



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:52 PM UTC

PDB ID : 1CGP / pdb_00001cgp
Title : CATABOLITE GENE ACTIVATOR PROTEIN (CAP)/DNA COMPLEX +
ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE
Authors : Schultz, S.C.; Shields, G.C.; Steitz, T.A.
Deposited on : 1991-08-12
Resolution : 3.00 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

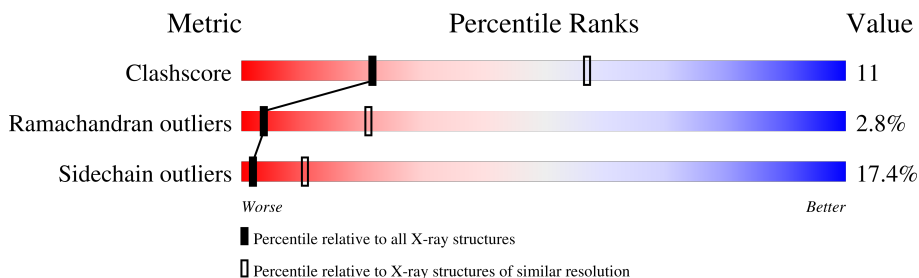
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)

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

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	18	
1	E	18	
2	D	13	
2	F	13	
3	A	205	
3	B	205	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5274 atoms, of which 856 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 DNA chain called DNA (5'-D(*GP*CP*GP*AP*AP*AP*AP*GP*TP*GP*TP*GP*AP*CP*AP*TP*AP*T)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	C	18	Total	C	H	N	O	P	0	0	0
			411	178	39	74	103	17			
1	E	18	Total	C	H	N	O	P	0	0	0
			411	178	39	74	103	17			

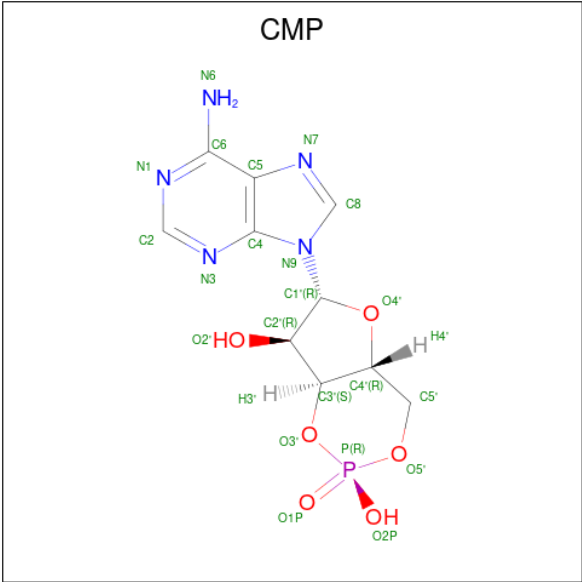
- Molecule 2 is a DNA chain called DNA (5'-D(*GP*TP*CP*AP*CP*AP*CP*TP*TP*TP*TP*CP*G)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	D	13	Total	C	H	N	O	P	0	0	0
			284	126	25	42	79	12			
2	F	13	Total	C	H	N	O	P	0	0	0
			284	126	25	42	79	12			

- Molecule 3 is a protein called PROTEIN (CATABOLITE GENE ACTIVATOR PROTEIN (CAP)).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	A	197	Total	C	H	N	O	S	0	0	0
			1917	986	361	273	288	9			
3	B	197	Total	C	H	N	O	S	0	0	0
			1917	986	361	273	288	9			

- Molecule 4 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (CCD ID: CMP) (formula: C₁₀H₁₂N₅O₆P).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	P	0	0
			25	10	3	5	6	1		
4	B	1	Total	C	H	N	O	P	0	0
			25	10	3	5	6	1		

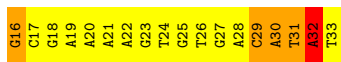
3 Residue-property plots [i](#)

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. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

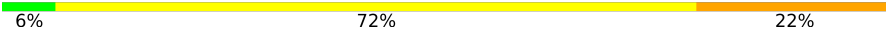
Note EDS was not executed.

- Molecule 1: DNA (5'-D(*GP*CP*GP*AP*AP*AP*AP*GP*TP*GP*TP*GP*AP*CP*AP*TP*AP*T)-3')

Chain C: 



- Molecule 1: DNA (5'-D(*GP*CP*GP*AP*AP*AP*AP*GP*TP*GP*TP*GP*AP*CP*AP*TP*AP*T)-3')

Chain E: 

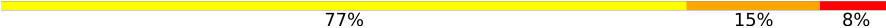


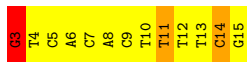
- Molecule 2: DNA (5'-D(*GP*TP*CP*AP*CP*AP*CP*TP*TP*TP*TP*CP*G)-3')

Chain D: 



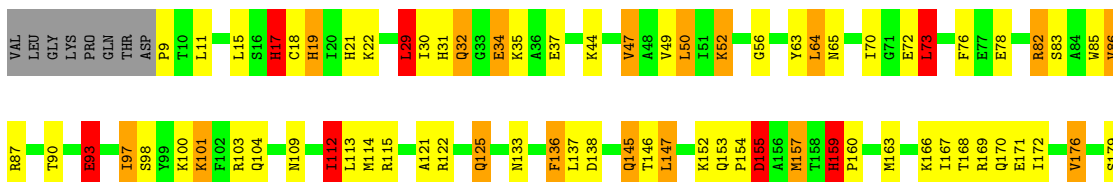
- Molecule 2: DNA (5'-D(*GP*TP*CP*AP*CP*AP*CP*TP*TP*TP*TP*CP*G)-3')

