE1	1:A:115:PHE:CD1	2.39	0.57
1:B:212:LYS:HD3	1:B:213:ARG:N	2.19	0.57
1:B:8:SER:C	1:B:10:PHE:N	2.58	0.57
1:B:53:MET:HE1	1:B:115:PHE:CD1	2.39	0.57
1:A:63:LEU:HD12	1:A:78:HIS:HD2	1.70	0.57
1:A:149:LEU:HD11	1:B:73:LEU:HB2	1.86	0.57
1:A:84:PHE:CD1	1:A:84:PHE:C	2.84	0.56
1:A:164:LEU:HD11	1:A:178:LEU:HD22	1.87	0.56
1:B:63:LEU:HD12	1:B:78:HIS:HD2	1.70	0.56
2:C:1:DC:H2"	2:C:2:DT:OP2	2.05	0.56
4:A:301:CMP:HN61	1:B:141:THR:CG2	2.19	0.56
1:A:15:PRO:O	1:A:16:ILE:HB	2.06	0.55
1:A:182:LEU:HD23	1:A:186:ASP:CB	2.35	0.55
1:B:164:LEU:HD11	1:B:178:LEU:HD22	1.87	0.55
2:C:3:DA:H2"	2:C:4:DG:H5"	1.88	0.55
1:A:156:VAL:HG12	1:A:205:LEU:HD11	1.88	0.54
1:A:231:PRO:HB3	1:B:69:GLN:HG3	1.89	0.54
1:A:164:LEU:HD11	1:A:178:LEU:HB2	1.90	0.54
1:B:17:PHE:C	1:B:18:GLU:O	2.50	0.54
1:B:7:SER:O	1:B:7:SER:OG	2.18	0.53
1:A:195:ARG:HE	3:D:9:DT:H73	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:ASP:C	1:B:69:GLN:HE21	2.15	0.53
1:A:17:PHE:C	1:A:18:GLU:O	2.50	0.53
1:A:29:SER:C	1:A:31:ILE:N	2.67	0.53
1:B:164:LEU:HD11	1:B:178:LEU:HB2	1.90	0.53
2:E:3:DA:H8	2:E:3:DA:H5''	1.74	0.52
1:A:19:GLU:OE2	1:A:131:ARG:NH2	2.42	0.52
1:A:60:ARG:HB2	1:A:102:VAL:CG2	2.40	0.52
3:D:2:DC:H4'	3:D:2:DC:OP1	2.08	0.52
1:B:19:GLU:OE2	1:B:131:ARG:NH2	2.42	0.52
1:B:156:VAL:HG12	1:B:205:LEU:HD11	1.92	0.52
1:A:84:PHE:C	1:A:84:PHE:HD1	2.18	0.52
1:B:221:ASN:CG	1:B:224:ARG:HG3	2.35	0.52
1:B:78:HIS:ND1	1:B:82:ALA:HB1	2.25	0.52
1:A:168:HIS:CD2	1:A:168:HIS:H	2.28	0.52
1:B:60:ARG:HB2	1:B:102:VAL:CG2	2.40	0.51
1:B:11:TRP:HE1	1:B:29:SER:CB	2.11	0.51
1:A:60:ARG:HB3	1:A:77:GLN:HE21	1.76	0.51
1:B:96:SER:OG	4:B:301:CMP:H8	2.11	0.50
1:B:168:HIS:H	1:B:168:HIS:CD2	2.28	0.50
1:A:78:HIS:ND1	1:A:82:ALA:HB1	2.25	0.50
2:C:11:DT:H2'	2:C:12:DT:C7	2.42	0.50
1:A:169:GLY:O	1:A:171:GLU:HG2	2.11	0.50
3:F:11:DT:C2'	3:F:12:DT:H71	2.42	0.50
2:C:4:DG:H2''	2:C:5:DG:H8	1.77	0.50
1:B:71:ARG:HG3	1:B:71:ARG:NH1	2.27	0.50
1:A:168:HIS:CB	1:A:171:GLU:OE2	2.52	0.49
