



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:33 AM UTC

PDB ID : 1ZRD / pdb_00001zrd
Title : 4 crystal structures of CAP-DNA with all base-pair substitutions at position 6, CAP-[6A;17T]ICAP38 DNA
Authors : Berman, H.M.; Napoli, A.A.
Deposited on : 2005-05-19
Resolution : 2.80 Å(reported)

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

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

A user guide is available at

<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

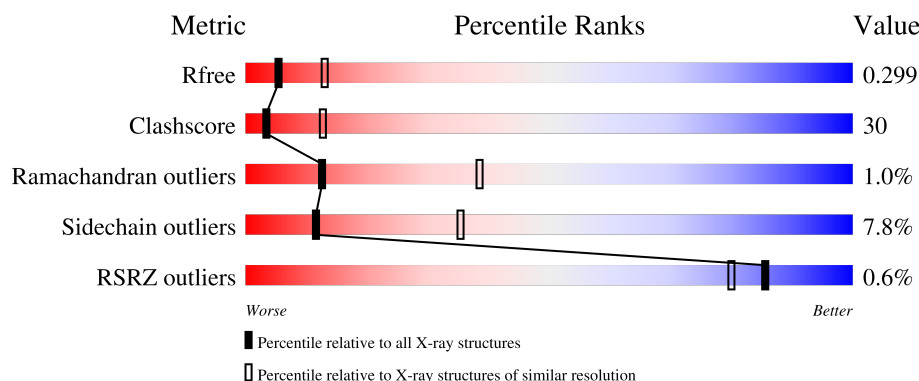
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	W	17	100%
1	Y	17	6% 94%
2	X	21	24% 76%
2	Z	21	5% 14% 86%
3	A	209	56% 34% 5% .

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Mol	Chain	Length	Quality of chain
3	B	209	48% 39% 8%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4781 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 DNA chain called 5'-D(*AP*TP*TP*TP*CP*GP*AP*AP*AP*AP*T
P*GP*AP*GP*AP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	W	17	Total	C	N	O	P	0	0	0
			350	169	68	97	16			
1	Y	17	Total	C	N	O	P	0	0	0
			350	169	68	97	16			

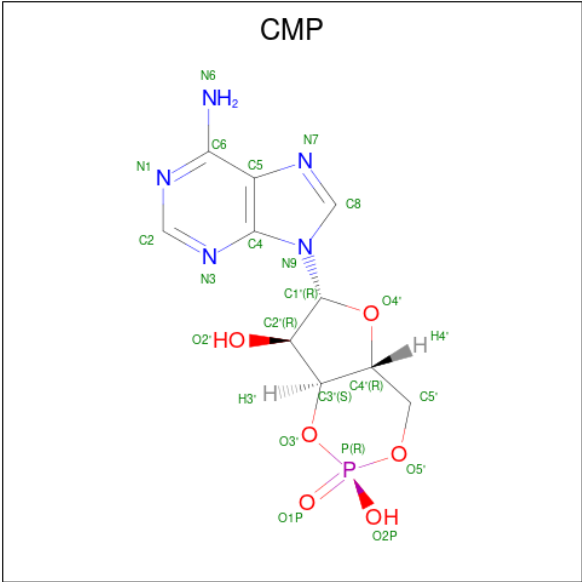
- Molecule 2 is a DNA chain called 5'-D(*CP*TP*AP*GP*AP*TP*CP*TP*CP*AP*TP*TP
*TP*TP*TP*CP*GP*AP*AP*AP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	X	21	Total	C	N	O	P	0	0	0
			423	206	70	127	20			
2	Z	21	Total	C	N	O	P	0	0	0
			423	206	70	127	20			

- Molecule 3 is a protein called Catabolite gene activator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	200	Total	C	N	O	S	0	0	0
			1580	1001	276	294	9			
3	B	200	Total	C	N	O	S	0	0	0
			1580	1001	276	294	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	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		

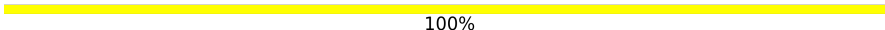
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	W	2	Total	O	0	0
			2	2		
5	X	3	Total	O	0	0
			3	3		
5	Z	1	Total	O	0	0
			1	1		
5	A	16	Total	O	0	0
			16	16		
5	B	9	Total	O	0	0
			9	9		