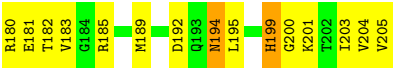
Chain F: 



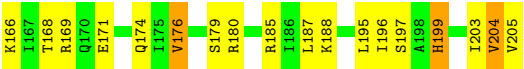
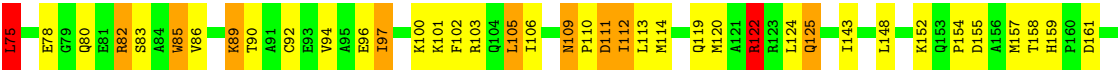
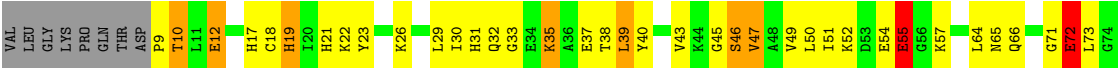
- Molecule 3: PROTEIN (CATALYTIC GENE ACTIVATOR PROTEIN (CAP))

Chain A: 





● Molecule 3: PROTEIN (CATABOLITE GENE ACTIVATOR PROTEIN (CAP))



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	138.00Å 152.60Å 76.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 3.00	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.235 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5274	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	C	0.89	0/419	3.36	46/646 (7.1%)
1	E	0.77	0/419	3.85	47/646 (7.3%)
2	D	0.73	0/288	4.06	29/442 (6.6%)
2	F	0.81	0/288	3.20	25/442 (5.7%)
3	A	1.02	6/1580 (0.4%)	1.87	37/2127 (1.7%)
3	B	1.03	7/1580 (0.4%)	1.87	41/2127 (1.9%)
All	All	0.96	13/4574 (0.3%)	2.60	225/6430 (3.5%)

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	C	0	1
2	D	0	1
2	F	0	2
3	A	0	1
All	All	0	5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	17	HIS	CD2-NE2	-8.32	1.28	1.37
3	B	17	HIS	CD2-NE2	-6.68	1.30	1.37
3	B	31	HIS	CD2-NE2	-6.15	1.31	1.37
3	A	19	HIS	CD2-NE2	-5.97	1.31	1.37
3	A	21	HIS	CD2-NE2	-5.93	1.31	1.37
3	B	159	HIS	CD2-NE2	-5.91	1.31	1.37
3	A	199	HIS	CD2-NE2	-5.91	1.31	1.37
3	B	19	HIS	CD2-NE2	-5.78	1.31	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	159	HIS	CD2-NE2	-5.72	1.31	1.37
3	B	21	HIS	CD2-NE2	-5.66	1.31	1.37
3	B	199	HIS	CD2-NE2	-5.43	1.31	1.37
3	A	31	HIS	CD2-NE2	-5.39	1.31	1.37
3	B	57	LYS	CA-CB	5.25	1.61	1.53