1:B:19:GLU:OE2	1:B:131:ARG:NH1	2.45	0.49
2:C:1:DC:H1'	2:C:2:DT:H5'	1.94	0.49
1:A:19:GLU:OE2	1:A:131:ARG:NH1	2.45	0.49
1:B:60:ARG:HB3	1:B:77:GLN:HE21	1.76	0.49
3:F:9:DT:O2	3:F:10:DG:C8	2.66	0.49
3:F:11:DT:H2'	3:F:12:DT:H71	1.95	0.49
1:A:27:GLU:C	1:A:29:SER:N	2.70	0.49
3:D:14:DC:C4	2:E:1:DC:C4	3.01	0.48
1:B:76:ARG:HE	4:B:301:CMP:H2	1.78	0.48
1:A:91:ASP:OD1	1:B:138:ARG:NE	2.44	0.47
3:F:10:DG:H2''	3:F:11:DT:O5'	2.15	0.47
4:A:301:CMP:HN61	1:B:141:THR:HG21	1.79	0.47
1:B:195:ARG:HB3	1:B:195:ARG:NH1	2.29	0.47
1:B:214:ALA:HB3	1:B:217:ILE:HG22	1.95	0.47
2:C:13:DA:C2	3:D:7:DA:C2	3.02	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:ILE:HG23	1:B:110:ILE:HG23	1.97	0.47
1:A:63:LEU:HD12	1:A:78:HIS:CD2	2.49	0.47
2:C:4:DG:H2''	2:C:5:DG:C8	2.50	0.47
1:A:61:ILE:HG23	1:A:100:THR:O	2.14	0.46
1:B:31:ILE:HG12	1:B:114:ASP:CG	2.41	0.46
1:A:31:ILE:HG12	1:A:114:ASP:CG	2.41	0.46
1:A:195:ARG:NH1	1:A:195:ARG:HB3	2.29	0.46
2:C:5:DG:H2'	2:C:6:DT:H71	1.98	0.46
1:B:61:ILE:HG23	1:B:100:THR:O	2.14	0.46
1:A:138:ARG:NH1	1:B:96:SER:HB3	2.31	0.46
1:A:31:ILE:HG12	1:A:114:ASP:CB	2.46	0.46
1:A:31:ILE:HG23	1:A:110:ILE:HG23	1.97	0.46
1:B:31:ILE:HG12	1:B:114:ASP:CB	2.46	0.45
1:A:118:LEU:HA	1:A:118:LEU:HD12	1.69	0.45
1:A:173:PRO:HB2	1:A:174:GLN:H	1.38	0.45
2:C:9:DC:H2'	2:C:10:DA:C8	2.52	0.45
1:B:137:LEU:HD12	1:B:137:LEU:HA	1.73	0.45
2:E:3:DA:C6	2:E:4:DG:C6	3.04	0.45
1:A:88:ALA:HB1	6:A:402:HOH:O	2.17	0.45
1:A:53:MET:HE2	1:A:89:LEU:HD22	1.98	0.45
1:A:8:SER:C	6:A:401:HOH:O	2.60	0.45
1:A:230:ASP:HB3	1:A:231:PRO:CD	2.47	0.45
1:B:32:ALA:HB1	1:B:108:TYR:HB3	1.98	0.45
1:B:230:ASP:HB3	1:B:231:PRO:CD	2.47	0.45
1:A:32:ALA:HB1	1:A:108:TYR:HB3	1.98	0.45
1:B:63:LEU:HD12	1:B:78:HIS:CD2	2.49	0.45
1:B:143:ARG:NH1	1:B:190:ILE:HG23	2.31	0.44
1:A:153:ASN:HD22	1:A:153:ASN:N	2.15	0.44
1:B:53:MET:HE2	1:B:89:LEU:HD22	1.98	0.44
1:A:90:LEU:H	1:A:90:LEU:HG	1.67	0.44
1:B:159:PHE:CE2	1:B:187:ILE:CD1	3.00	0.44
1:A:141:THR:HG23	1:B:140:THR:HG21	1.99	0.44