3 Residue-property plots

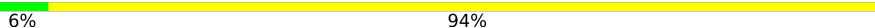
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 5'-D(*AP*TP*TP*TP*CP*GP*AP*AP*AP*AP*TP*GP*AP*GP*AP*T)-3',

Chain W:  100%

A-8
T-7
T-6
T-5
C-4
G-3
A-2
A-1
A1
A2
A3
T4
G5
A6
G7
A8
T9

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

Chain Y:  6% 94%

A-8
T-7
T-6
T-5
C-4
G-3
A-2
A-1
A1
A2
A3
T4
G5
A6
G7
A8
T9

- Molecule 2: 5'-D(*CP*TP*AP*GP*AP*TP*CP*TP*CP*AP*TP*TP*TP*TP*CP*GP*AP*AP*AP*T)-3'

Chain X:  24% 76%

C13
T12
A11
G10
A9
T8
C7
T6
C5
A4
T3
T2
T1
T-2
A-6
A-7
T-8

- Molecule 2: 5'-D(*CP*TP*AP*GP*AP*TP*CP*TP*CP*AP*TP*TP*TP*TP*CP*GP*AP*AP*AP*T)-3'

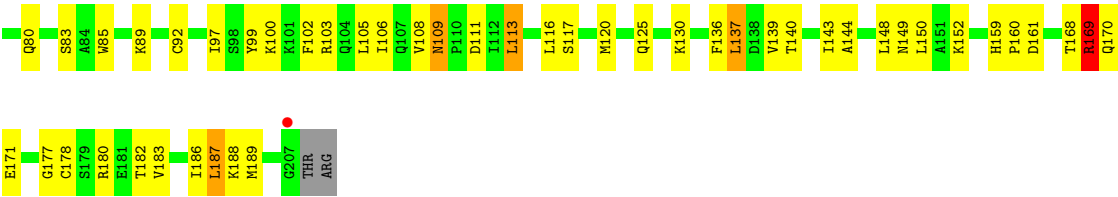
Chain Z:  5% 14% 86%

C13
T12
A11
G10
A9
T8
C7
T6
C5
A4
T3
T2
T1
T-2
C-3
G-4
A-5
A-6
A-7
T-8

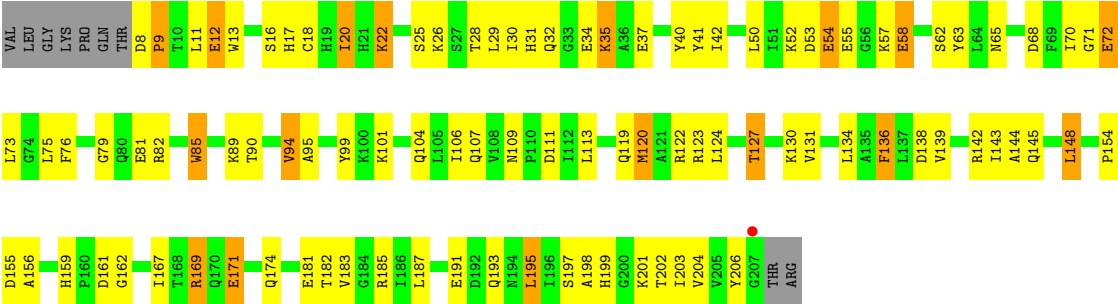
- Molecule 3: Catabolite gene activator

Chain A:  56% 34% 5%

VAL
LEU
GLY
LYS
PRO
GLN
THR
D8
P9
T10
L11
E12
M13
F14
L15
S16
H17
C18
H19
I20
Y23
P24
S25
K26
L29
I30
A36
E37
T38
L39
Y40
Y41
I42
S46
V47
L50
I51
K52
D53
K57
E58
N59
Y63
L64
N65
F69
I70
G71
E72
L73
G74
L75



● Molecule 3: Catabolite gene activator



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.29Å 75.34Å 178.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.63 – 2.80 46.63 – 2.80	Depositor EDS
% Data completeness (in resolution range)	77.9 (46.63-2.80) 77.9 (46.63-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.77Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.224 , 0.297 0.229 , 0.299	Depositor DCC
R_{free} test set	1599 reflections (8.93%)	wwPDB-VP
Wilson B-factor (Å ²)	67.3	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 57.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4781	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% 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