All (225) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	12	DT	O3'-P-O5'	38.68	162.02	104.00
1	E	19	DA	O3'-P-O5'	36.55	158.82	104.00
2	D	9	DC	O3'-P-O5'	35.83	157.74	104.00
1	E	18	DG	O3'-P-O5'	33.78	154.67	104.00
2	F	14	DC	O3'-P-O5'	28.91	147.37	104.00
1	E	25	DG	O3'-P-O5'	26.68	144.02	104.00
1	C	30	DA	O3'-P-O5'	24.54	140.81	104.00
2	D	10	DT	O3'-P-O5'	24.19	140.29	104.00
1	C	31	DT	O3'-P-O5'	22.18	137.28	104.00
1	C	16	DG	O3'-P-O5'	21.76	136.64	104.00
1	E	24	DT	OP2-P-O3'	21.29	171.86	108.00
2	D	4	DT	OP2-P-O3'	18.97	164.91	108.00
1	E	24	DT	O3'-P-O5'	-18.64	76.04	104.00
1	C	27	DG	O3'-P-O5'	-18.63	76.05	104.00
1	E	25	DG	OP2-P-O3'	-18.55	52.34	108.00
2	D	3	DG	OP1-P-O3'	-17.69	54.94	108.00
1	E	18	DG	OP2-P-O3'	-17.52	55.45	108.00
2	F	13	DT	O3'-P-O5'	-17.35	77.97	104.00
1	C	31	DT	OP2-P-O3'	-17.00	56.98	108.00
1	E	23	DG	O3'-P-O5'	16.46	128.69	104.00
2	F	11	DT	O3'-P-O5'	16.15	128.23	104.00
1	E	17	DC	O3'-P-O5'	15.80	127.70	104.00
1	C	17	DC	OP1-P-O3'	-15.49	61.53	108.00
1	E	30	DA	OP1-P-O3'	-15.22	62.34	108.00
2	D	4	DT	OP1-P-O3'	-15.21	62.38	108.00
2	D	5	DC	OP2-P-O3'	15.09	153.27	108.00
1	C	27	DG	OP2-P-O3'	14.97	152.90	108.00
2	F	10	DT	O3'-P-O5'	14.94	126.40	104.00
1	C	26	DT	OP2-P-O3'	14.78	152.35	108.00
2	D	3	DG	O3'-P-O5'	14.59	125.88	104.00
1	E	26	DT	O3'-P-O5'	-14.57	82.15	104.00
2	D	5	DC	O3'-P-O5'	-14.24	82.64	104.00
1	E	24	DT	OP1-P-O3'	-13.88	66.37	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	7	DC	O3'-P-O5'	-13.61	83.59	104.00
1	C	26	DT	O3'-P-O5'	-13.42	83.87	104.00
1	C	25	DG	O3'-P-O5'	-13.28	84.08	104.00
2	D	4	DT	O3'-P-O5'	-13.18	84.22	104.00
1	C	18	DG	OP1-P-O3'	-13.13	68.60	108.00
1	C	18	DG	OP2-P-O3'	13.01	147.04	108.00
1	C	29	DC	OP2-P-O3'	12.94	146.83	108.00
2	D	12	DT	OP2-P-O3'	-12.69	69.93	108.00
1	E	30	DA	OP2-P-O3'	12.63	145.90	108.00
1	E	22	DA	OP1-P-O3'	-12.52	70.45	108.00
1	C	16	DG	OP2-P-O3'	-12.32	71.03	108.00
2	F	4	DT	OP1-P-O3'	-12.08	71.77	108.00
2	F	12	DT	O3'-P-O5'	-11.93	86.11	104.00
1	E	31	DT	O3'-P-O5'	11.73	121.60	104.00
2	F	10	DT	OP1-P-O3'	-11.44	73.67	108.00
1	E	32	DA	OP1-P-O3'	-11.37	73.88	108.00
1	C	20	DA	OP1-P-O3'	-11.33	74.00	108.00
1	E	29	DC	O3'-P-O5'	-11.05	87.43	104.00
2	F	6	DA	OP2-P-O3'	11.04	141.13	108.00
1	C	26	DT	OP1-P-O3'	-11.02	74.93	108.00
2	D	14	DC	O3'-P-O5'	10.80	120.20	104.00
2	F	6	DA	O3'-P-O5'	-10.79	87.82	104.00
2	F	14	DC	OP2-P-O3'	-10.71	75.86	108.00
1	C	23	DG	OP1-P-O3'	-10.66	76.03	108.00
1	E	22	DA	OP2-P-O3'	10.64	139.92	108.00
1	E	20	DA	O3'-P-O5'	-10.64	88.05	104.00
1	C	17	DC	O3'-P-O5'	10.55	119.82	104.00
2	F	4	DT	OP2-P-O3'	10.17	138.50	108.00
1	C	29	DC	OP1-P-O3'	-9.99	78.02	108.00
2	F	8	DA	OP2-P-O3'	9.99	137.97	108.00
3	A	138	ASP	CA-CB-CG	9.92	122.52	112.60
2	D	5	DC	OP1-P-O3'	-9.85	78.45	108.00
1	E	27	DG	OP2-P-O3'	9.82	137.46	108.00
1	C	32	DA	OP2-P-O3'	9.73	137.18	108.00
1	E	32	DA	OP2-P-O3'	9.70	137.10	108.00
2	D	8	DA	O3'-P-O5'	9.54	118.31	104.00
1	C	27	DG	OP1-P-O3'	-9.51	79.46	108.00
3	A	17	HIS	CA-CB-CG	9.40	123.20	113.80
1	E	20	DA	OP2-P-O3'	9.38	136.16	108.00
1	E	21	DA	OP1-P-O3'	-9.37	79.90	108.00
2	D	9	DC	OP2-P-O3'	-9.34	79.98	108.00