1:A:143:ARG:NH1	1:A:190:ILE:HG23	2.31	0.44
1:A:145:GLU:HB2	1:B:65:LEU:HD13	2.00	0.44
1:A:150:TYR:CE1	1:B:67:THR:HG21	2.52	0.44
1:B:118:LEU:HD12	1:B:118:LEU:HA	1.69	0.44
1:A:31:ILE:HG12	1:A:114:ASP:HB3	2.00	0.43
1:A:200:ARG:HA	1:A:200:ARG:HD3	1.56	0.43
1:A:137:LEU:HA	1:A:137:LEU:HD12	1.73	0.43
1:B:153:ASN:N	1:B:153:ASN:HD22	2.15	0.43
2:C:11:DT:H2'	2:C:12:DT:H72	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:GLY:N	1:B:225:LEU:H	2.06	0.43
1:B:180:LEU:HD21	1:B:187:ILE:HD11	1.99	0.43
1:B:173:PRO:HB2	1:B:174:GLN:H	1.38	0.43
1:B:206:GLU:OE2	1:B:213:ARG:NH1	2.51	0.43
1:B:31:ILE:HG12	1:B:114:ASP:HB3	2.00	0.43
1:A:184:GLN:HG3	3:D:9:DT:OP2	2.19	0.43
1:A:206:GLU:OE2	1:A:213:ARG:NH1	2.51	0.43
1:B:25:LEU:O	1:B:28:LEU:HB2	2.18	0.42
1:A:223:GLY:H	1:A:225:LEU:N	2.15	0.42
1:B:28:LEU:O	1:B:30:GLY:N	2.53	0.42
3:F:9:DT:N3	3:F:10:DG:N7	2.68	0.42
1:A:45:GLN:HA	1:A:98:ASP:OD1	2.20	0.41
1:B:63:LEU:CD1	1:B:78:HIS:CD2	3.04	0.41
1:A:63:LEU:HD22	4:A:301:CMP:O4'	2.20	0.41
2:E:7:DA:C2	3:F:13:DA:C2	3.08	0.41
1:B:74:MET:HE3	1:B:74:MET:HB2	1.82	0.41
2:C:9:DC:H2'	2:C:10:DA:H8	1.85	0.41
1:A:74:MET:HE3	1:A:74:MET:HB2	1.82	0.41
1:A:8:SER:C	1:A:10:PHE:H	2.29	0.41
1:A:27:GLU:O	1:A:29:SER:N	2.45	0.41
1:A:168:HIS:HB3	1:A:171:GLU:CD	2.41	0.41
1:B:195:ARG:HE	3:F:9:DT:H72	1.85	0.41
1:A:136:GLN:OE1	1:A:136:GLN:HA	2.21	0.41
1:B:45:GLN:HA	1:B:98:ASP:OD1	2.20	0.41
1:B:136:GLN:HA	1:B:136:GLN:OE1	2.21	0.41
2:E:3:DA:H5''	2:E:3:DA:C8	2.54	0.41
3:F:8:DA:H2'	3:F:9:DT:H71	2.02	0.41
1:A:63:LEU:CD1	1:A:78:HIS:CD2	3.04	0.40
1:A:184:GLN:HG3	3:D:9:DT:P	2.61	0.40
1:A:55:VAL:HG12	1:A:108:TYR:HB2	2.03	0.40
1:B:182:LEU:HD23	1:B:186:ASP:CB	2.47	0.40
1:A:138:ARG:HG3	1:B:87:MET:CE	2.51	0.40
1:B:129:VAL:O	1:B:132:PHE:HB3	2.22	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	223/244 (91%)	198 (89%)	15 (7%)	10 (4%)	2	3
1	B	224/244 (92%)	197 (88%)	15 (7%)	12 (5%)	1	2
All	All	447/488 (92%)	395 (88%)	30 (7%)	22 (5%)	1	3