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	W	0.29	0/394	0.81	0/607
1	Y	0.28	0/394	0.73	0/607
2	X	0.31	0/472	0.76	0/726
2	Z	0.31	0/472	0.75	0/726
3	A	0.56	0/1605	1.07	7/2162 (0.3%)
3	B	0.47	0/1605	1.02	10/2162 (0.5%)
All	All	0.46	0/4942	0.95	17/6990 (0.2%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	73	LEU	N-CA-C	9.49	121.62	111.28
3	B	94	VAL	N-CA-C	7.25	118.26	108.11
3	B	136	PHE	N-CA-C	7.02	119.01	111.36
3	B	195	LEU	N-CA-C	-6.09	106.82	114.56
3	B	58	GLU	N-CA-C	5.86	119.03	109.24
3	B	120	MET	N-CA-C	-5.83	104.84	111.14
3	A	169	ARG	N-CA-C	-5.80	105.04	111.36
3	B	85	TRP	N-CA-C	-5.60	101.74	110.42
3	A	50	LEU	N-CA-C	5.53	117.47	109.07
3	A	137	LEU	N-CA-C	5.49	118.51	109.72
3	A	75	LEU	N-CA-C	5.33	119.27	112.87
3	B	169	ARG	N-CA-C	-5.31	105.60	111.71
3	A	20	ILE	N-CA-C	5.31	116.43	109.21
3	B	50	LEU	N-CA-C	5.30	117.32	108.99
3	A	125	GLN	N-CA-C	-5.27	105.43	111.07
3	B	35	LYS	N-CA-C	-5.15	102.66	110.28
3	A	89	LYS	N-CA-C	-5.00	104.83	112.04

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	W	350	0	194	26	0
1	Y	350	0	194	20	0
2	X	423	0	242	28	0
2	Z	423	0	242	40	0
3	A	1580	0	1615	76	0
3	B	1580	0	1615	89	0
4	A	22	0	11	4	0
4	B	22	0	11	3	0
5	A	16	0	0	1	0
5	B	9	0	0	0	0
5	W	2	0	0	0	0
5	X	3	0	0	0	0
5	Z	1	0	0	0	0
All	All	4781	0	4124	266	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 30.