1	E	27	DG	O3'-P-O5'	-9.25	90.12	104.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	14	DC	OP2-P-O3'	-9.18	80.47	108.00
1	C	32	DA	OP1-P-O3'	-8.93	81.22	108.00
1	E	31	DT	OP2-P-O3'	-8.88	81.35	108.00
1	E	29	DC	OP2-P-O3'	8.87	134.62	108.00
1	E	26	DT	C1'-O4'-C4'	-8.80	96.50	109.70
3	B	204	VAL	N-CA-C	-8.79	94.41	107.51
1	E	19	DA	OP1-P-O3'	-8.71	81.86	108.00
2	F	8	DA	OP1-P-O3'	-8.71	81.88	108.00
2	F	11	DT	C1'-O4'-C4'	-8.53	96.91	109.70
3	B	109	ASN	CA-CB-CG	-8.52	104.08	112.60
1	E	19	DA	OP2-P-O3'	-8.50	82.49	108.00
3	B	176	VAL	N-CA-C	8.49	118.52	110.53
1	C	21	DA	O3'-P-O5'	-8.46	91.31	104.00
3	A	76	PHE	CA-CB-CG	-8.43	105.38	113.80
1	C	22	DA	O3'-P-O5'	8.40	116.60	104.00
2	D	13	DT	C1'-O4'-C4'	-8.29	97.27	109.70
3	A	145	GLN	OE1-CD-NE2	-8.23	114.36	122.60
2	F	11	DT	OP1-P-O3'	-8.17	83.50	108.00
2	F	6	DA	OP1-P-O3'	-8.16	83.50	108.00
3	B	43	VAL	N-CA-C	-8.12	102.78	110.42
2	D	10	DT	OP1-P-O3'	-7.83	84.50	108.00
3	A	109	ASN	OD1-CG-ND2	-7.82	114.78	122.60
3	A	155	ASP	CA-CB-CG	7.77	120.37	112.60
3	A	31	HIS	CB-CG-CD2	-7.74	121.14	131.20
3	A	47	VAL	N-CA-CB	-7.70	97.69	111.93
2	D	9	DC	OP1-P-O3'	-7.64	85.08	108.00
2	D	6	DA	OP2-P-O3'	7.56	130.69	108.00
1	C	18	DG	P-O5'-C5'	7.50	131.26	120.00
3	A	34	GLU	CA-CB-CG	7.48	129.06	114.10
2	D	11	DT	C1'-O4'-C4'	-7.39	98.62	109.70
1	C	20	DA	OP2-P-O3'	7.33	129.98	108.00
1	C	17	DC	OP2-P-O3'	7.31	129.93	108.00
1	E	28	DA	P-O5'-C5'	7.30	130.96	120.00
2	F	7	DC	OP2-P-O3'	7.27	129.80	108.00
1	E	26	DT	OP2-P-O3'	7.22	129.65	108.00
1	C	30	DA	P-O3'-C3'	7.21	131.01	120.20
1	C	28	DA	O3'-P-O5'	7.16	114.73	104.00
3	B	31	HIS	CB-CG-CD2	-7.15	121.90	131.20
2	F	3	DG	O3'-P-O5'	7.12	114.68	104.00
3	B	47	VAL	N-CA-CB	-7.07	100.56	112.44
2	F	9	DC	OP2-P-O3'	7.06	129.19	108.00
3	A	97	ILE	CB-CA-C	-7.01	99.47	110.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	9	DC	O3'-P-O5'	-7.00	93.49	104.00
1	C	28	DA	OP1-P-O3'	-6.88	87.37	108.00
1	C	22	DA	P-O5'-C5'	6.84	130.26	120.00
3	B	10	THR	N-CA-C	-6.80	103.87	111.28
1	E	23	DG	OP1-P-O3'	-6.73	87.81	108.00
2	D	13	DT	O3'-P-O5'	-6.68	93.98	104.00
1	C	23	DG	OP2-P-O3'	6.64	127.91	108.00
1	E	18	DG	P-O3'-C3'	6.60	130.10	120.20
2	D	3	DG	OP2-P-O3'	6.57	127.72	108.00
1	C	23	DG	C4'-C3'-O3'	-6.54	100.19	110.00
3	A	93	GLU	N-CA-C	-6.53	97.04	108.20
3	A	194	ASN	OD1-CG-ND2	-6.51	116.09	122.60
2	F	12	DT	OP2-P-O3'	6.46	127.39	108.00
3	A	31	HIS	N-CA-C	-6.46	99.19	109.59
3	A	83	SER	N-CA-C	6.44	120.39	112.54
3	B	122	ARG	CB-CA-C	-6.38	100.86	110.88
2	D	8	DA	OP1-P-O3'	-6.36	88.92	108.00
1	E	28	DA	O3'-P-O5'	-6.34	94.49	104.00
3	B	17	HIS	CB-CG-CD2	-6.30	123.01	131.20
1	C	18	DG	O5'-C5'-C4'	6.27	120.21	110.80
3	B	65	ASN	OD1-CG-ND2	-6.26	116.34	122.60
3	B	65	ASN	CB-CG-ND2	6.25	125.77	116.40
1	E	21	DA	OP2-P-O3'	6.23	126.69	108.00
1	C	31	DT	C1'-O4'-C4'	-6.23	100.36	109.70
3	A	63	TYR	N-CA-C	-6.19	99.60	109.76
1	E	20	DA	OP1-P-O3'	-6.19	89.44	108.00
3	B	166	LYS	N-CA-C	-6.10	99.29	109.24
1	C	30	DA	OP1-P-O3'	-6.10	89.70	108.00
2	F	14	DC	P-O3'-C3'	6.08	129.33	120.20
3	A	112	ILE	CA-C-O	-6.06	114.42	120.85
3	A	65	ASN	OD1-CG-ND2	-6.03	116.57	122.60
3	A	56	GLY	N-CA-C	5.98	122.03	114.25
3	A	73	LEU	N-CA-C	5.97	120.03	112.87