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	27	GLU
1	A	30	GLY
1	A	222	VAL
1	A	230	ASP
1	A	231	PRO
1	B	9	ALA
1	B	27	GLU
1	B	171	GLU
1	B	222	VAL
1	B	224	ARG
1	B	230	ASP
1	A	18	GLU
1	A	168	HIS
1	A	173	PRO
1	A	229	ALA
1	B	18	GLU
1	B	29	SER
1	B	168	HIS
1	B	173	PRO
1	B	229	ALA
1	A	172	MET
1	B	172	MET

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	181/201 (90%)	156 (86%)	25 (14%)	3	8
1	B	180/201 (90%)	155 (86%)	25 (14%)	3	8
All	All	361/402 (90%)	311 (86%)	50 (14%)	3	8

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ILE
1	A	26	CYS
1	A	31	ILE
1	A	38	SER
1	A	41	THR
1	A	49	GLN
1	A	52	TYR
1	A	62	LYS
1	A	63	LEU
1	A	73	LEU
1	A	84	PHE
1	A	141	THR
1	A	143	ARG
1	A	144	LEU
1	A	156	VAL
1	A	175	SER
1	A	182	LEU
1	A	187	ILE
1	A	189	SER
1	A	195	ARG
1	A	208	SER
1	A	217	ILE
1	A	219	CYS
1	A	222	VAL
1	A	225	LEU
1	B	22	SER
1	B	25	LEU

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Mol	Chain	Res	Type
1	B	26	CYS
1	B	31	ILE
1	B	35	ARG
1	B	36	LYS
1	B	38	SER
1	B	41	THR
1	B	42	VAL
1	B	52	TYR
1	B	55	VAL
1	B	63	LEU
1	B	71	ARG
1	B	73	LEU
1	B	109	VAL
1	B	129	VAL
1	B	143	ARG
1	B	144	LEU
1	B	175	SER
1	B	182	LEU
1	B	189	SER
1	B	195	ARG
1	B	208	SER
1	B	222	VAL
1	B	225	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	HIS
1	A	153	ASN
1	B	69	GLN
1	B	78	HIS
1	B	153	ASN
1	B	168	HIS
1	B	199	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 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$
4	CMP	B	301	-	25,25,25	1.73	7 (28%)	37,39,39	2.12	12 (32%)
4	CMP	A	301	-	25,25,25	1.72	7 (28%)	37,39,39	2.05	12 (32%)

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	CMP	B	301	-	-	0/4/31/31	0/4/4/4
4	CMP	A	301	-	-	0/4/31/31	0/4/4/4

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	301	CMP	C5-C4	4.24	1.46	1.39
4	B	301	CMP	C5-C4	4.24	1.46	1.39
4	A	301	CMP	C4-N9	-3.04	1.31	1.37
4	B	301	CMP	C4-N9	-3.04	1.31	1.37
4	B	301	CMP	P-O3'	2.85	1.62	1.57
4	A	301	CMP	P-O3'	2.84	1.62	1.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	301	CMP	C5-C6	2.58	1.48	1.41
4	B	301	CMP	C5-C6	2.58	1.48	1.41
4	B	301	CMP	O5'-C5'	-2.50	1.42	1.46
4	A	301	CMP	O5'-C5'	-2.50	1.42	1.46
4	A	301	CMP	O3'-C3'	-2.42	1.40	1.44
4	B	301	CMP	O3'-C3'	-2.42	1.40	1.44
4	A	301	CMP	C8-N7	2.36	1.36	1.31
4	B	301	CMP	C8-N7	2.36	1.36	1.31

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	301	CMP	O5'-P-O1P	-4.60	99.82	110.44
4	A	301	CMP	C5-C4-N3	-4.60	120.38	126.72
4	B	301	CMP	C5-C4-N3	-4.60	120.39	126.72
4	B	301	CMP	C4-N9-C8	4.07	110.01	105.74
4	A	301	CMP	C4-N9-C8	4.07	110.01	105.74
4	A	301	CMP	O2P-P-O1P	3.90	120.56	108.56
4	B	301	CMP	O2P-P-O1P	3.90	120.55	108.56
4	A	301	CMP	N3-C4-N9	3.78	133.60	127.17
4	B	301	CMP	N3-C4-N9	3.78	133.60	127.17
4	A	301	CMP	C2-N3-C4	3.47	120.31	111.83
4	B	301	CMP	C2-N3-C4	3.47	120.31	111.83
4	A	301	CMP	N3-C2-N1	-3.44	123.37	128.58
4	B	301	CMP	N3-C2-N1	-3.44	123.37	128.58
4	A	301	CMP	C4-C5-N7	-3.31	106.80	110.58
4	B	301	CMP	C4-C5-N7	-3.31	106.80	110.58
4	A	301	CMP	C6-C5-N7	3.17	138.21	132.09
4	B	301	CMP	C6-C5-N7	3.17	138.21	132.09
4	A	301	CMP	O2P-P-O5'	-3.00	99.82	107.16
4	A	301	CMP	N9-C8-N7	-2.50	110.39	113.94
4	B	301	CMP	N9-C8-N7	-2.50	110.39	113.94
4	A	301	CMP	O3'-C3'-C2'	2.42	117.98	115.61
4	B	301	CMP	O3'-C3'-C2'	2.42	117.98	115.61
4	A	301	CMP	C5-N7-C8	2.13	106.79	103.45
4	B	301	CMP	C5-N7-C8	2.12	106.79	103.45

There are no chirality outliers.