All (266) 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:210:CMP:H2	4:A:210:CMP:C2	0.97	1.49
4:B:210:CMP:C2	4:B:210:CMP:H2	0.97	1.48
2:X:11:DA:H2''	2:X:10:DG:H5''	1.08	1.08
2:X:11:DA:C2'	2:X:10:DG:H5''	1.94	0.96
3:B:42:ILE:HD11	3:B:70:ILE:HD11	1.50	0.92
3:A:116:LEU:HD13	3:A:120:MET:HE2	1.53	0.90
2:X:2:DT:H2''	2:X:1:DT:H5''	1.54	0.89
3:A:113:LEU:HD13	3:B:113:LEU:HD22	1.52	0.87
2:X:2:DT:H2''	2:X:1:DT:C5'	2.04	0.86
2:X:6:DT:H2''	2:X:5:DC:H5''	1.55	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:104:GLN:HA	3:B:107:GLN:HE21	1.40	0.86
3:B:106:ILE:HD13	3:B:113:LEU:HD12	1.58	0.84
3:A:186:ILE:HA	3:A:189:MET:HE2	1.60	0.84
2:X:4:DA:H1'	2:X:3:DT:H5''	1.59	0.83
3:A:42:ILE:HD11	3:A:70:ILE:HD11	1.60	0.83
1:W:-7:DT:H2''	1:W:-6:DT:H71	1.59	0.83
2:X:5:DC:H2''	2:X:4:DA:C8	2.14	0.82
2:X:11:DA:H2''	2:X:10:DG:C5'	2.03	0.80
3:A:40:TYR:HB2	3:A:70:ILE:HB	1.67	0.77
3:A:73:LEU:HD21	3:B:124:LEU:HD23	1.65	0.77
3:A:144:ALA:O	3:A:148:LEU:HD13	1.85	0.77
2:Z:8:DT:H2''	2:Z:7:DC:H5'	1.67	0.76
2:X:-6:DA:H2''	2:X:-7:DA:OP2	1.85	0.76
2:Z:-5:DA:H2''	2:Z:-6:DA:H5'	1.67	0.75
1:W:1:DA:H1'	1:W:2:DA:H5''	1.67	0.75
3:B:35:LYS:HZ2	3:B:79:GLY:HA2	1.50	0.75
3:B:31:HIS:HB2	3:B:34:GLU:HG3	1.69	0.74
3:B:31:HIS:O	3:B:34:GLU:HB2	1.89	0.72
3:B:17:HIS:HB3	3:B:101:LYS:NZ	2.03	0.72
3:A:159:HIS:CD2	3:A:161:ASP:H	2.07	0.71
2:Z:9:DA:H2''	2:Z:8:DT:H5'	1.71	0.71
2:Z:7:DC:H2''	2:Z:6:DT:H5'	1.71	0.71
1:W:-7:DT:H2''	1:W:-6:DT:C7	2.19	0.71
2:Z:6:DT:H2''	2:Z:5:DC:C5'	2.21	0.70
3:A:52:LYS:HD2	3:A:85:TRP:NE1	2.07	0.70
1:W:4:DT:P	3:A:169:ARG:HD2	2.34	0.68
1:W:8:DA:H1'	1:W:9:DT:H5''	1.75	0.68
3:B:54:GLU:OE2	3:B:55:GLU:HG3	1.93	0.68
2:Z:4:DA:H1'	2:Z:3:DT:H5''	1.76	0.67
3:B:197:SER:HB3	3:B:204:VAL:HB	1.76	0.67
3:B:104:GLN:HA	3:B:107:GLN:NE2	2.09	0.67
2:Z:6:DT:H2''	2:Z:5:DC:H5''	1.77	0.66
1:W:4:DT:OP1	3:A:169:ARG:HD2	1.95	0.66
1:W:-4:DC:H2''	1:W:-3:DG:OP2	1.95	0.66
1:W:-1:DA:H2''	1:W:1:DA:OP2	1.96	0.66
3:B:159:HIS:HD2	3:B:161:ASP:HB2	1.60	0.66
3:B:104:GLN:CA	3:B:107:GLN:HE21	2.08	0.65
3:B:109:ASN:HD21	3:B:111:ASP:HB3	1.61	0.65
2:X:2:DT:H2''	2:X:1:DT:H5'	1.77	0.65
2:Z:-2:DT:H2''	2:Z:-3:DC:H5'	1.78	0.65
3:B:119:GLN:O	3:B:123:ARG:HG3	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:5:DG:H4'	1:W:6:DA:OP1	1.97	0.64