2	D	12	DT	OP1-P-O3'	-5.96	90.11	108.00
3	A	145	GLN	CG-CD-NE2	5.95	125.32	116.40
3	B	119	GLN	CA-CB-CG	-5.92	102.26	114.10
2	D	6	DA	O3'-P-O5'	-5.89	95.16	104.00
3	B	152	LYS	N-CA-CB	-5.89	101.88	110.53
3	B	19	HIS	CB-CG-CD2	-5.84	123.60	131.20
3	B	174	GLN	CG-CD-NE2	5.84	125.16	116.40
3	A	29	LEU	N-CA-C	-5.83	98.38	110.80
3	B	174	GLN	OE1-CD-NE2	-5.79	116.81	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	17	DC	OP2-P-O3'	-5.75	90.76	108.00
3	B	112	ILE	CB-CA-C	-5.74	104.30	112.22
3	B	66	GLN	OE1-CD-NE2	-5.71	116.89	122.60
3	B	31	HIS	CB-CG-ND1	5.71	131.26	122.70
1	E	16	DG	OP1-P-O3'	5.68	125.05	108.00
3	B	23	TYR	CA-C-N	5.67	125.65	120.21
3	B	23	TYR	C-N-CA	5.67	125.65	120.21
3	B	188	LYS	N-CA-C	-5.66	105.03	111.14
2	F	5	DC	O3'-P-O5'	5.64	112.47	104.00
3	A	86	VAL	N-CA-CB	-5.63	104.89	112.10
1	E	18	DG	C4'-C3'-O3'	-5.63	101.55	110.00
1	C	22	DA	P-O3'-C3'	5.63	128.64	120.20
3	B	75	LEU	N-CA-C	5.61	122.74	110.80
3	A	136	PHE	N-CA-C	5.57	120.21	113.41
1	C	29	DC	O3'-P-O5'	-5.56	95.66	104.00
2	D	12	DT	P-O3'-C3'	5.50	128.45	120.20
3	A	166	LYS	N-CA-C	-5.50	99.56	108.52
2	D	6	DA	OP1-P-O3'	-5.49	91.53	108.00
3	B	55	GLU	N-CA-C	-5.45	99.19	110.80
3	A	152	LYS	N-CA-C	-5.43	106.51	114.39
1	C	26	DT	C5'-C4'-C3'	-5.43	106.76	114.90
1	E	22	DA	C1'-O4'-C4'	-5.43	101.56	109.70
3	B	159	HIS	CB-CG-CD2	-5.42	124.16	131.20
1	C	19	DA	P-O3'-C3'	5.41	128.32	120.20
3	A	31	HIS	CB-CG-ND1	5.40	130.80	122.70
1	E	31	DT	N1-C1'-C2'	-5.38	105.42	113.50
3	A	192	ASP	CA-CB-CG	5.37	117.97	112.60
3	B	122	ARG	N-CA-CB	5.37	117.80	110.01
3	B	111	ASP	CA-CB-CG	5.37	117.97	112.60
3	A	157	MET	N-CA-C	-5.37	101.79	110.32
1	C	16	DG	C2'-C3'-O3'	-5.34	103.49	111.50
1	E	16	DG	OP2-P-O3'	-5.33	92.01	108.00
3	A	204	VAL	N-CA-C	-5.32	100.09	108.23
1	E	28	DA	O5'-C5'-C4'	5.32	118.78	110.80
1	C	29	DC	O4'-C1'-N1	5.31	116.36	108.40
1	E	31	DT	P-O3'-C3'	5.31	128.16	120.20
3	B	72	GLU	N-CA-C	5.30	122.09	110.80
3	A	65	ASN	CB-CG-ND2	5.30	124.34	116.40
3	B	161	ASP	N-CA-C	-5.28	96.66	107.70
1	C	30	DA	O5'-C5'-C4'	5.27	118.71	110.80
3	B	112	ILE	N-CA-CB	5.27	118.48	110.58
3	B	12	GLU	CB-CG-CD	5.26	121.55	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	25	DG	OP2-P-O3'	5.26	123.78	108.00
3	B	158	THR	N-CA-CB	-5.26	101.33	109.69
2	F	8	DA	O3'-P-O5'	-5.25	96.13	104.00
3	A	86	VAL	N-CA-C	-5.24	98.43	107.24
3	B	89	LYS	O-C-N	5.23	127.74	122.09
3	A	176	VAL	N-CA-CB	-5.20	101.88	110.56
3	A	146	THR	CA-CB-OG1	-5.19	101.81	109.60
3	A	64	LEU	N-CA-C	-5.19	101.24	109.59
3	B	39	LEU	N-CA-C	-5.17	101.29	109.76
3	B	83	SER	N-CA-C	5.16	118.73	112.23
3	B	57	LYS	CB-CG-CD	5.16	123.16	111.30
3	B	85	TRP	CE2-CD2-CG	-5.15	101.02	107.20
1	E	31	DT	C1'-O4'-C4'	-5.14	101.99	109.70
3	A	47	VAL	CB-CA-C	5.13	118.90	110.69
1	E	27	DG	OP1-P-O3'	-5.13	92.61	108.00
1	E	21	DA	O3'-P-O5'	5.12	111.68	104.00
1	C	28	DA	P-O5'-C5'	5.10	127.64	120.00
3	A	112	ILE	N-CA-CB	5.09	118.21	110.58
2	D	7	DC	O3'-P-O5'	-5.08	96.37	104.00
3	B	85	TRP	CG-CD2-CE3	5.07	138.97	133.90
3	A	15	LEU	N-CA-C	5.04	118.19	111.75
3	B	64	LEU	N-CA-C	-5.02	101.21	109.40
3	B	82	ARG	NE-CZ-NH2	-5.01	114.69	119.20
1	C	31	DT	C2'-C3'-O3'	-5.01	103.99	111.50