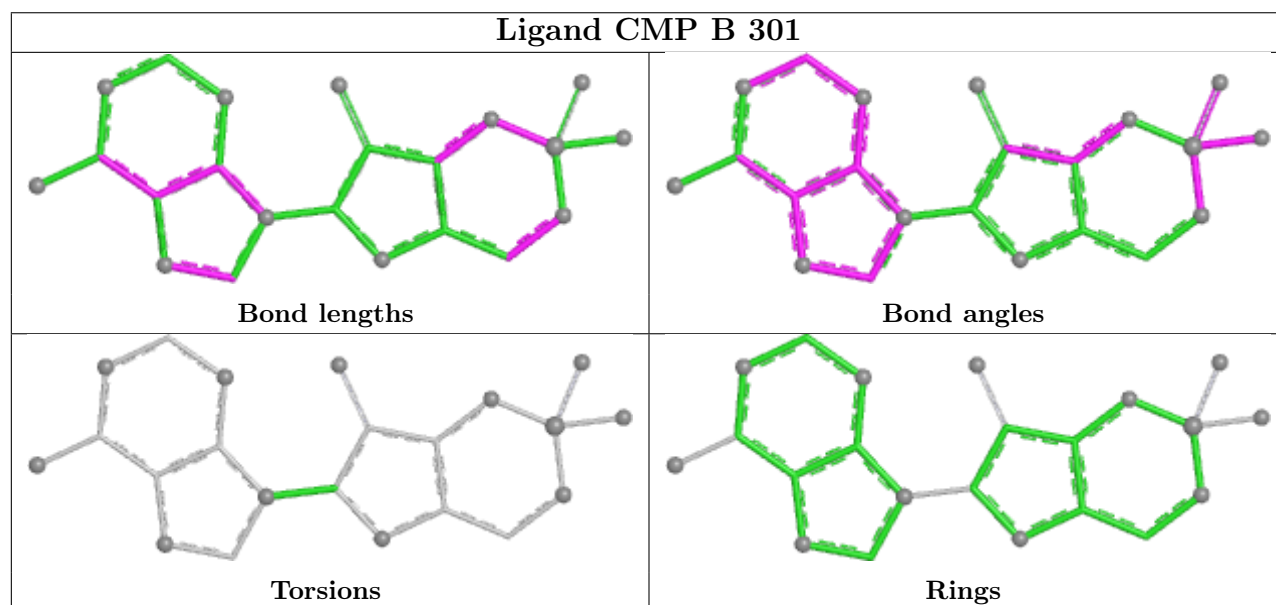
There are no torsion outliers.