3:B:181:GLU:HB3	3:B:185:ARG:NH1	2.12	0.64
3:A:17:HIS:ND1	3:A:105:LEU:HD21	2.12	0.64
3:A:116:LEU:CD1	3:A:120:MET:HE2	2.25	0.64
1:Y:-5:DT:H1'	1:Y:-4:DC:H5'	1.81	0.63
2:Z:11:DA:H2''	2:Z:10:DG:H5'	1.81	0.63
2:X:13:DC:H5'	1:Y:9:DT:O3'	1.99	0.63
3:A:102:PHE:O	3:A:106:ILE:HG13	1.99	0.63
2:Z:2:DT:H2'	2:Z:1:DT:H72	1.79	0.63
3:A:108:VAL:O	3:A:109:ASN:HB2	1.99	0.61
1:Y:3:DA:H1'	1:Y:4:DT:H5''	1.82	0.61
2:Z:1:DT:H2''	2:Z:-1:DT:H5'	1.82	0.61
3:B:37:GLU:HA	3:B:99:TYR:CD2	2.36	0.61
2:Z:9:DA:H2''	2:Z:8:DT:C5'	2.31	0.60
3:B:18:CYS:HB3	3:B:95:ALA:HB1	1.83	0.60
3:A:63:TYR:OH	3:A:171:GLU:OE2	2.18	0.60
3:A:188:LYS:HD3	5:A:224:HOH:O	2.01	0.60
2:X:1:DT:H1'	2:X:-1:DT:H5''	1.84	0.60
3:A:72:GLU:HG2	3:A:116:LEU:HD21	1.82	0.60
2:X:2:DT:C2'	2:X:1:DT:H5''	2.28	0.60
3:A:59:MET:HE3	3:B:131:VAL:HG12	1.82	0.60
3:B:40:TYR:HB2	3:B:70:ILE:HB	1.84	0.60
2:Z:2:DT:H2''	2:Z:1:DT:C5'	2.32	0.59
2:X:1:DT:H2''	2:X:-1:DT:H5'	1.84	0.59
1:W:-6:DT:H1'	1:W:-5:DT:H5'	1.83	0.59
3:A:59:MET:HB2	3:B:136:PHE:CZ	2.36	0.59
3:B:25:SER:O	3:B:26:LYS:HB2	2.02	0.59
3:A:83:SER:OG	4:A:210:CMP:H5'1	2.03	0.59
3:B:35:LYS:NZ	3:B:79:GLY:HA2	2.18	0.59
1:W:-6:DT:H2''	1:W:-5:DT:OP2	2.03	0.58
2:Z:6:DT:C2'	2:Z:5:DC:H5''	2.34	0.58
3:B:53:ASP:CG	3:B:57:LYS:HB3	2.29	0.58
3:B:144:ALA:O	3:B:148:LEU:HD13	2.03	0.58
3:A:53:ASP:OD1	3:A:57:LYS:HB2	2.04	0.58
1:W:4:DT:H2''	1:W:5:DG:C8	2.39	0.58
3:B:113:LEU:HD23	3:B:113:LEU:O	2.04	0.58
2:X:9:DA:H2''	2:X:8:DT:H5'	1.86	0.57
1:Y:-1:DA:H2''	1:Y:1:DA:H5'	1.86	0.57
1:Y:7:DG:H2''	1:Y:8:DA:OP2	2.03	0.57
3:A:99:TYR:O	3:A:103:ARG:HG3	2.05	0.57
2:Z:10:DG:H2''	2:Z:9:DA:OP2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:-6:DT:H2''	1:Y:-5:DT:H71	1.87	0.57
3:A:42:ILE:CD1	3:A:70:ILE:HD11	2.34	0.57
3:A:11:LEU:O	3:A:14:PHE:HB3	2.05	0.57
2:Z:1:DT:H1'	2:Z:-1:DT:H5''	1.87	0.56
2:Z:1:DT:H2''	2:Z:-1:DT:C5'	2.35	0.56
2:Z:-3:DC:H2''	2:Z:-4:DG:C8	2.40	0.56
3:A:52:LYS:HD2	3:A:85:TRP:CD1	2.41	0.56
3:A:100:LYS:HE3	3:A:103:ARG:HH11	1.71	0.56
1:W:8:DA:H1'	1:W:9:DT:C5'	2.35	0.55
3:A:183:VAL:O	3:A:187:LEU:HB2	2.07	0.55
2:Z:-4:DG:H2''	2:Z:-5:DA:OP2	2.07	0.55
3:A:100:LYS:HE3	3:A:100:LYS:HA	1.89	0.55
3:B:31:HIS:HD1	3:B:85:TRP:HZ3	1.55	0.55
3:B:53:ASP:OD1	3:B:57:LYS:HB3	2.07	0.55