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	159	HIS	Peptide
1	C	32	DA	Sidechain
2	D	3	DG	Sidechain
2	F	11	DT	Sidechain
2	F	3	DG	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	372	39	204	6	1
1	E	372	39	204	3	0
2	D	259	25	150	3	0
2	F	259	25	150	2	1
3	A	1556	361	1600	42	0
3	B	1556	361	1600	43	0
4	A	22	3	11	3	0
4	B	22	3	11	3	0
All	All	4418	856	3930	89	1

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 11.

All (89) 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:206:CMP:H2	4:A:206:CMP:C2	0.97	1.49
4:B:206:CMP:H2	4:B:206:CMP:C2	0.97	1.48
3:A:17:HIS:HB2	3:A:101:LYS:HZ3	1.40	0.86
3:B:30:ILE:HG23	3:B:82:ARG:HD3	1.63	0.79
3:A:30:ILE:HG23	3:A:82:ARG:HD3	1.69	0.75
3:A:17:HIS:HB2	3:A:101:LYS:NZ	2.04	0.71
3:A:122:ARG:HH21	3:A:125:GLN:HG3	1.56	0.71
3:A:34:GLU:HB2	3:A:82:ARG:HG3	1.73	0.70
3:A:133:ASN:HA	3:A:137:LEU:HD13	1.72	0.70
3:B:46:SER:HB2	3:B:89:LYS:HE3	1.72	0.69
3:B:122:ARG:HH21	3:B:125:GLN:HG3	1.60	0.66
3:B:18:CYS:SG	3:B:97:ILE:HG23	2.38	0.64
1:C:24:DT:H2'	3:B:180:ARG:HH21	1.63	0.63
1:C:24:DT:H2'	3:B:180:ARG:NH2	2.14	0.62
1:E:23:DG:H5''	3:A:168:THR:HG21	1.81	0.61
3:B:39:LEU:HD23	3:B:97:ILE:HD11	1.82	0.61
3:A:50:LEU:HD12	3:A:85:TRP:HB2	1.83	0.60
4:A:206:CMP:H2	4:A:206:CMP:N3	1.93	0.60
3:B:45:GLY:HA3	3:B:92:CYS:SG	2.42	0.60
3:A:112:ILE:HD12	3:A:113:LEU:H	1.67	0.59
3:B:197:SER:HB2	3:B:203:ILE:HD13	1.85	0.59
3:B:71:GLY:HA2	4:B:206:CMP:O2P	2.04	0.57
3:A:18:CYS:SG	3:A:97:ILE:HG13	2.45	0.57
3:A:72:GLU:HG2	3:A:73:LEU:HD13	1.87	0.57
3:B:122:ARG:HH21	3:B:125:GLN:HE21	1.51	0.56
3:A:70:ILE:HG21	3:A:86:VAL:HG11	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:195:LEU:O	3:A:205:VAL:HG22	2.06	0.56
3:A:121:ALA:HB2	3:B:120:MET:HE1	1.88	0.55
3:B:168:THR:HG23	3:B:171:GLU:H	1.71	0.54
3:B:109:ASN:OD1	3:B:111:ASP:HB3	2.09	0.53
3:B:122:ARG:NH2	3:B:125:GLN:HE21	2.08	0.52
3:B:122:ARG:HH21	3:B:125:GLN:CG	2.23	0.51
3:A:17:HIS:CD2	3:A:17:HIS:N	2.79	0.51
3:B:97:ILE:HD13	3:B:102:PHE:HB2	1.93	0.51
3:A:147:LEU:HD21	3:A:172:ILE:HD13	1.92	0.51
3:A:163:MET:O	3:A:203:ILE:HA	2.12	0.50
3:A:112:ILE:HD12	3:A:113:LEU:N	2.26	0.49
3:A:113:LEU:HD21	3:B:114:MET:SD	2.51	0.49
3:B:195:LEU:O	3:B:205:VAL:HG22	2.12	0.49
3:A:49:VAL:HG11	4:A:206:CMP:C4	2.48	0.49
3:A:179:SER:O	3:A:183:VAL:HG23	2.13	0.48
1:C:31:DT:H2''	1:C:32:DA:N7	2.28	0.48
1:E:24:DT:OP2	3:A:170:GLN:HG2	2.14	0.48
1:C:33:DT:H2''	3:A:182:THR:HG21	1.97	0.47
2:D:14:DC:H2''	2:D:15:DG:C8	2.50	0.46
3:A:155:ASP:O	3:A:157:MET:SD	2.73	0.46
3:B:122:ARG:HH21	3:B:125:GLN:NE2	2.13	0.46
2:D:10:DT:H2''	2:D:11:DT:O4'	2.16	0.46
2:F:14:DC:OP1	3:B:26:LYS:HE2	2.16	0.46
3:A:169:ARG:HH12	3:A:200:GLY:H	1.64	0.46
3:A:9:PRO:HB3	3:A:115:ARG:HH22	1.80	0.46
3:A:52:LYS:NZ	3:A:52:LYS:HB2	2.31	0.46
3:B:40:TYR:HD2	3:B:94:VAL:HG11	1.81	0.46
3:B:50:LEU:O	3:B:85:TRP:HE3	1.99	0.45
3:B:73:LEU:HD22	4:B:206:CMP:H3'	1.98	0.45
2:D:7:DC:H2''	2:D:8:DA:O4'	2.17	0.45
3:A:114:MET:SD	3:B:113:LEU:HD21	2.57	0.45
3:A:136:PHE:HB2	3:A:137:LEU:HD12	1.98	0.45
3:B:103:ARG:O	3:B:106:ILE:HB	2.18	0.44
3:B:122:ARG:NH2	3:B:125:GLN:HG3	2.31	0.44
3:A:101:LYS:O	3:A:104:GLN:HB2	2.18	0.44
3:B:101:LYS:O	3:B:105:LEU:HD13	2.19	0.43
3:A:167:ILE:HD11	3:A:171:GLU:HB3	1.99	0.43
3:B:72:GLU:O	3:B:75:LEU:HB3	2.18	0.43
1:C:24:DT:H5''	3:B:169:ARG:HD3	2.01	0.43
3:B:46:SER:CB	3:B:89:LYS:HE3	2.44	0.43
3:A:113:LEU:CD2	3:B:114:MET:SD	3.07	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:3:DG:H3'	3:B:179:SER:OG	2.19	0.42
3:A:114:MET:SD	3:B:113:LEU:HD11	2.59	0.42
3:A:122:ARG:NH2	3:A:125:GLN:HG3	2.31	0.42
3:A:159:HIS:CG	3:A:203:ILE:HG12	2.54	0.42
3:A:201:LYS:HE3	3:A:203:ILE:HD11	2.00	0.42
3:A:72:GLU:HG2	3:A:73:LEU:N	2.35	0.41
3:B:122:ARG:HH11	3:B:122:ARG:HG2	1.85	0.41
3:B:157:MET:SD	3:B:157:MET:N	2.93	0.41
3:B:32:GLN:HG2	3:B:33:GLY:N	2.35	0.41
1:E:16:DG:H4'	1:E:17:DC:O5'	2.20	0.41
3:B:9:PRO:N	3:B:12:GLU:HG2	2.36	0.41
1:C:29:DC:H2'	1:C:30:DA:C8	2.56	0.41
3:A:22:LYS:HG2	3:A:93:GLU:HG3	2.03	0.41
3:A:29:LEU:HB2	3:A:30:ILE:H	1.79	0.41
3:B:168:THR:HG22	3:B:171:GLU:HG2	2.02	0.41
3:A:101:LYS:HE3	3:A:101:LYS:HB3	1.92	0.40
3:B:49:VAL:HG22	3:B:86:VAL:HG22	2.03	0.40
3:A:32:GLN:HG2	3:A:82:ARG:O	2.21	0.40
3:B:35:LYS:HD3	3:B:35:LYS:H	1.86	0.40
3:B:109:ASN:HA	3:B:110:PRO:HD3	1.84	0.40
3:B:143:ILE:HG12	3:B:176:VAL:HG21	2.03	0.40
3:A:180:ARG:HG3	3:A:181:GLU:N	2.36	0.40