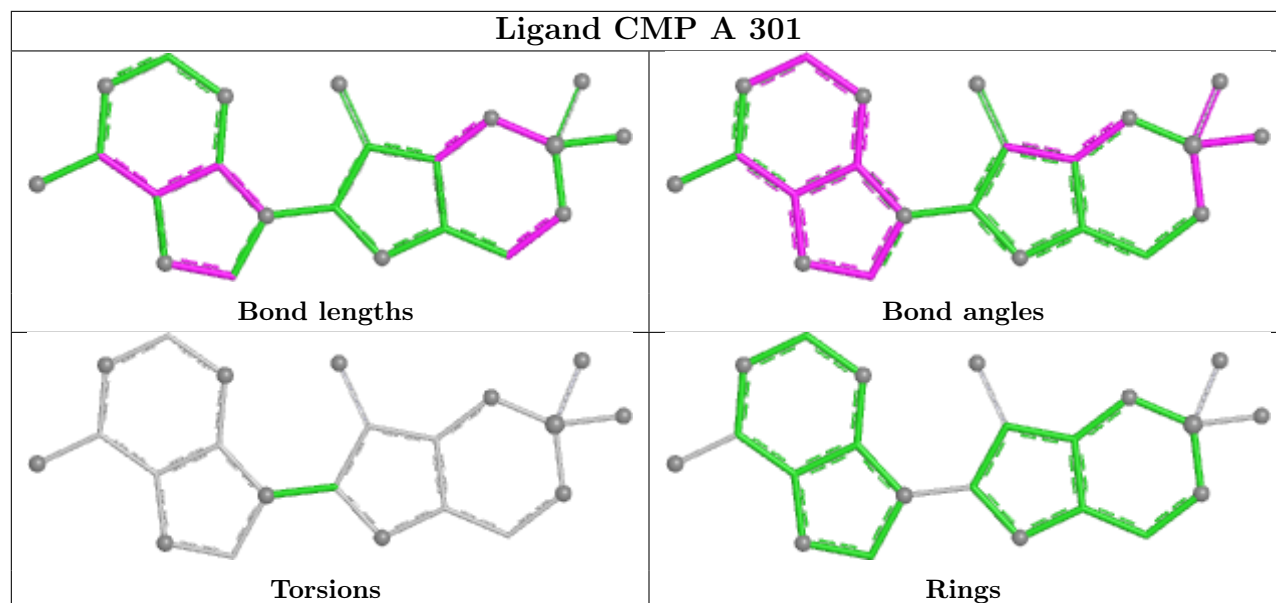
There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	301	CMP	5	0
4	A	301	CMP	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	225/244 (92%)	0.79	16 (7%) 22 19	33, 49, 79, 108	0
1	B	226/244 (92%)	1.35	52 (23%) 2 2	35, 77, 108, 131	0
2	C	17/18 (94%)	0.36	1 (5%) 28 25	47, 72, 114, 118	0
2	E	13/18 (72%)	0.50	0 100 100	53, 74, 95, 109	0
3	D	14/14 (100%)	0.68	1 (7%) 22 19	51, 73, 119, 120	0
3	F	9/14 (64%)	0.93	1 (11%) 10 9	73, 85, 101, 113	0
All	All	504/552 (91%)	1.02	71 (14%) 6 5	33, 63, 108, 131	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	231	PRO	6.6
1	A	28	LEU	4.4
1	B	57	VAL	4.2
1	B	225	LEU	4.2
1	A	180	LEU	4.1
1	B	214	ALA	4.0
1	B	180	LEU	3.8
1	B	167	ILE	3.4
1	B	12	ARG	3.4
1	B	34	TYR	3.3
1	B	215	ASP	3.3
1	B	56	VAL	3.2
1	B	129	VAL	3.2
1	B	14	PHE	3.1
1	B	216	GLY	3.0
1	B	39	ALA	3.0
1	B	200	ARG	3.0
1	B	28	LEU	2.9
1	B	174	GLN	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	71	ARG	2.9
1	B	115	PHE	2.8
1	A	170	SER	2.8
1	B	109	VAL	2.8
1	A	122	ARG	2.8
1	A	26	CYS	2.7
1	B	217	ILE	2.7
1	B	70	GLY	2.7
1	B	55	VAL	2.7
1	A	221	ASN	2.6
1	B	210	ALA	2.6
1	A	167	ILE	2.5
1	B	176	ALA	2.5
1	B	36	LYS	2.5
1	B	69	GLN	2.5
1	B	130	ILE	2.4
1	B	11	TRP	2.4
2	C	17	DG	2.4
1	B	227	SER	2.4
1	B	31	ILE	2.4
1	B	17	PHE	2.4
1	B	168	HIS	2.4
1	B	32	ALA	2.4
1	B	178	LEU	2.3
1	B	38	SER	2.3
1	B	222	VAL	2.3
1	A	31	ILE	2.3
1	A	130	ILE	2.2
1	B	9	ALA	2.2
1	B	26	CYS	2.2
1	B	52	TYR	2.2
3	F	14	DC	2.2
1	B	131	ARG	2.2
3	D	1	DG	2.2
1	B	160	PHE	2.2
1	A	51	ASP	2.2
1	B	204	SER	2.2
1	A	147	ILE	2.2
1	B	213	ARG	2.1
1	A	231	PRO	2.1
1	B	128	ALA	2.1
1	A	181	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	83	LEU	2.1
1	B	37	TRP	2.1
1	B	19	GLU	2.0
1	A	124	LYS	2.0
1	B	152	LEU	2.0
1	B	16	ILE	2.0
1	B	10	PHE	2.0
1	A	215	ASP	2.0
1	A	70	GLY	2.0
1	B	119	ILE	2.0