3:B:202:THR:C	3:B:203:ILE:HD12	2.32	0.55
3:B:76:PHE:CE1	3:B:120:MET:HE1	2.43	0.54
2:Z:6:DT:H2''	2:Z:5:DC:H5'	1.88	0.54
2:X:3:DT:H2''	2:X:2:DT:O5'	2.08	0.54
3:B:104:GLN:HA	3:B:107:GLN:HG3	1.88	0.54
3:A:53:ASP:CG	3:A:57:LYS:HB2	2.33	0.54
1:Y:-6:DT:C2'	1:Y:-5:DT:H71	2.36	0.54
2:Z:-1:DT:C2'	2:Z:-2:DT:H71	2.38	0.54
3:B:29:LEU:O	3:B:30:ILE:HG13	2.07	0.54
3:B:142:ARG:HG3	3:B:142:ARG:HH11	1.73	0.54
2:X:6:DT:C2'	2:X:5:DC:H5''	2.34	0.54
1:Y:-7:DT:H2''	1:Y:-6:DT:OP2	2.08	0.54
1:Y:3:DA:C2'	1:Y:4:DT:H5''	2.37	0.54
3:B:20:ILE:HA	3:B:95:ALA:HA	1.89	0.54
3:B:145:GLN:HA	3:B:148:LEU:HD22	1.90	0.54
3:A:69:PHE:CG	3:A:116:LEU:HD23	2.42	0.53
1:Y:3:DA:H2''	1:Y:4:DT:H5''	1.90	0.53
2:Z:-6:DA:H1'	2:Z:-7:DA:H5'	1.91	0.53
1:W:-7:DT:H4'	1:W:-6:DT:OP1	2.09	0.53
2:Z:2:DT:H2''	2:Z:1:DT:H5''	1.89	0.53
3:B:109:ASN:HD21	3:B:111:ASP:CB	2.23	0.52
3:A:20:ILE:HD12	3:A:20:ILE:N	2.25	0.52
3:B:109:ASN:ND2	3:B:111:ASP:HB3	2.24	0.52
2:Z:-5:DA:H1'	2:Z:-6:DA:H5''	1.93	0.51
3:A:105:LEU:HA	3:A:108:VAL:HG22	1.91	0.51
3:A:109:ASN:OD1	3:A:111:ASP:HB3	2.09	0.51
3:B:197:SER:HB2	3:B:206:TYR:OH	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:52:LYS:HG2	3:B:58:GLU:HB3	1.93	0.50
1:W:2:DA:H1'	1:W:3:DA:H5'	1.93	0.50
2:Z:9:DA:OP2	3:B:182:THR:OG1	2.27	0.50
2:Z:3:DT:H6	2:Z:3:DT:H5'	1.77	0.50
3:A:36:ALA:O	3:A:37:GLU:HG2	2.12	0.49
2:Z:-1:DT:H2''	2:Z:-2:DT:H71	1.94	0.49
3:B:63:TYR:CE1	3:B:89:LYS:HD3	2.47	0.49
3:A:25:SER:O	3:A:26:LYS:HB2	2.12	0.49
2:X:9:DA:OP1	3:A:178:CYS:HA	2.12	0.49
1:Y:5:DG:H2''	1:Y:6:DA:H5'	1.94	0.49
3:B:113:LEU:HD23	3:B:113:LEU:C	2.38	0.49
3:A:159:HIS:HD2	3:A:161:ASP:H	1.56	0.49
1:W:8:DA:H2''	1:W:9:DT:H5'	1.94	0.48
2:Z:2:DT:H2''	2:Z:1:DT:H5'	1.95	0.48
2:Z:-5:DA:H2''	2:Z:-6:DA:C5'	2.41	0.48
3:B:76:PHE:CZ	3:B:120:MET:HE1	2.49	0.48
1:Y:-8:DA:H2''	1:Y:-7:DT:C6	2.48	0.48
3:B:72:GLU:OE2	4:B:210:CMP:H2'	2.14	0.48
3:B:13:TRP:CH2	3:B:109:ASN:HB3	2.48	0.48
3:A:75:LEU:HG	3:A:75:LEU:O	2.13	0.48
3:B:12:GLU:O	3:B:12:GLU:OE1	2.31	0.48
3:B:22:LYS:NZ	3:B:22:LYS:HB3	2.29	0.48
2:X:1:DT:H2''	2:X:-1:DT:C5'	2.44	0.48
3:B:159:HIS:O	3:B:162:GLY:O	2.32	0.48
3:B:191:GLU:OE1	3:B:198:ALA:HB3	2.13	0.48
3:B:142:ARG:HG3	3:B:142:ARG:NH1	2.28	0.48
1:Y:8:DA:H1'	1:Y:9:DT:H5'	1.96	0.47
3:A:52:LYS:HA	3:A:57:LYS:O	2.14	0.47