All (1) 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)
1:C:16:DG:H1	2:F:15:DG:O6[3_554]	1.50	0.10

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	
3	A	195/205 (95%)	176 (90%)	15 (8%)	4 (2%)	5	27
3	B	195/205 (95%)	184 (94%)	4 (2%)	7 (4%)	2	16
All	All	390/410 (95%)	360 (92%)	19 (5%)	11 (3%)	4	21

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	155	ASP
3	A	160	PRO
3	B	55	GLU
3	B	75	LEU
3	B	72	GLU
3	B	154	PRO
3	B	199	HIS
3	A	199	HIS
3	B	54	GLU
3	B	155	ASP
3	A	154	PRO

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	
3	A	170/177 (96%)	139 (82%)	31 (18%)	2	9
3	B	170/177 (96%)	142 (84%)	28 (16%)	2	12
All	All	340/354 (96%)	281 (83%)	59 (17%)	2	10

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	11	LEU
3	A	17	HIS
3	A	19	HIS
3	A	29	LEU
3	A	32	GLN

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Mol	Chain	Res	Type
3	A	35	LYS
3	A	37	GLU
3	A	44	LYS
3	A	47	VAL
3	A	50	LEU
3	A	52	LYS
3	A	64	LEU
3	A	73	LEU
3	A	78	GLU
3	A	82	ARG
3	A	87	ARG
3	A	90	THR
3	A	93	GLU
3	A	98	SER
3	A	100	LYS
3	A	101	LYS
3	A	103	ARG
3	A	112	ILE
3	A	125	GLN
3	A	145	GLN
3	A	147	LEU
3	A	153	GLN
3	A	176	VAL
3	A	185	ARG
3	A	189	MET
3	A	194	ASN
3	B	10	THR
3	B	19	HIS
3	B	22	LYS
3	B	29	LEU
3	B	35	LYS
3	B	37	GLU
3	B	38	THR
3	B	46	SER
3	B	47	VAL
3	B	51	ILE
3	B	52	LYS
3	B	55	GLU
3	B	78	GLU
3	B	80	GLN
3	B	90	THR
3	B	96	GLU

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Mol	Chain	Res	Type
3	B	97	ILE
3	B	100	LYS
3	B	105	LEU
3	B	112	ILE
3	B	122	ARG
3	B	124	LEU
3	B	125	GLN
3	B	148	LEU
3	B	185	ARG
3	B	187	LEU
3	B	196	ILE
3	B	204	VAL

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

Mol	Chain	Res	Type
3	A	32	GLN
3	A	125	GLN
3	A	174	GLN
3	B	65	ASN
3	B	80	GLN
3	B	125	GLN
3	B	153	GLN

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 ⓘ

2 ligands are modelled in this entry.

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	206	-	25,25,25	1.29	3 (12%)	37,39,39	1.58	7 (18%)
4	CMP	A	206	-	25,25,25	1.15	2 (8%)	37,39,39	1.51	4 (10%)

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	206	-	-	0/4/31/31	0/4/4/4
4	CMP	A	206	-	-	0/4/31/31	0/4/4/4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	206	CMP	P-O5'	3.71	1.61	1.57
4	A	206	CMP	C5'-C4'	-2.79	1.47	1.51
4	A	206	CMP	O5'-C5'	-2.73	1.42	1.46
4	B	206	CMP	O3'-C3'	-2.53	1.40	1.44
4	B	206	CMP	P-O3'	2.39	1.61	1.57

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	206	CMP	O5'-P-O3'	5.33	112.84	105.70
4	A	206	CMP	O3'-C3'-C4'	-4.30	107.46	110.71
4	A	206	CMP	O2P-P-O1P	3.72	119.99	108.56
4	A	206	CMP	O5'-P-O3'	-3.71	100.73	105.70
4	B	206	CMP	O2P-P-O1P	3.41	119.05	108.56
4	A	206	CMP	C5'-C4'-C3'	-3.29	106.38	112.56
4	B	206	CMP	O3'-C3'-C4'	-3.10	108.37	110.71
4	B	206	CMP	O3'-C3'-C2'	-2.87	112.79	115.61
4	B	206	CMP	O5'-P-O1P	-2.67	104.29	110.44
4	B	206	CMP	C2'-C3'-C4'	-2.31	99.20	103.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	206	CMP	O5'-C5'-C4'	2.04	110.47	105.71

There are no chirality outliers.