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](#)

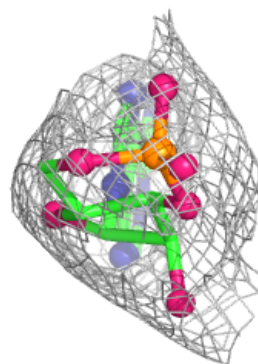
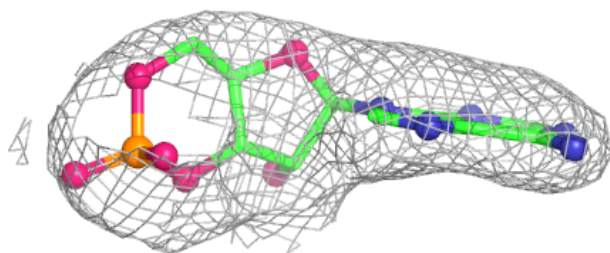
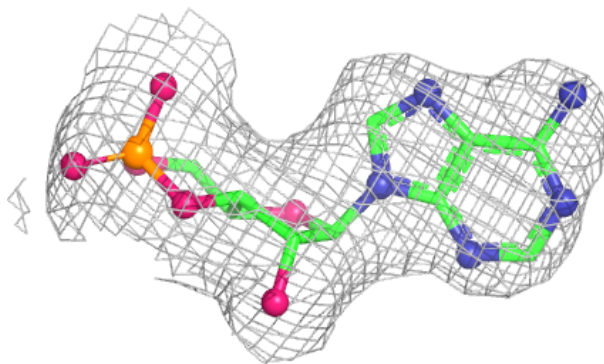
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MG	B	302	1/1	0.93	0.11	40,40,40,40	0
4	CMP	B	301	22/22	0.94	0.09	36,55,64,75	0
4	CMP	A	301	22/22	0.94	0.10	24,39,47,48	0
5	MG	A	302	1/1	0.95	0.15	28,28,28,28	0

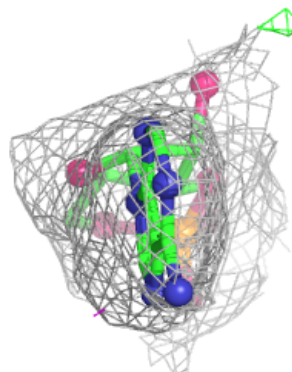
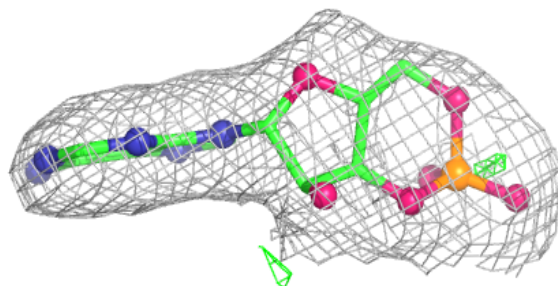
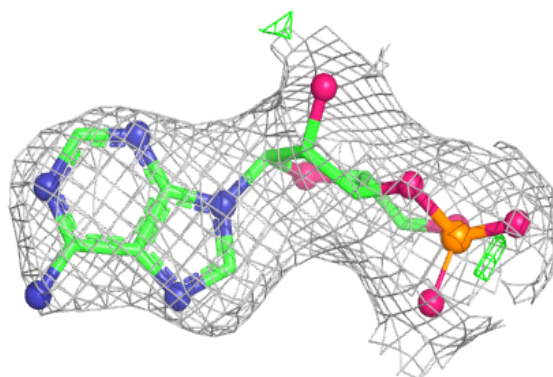
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CMP B 301:

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

**Electron density around CMP A 301:**

$2mF_o-DF_c$ (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.