3:A:46:SER:HB2	3:A:65:ASN:OD1	2.14	0.47
1:W:3:DA:H1'	1:W:4:DT:H5''	1.95	0.47
2:Z:9:DA:H1'	2:Z:8:DT:H5''	1.97	0.47
3:A:11:LEU:HD21	3:A:41:TYR:OH	2.13	0.47
1:W:4:DT:OP1	3:A:168:THR:HB	2.14	0.47
2:Z:3:DT:C6	2:Z:2:DT:H72	2.49	0.47
3:A:18:CYS:SG	3:A:97:ILE:HB	2.54	0.47
3:A:149:ASN:O	3:A:152:LYS:N	2.34	0.47
1:W:1:DA:H1'	1:W:2:DA:C5'	2.40	0.47
1:Y:3:DA:H2''	1:Y:4:DT:C5'	2.44	0.47
3:B:8:ASP:HB2	3:B:9:PRO:CD	2.45	0.47
2:X:13:DC:H5'	1:Y:9:DT:H2''	1.98	0.46
1:Y:2:DA:H1'	1:Y:3:DA:H5'	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:75:LEU:HA	3:A:99:TYR:CE2	2.50	0.46
3:A:180:ARG:HH11	3:A:180:ARG:HG3	1.81	0.46
3:B:154:PRO:C	3:B:156:ALA:H	2.23	0.46
3:B:71:GLY:HA2	4:B:210:CMP:O3'	2.15	0.46
3:B:65:ASN:O	3:B:68:ASP:HB2	2.16	0.46
3:B:72:GLU:H	3:B:72:GLU:HG3	1.48	0.46
3:A:83:SER:HG	4:A:210:CMP:H5'1	1.81	0.46
3:B:193:GLN:HE21	3:B:195:LEU:HD12	1.80	0.46
3:B:193:GLN:HE21	3:B:195:LEU:CD1	2.28	0.46
1:W:4:DT:O5'	3:A:169:ARG:HD2	2.16	0.46
2:Z:-1:DT:H2'	2:Z:-2:DT:H71	1.98	0.46
3:B:193:GLN:HB2	3:B:195:LEU:HD12	1.97	0.46
3:A:102:PHE:CE2	3:A:106:ILE:HD11	2.51	0.45
3:A:159:HIS:CG	3:A:160:PRO:HD2	2.51	0.45
2:Z:3:DT:H2'	2:Z:2:DT:H72	1.99	0.45
2:X:7:DC:H2''	2:X:6:DT:OP2	2.17	0.45
3:B:29:LEU:N	3:B:29:LEU:HD12	2.32	0.45
3:B:31:HIS:HB2	3:B:34:GLU:CG	2.44	0.45
3:B:62:SER:OG	3:B:63:TYR:N	2.49	0.44
3:A:159:HIS:ND1	3:A:160:PRO:HD2	2.32	0.44
3:B:53:ASP:OD1	3:B:57:LYS:N	2.50	0.44
3:A:10:THR:HG21	3:A:111:ASP:OD2	2.17	0.44
3:A:102:PHE:CD2	3:A:106:ILE:HD11	2.53	0.44
3:A:106:ILE:HG23	3:A:113:LEU:HB2	1.99	0.44
3:B:32:GLN:C	3:B:34:GLU:H	2.26	0.44
3:B:40:TYR:CD1	3:B:40:TYR:N	2.85	0.44
3:B:169:ARG:HG2	3:B:183:VAL:HG12	1.99	0.44
1:Y:-7:DT:C6	1:Y:-6:DT:H72	2.52	0.44
1:W:1:DA:H2''	1:W:2:DA:OP2	2.17	0.44
3:B:183:VAL:O	3:B:187:LEU:HD13	2.17	0.44
2:X:4:DA:C1'	2:X:3:DT:H5''	2.40	0.44
3:B:75:LEU:O	3:B:75:LEU:HG	2.17	0.44
3:B:167:ILE:HD11	3:B:171:GLU:HB3	1.99	0.44
3:A:136:PHE:O	3:A:137:LEU:HD23	2.18	0.44
3:B:25:SER:HB2	3:B:90:THR:C	2.43	0.44
2:Z:-2:DT:H1'	2:Z:-3:DC:H5''	2.00	0.44
3:A:53:ASP:OD2	3:A:53:ASP:C	2.62	0.43
3:A:53:ASP:OD2	3:A:57:LYS:N	2.51	0.43
3:B:11:LEU:HD21	3:B:41:TYR:OH	2.17	0.43
3:A:117:SER:HA	3:A:120:MET:HE3	2.01	0.43