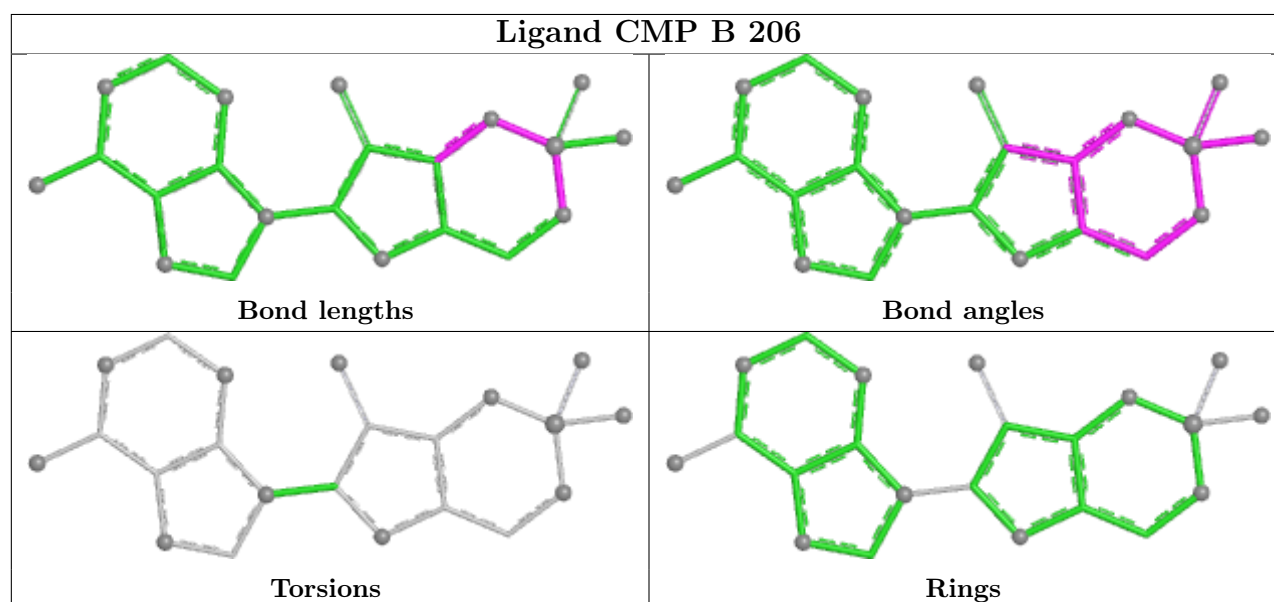
There are no torsion outliers.

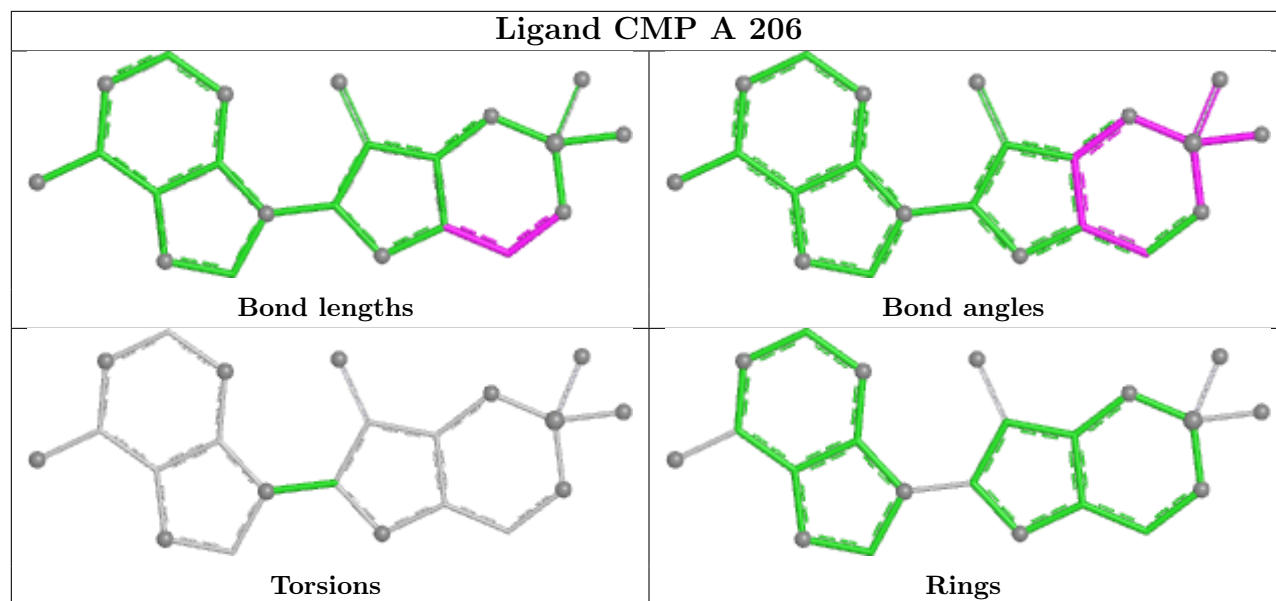
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	206	CMP	3	0
4	A	206	CMP	3	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.