1:Y:-3:DG:N2	2:Z:-4:DG:N2	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Z:11:DA:H2''	2:Z:10:DG:C5'	2.48	0.43
3:A:136:PHE:CG	3:B:53:ASP:HB3	2.53	0.43
3:B:201:LYS:O	3:B:201:LYS:HG2	2.19	0.43
2:X:-1:DT:H2''	2:X:-2:DT:C6	2.53	0.43
3:B:58:GLU:HG3	3:B:174:GLN:NE2	2.34	0.43
3:B:127:THR:O	3:B:130:LYS:HB3	2.19	0.43
1:W:-8:DA:H2''	1:W:-7:DT:OP2	2.19	0.43
2:X:5:DC:H2''	2:X:4:DA:N7	2.32	0.43
3:A:139:VAL:O	3:A:143:ILE:HG13	2.19	0.43
3:A:180:ARG:HG3	3:A:180:ARG:NH1	2.34	0.43
3:B:70:ILE:HG22	3:B:71:GLY:N	2.34	0.43
1:W:8:DA:H2''	1:W:9:DT:OP2	2.17	0.42
3:B:41:TYR:O	3:B:94:VAL:HG13	2.19	0.42
3:B:13:TRP:O	3:B:16:SER:HB2	2.19	0.42
3:B:169:ARG:CG	3:B:187:LEU:HD22	2.49	0.42
3:B:8:ASP:HB2	3:B:9:PRO:HD2	2.00	0.42
3:B:134:LEU:HA	3:B:142:ARG:HD3	2.01	0.42
3:B:139:VAL:O	3:B:143:ILE:HG13	2.19	0.42
3:B:169:ARG:HG2	3:B:187:LEU:HD22	2.01	0.42
1:W:7:DG:H1'	1:W:8:DA:C8	2.55	0.42
2:X:6:DT:H6	2:X:6:DT:H2'	1.67	0.42
3:A:143:ILE:HD12	3:A:186:ILE:HD12	2.02	0.42
3:B:29:LEU:C	3:B:30:ILE:HG13	2.45	0.42
3:B:159:HIS:CD2	3:B:204:VAL:HG11	2.54	0.42
3:A:63:TYR:HE2	3:A:150:LEU:HD21	1.85	0.42
3:B:104:GLN:CB	3:B:107:GLN:HE21	2.32	0.42
1:W:-3:DG:H2''	1:W:-2:DA:O5'	2.20	0.42
2:Z:11:DA:H1'	2:Z:10:DG:H5''	2.02	0.41
1:Y:5:DG:H1'	1:Y:6:DA:H5''	2.01	0.41
3:A:29:LEU:C	3:A:30:ILE:HG13	2.45	0.41
3:B:109:ASN:ND2	3:B:111:ASP:H	2.18	0.41
1:W:3:DA:C8	1:W:4:DT:H72	2.56	0.41
2:X:13:DC:H6	2:X:13:DC:O5'	2.04	0.41
3:A:83:SER:OG	4:A:210:CMP:H8	2.21	0.41
2:Z:-3:DC:H6	2:Z:-3:DC:H2'	1.76	0.41
3:A:23:TYR:CD2	3:A:29:LEU:HG	2.55	0.41
3:A:149:ASN:O	3:A:150:LEU:C	2.63	0.41
1:Y:4:DT:H2''	1:Y:5:DG:C8	2.55	0.41
2:Z:2:DT:H2'	2:Z:1:DT:C7	2.48	0.41
3:A:106:ILE:O	3:A:106:ILE:HG22	2.20	0.41
3:A:13:TRP:O	3:A:16:SER:OG	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:39:LEU:HB3	3:A:97:ILE:CG2	2.51	0.40
3:A:47:VAL:HG22	3:A:64:LEU:HB2	2.03	0.40
3:B:30:ILE:CG2	3:B:82:ARG:HD2	2.51	0.40
2:X:10:DG:P	3:A:139:VAL:HG23	2.62	0.40
2:X:2:DT:H2'	2:X:1:DT:H72	2.03	0.40
2:Z:4:DA:C1'	2:Z:3:DT:H5''	2.48	0.40
3:A:69:PHE:CD1	3:A:116:LEU:HD23	2.56	0.40
3:A:57:LYS:HB3	3:A:57:LYS:HE2	1.86	0.40
3:B:35:LYS:NZ	3:B:79:GLY:CA	2.84	0.40
3:A:143:ILE:CD1	3:A:186:ILE:HD12	2.51	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	
3	A	198/209 (95%)	183 (92%)	12 (6%)	3 (2%)	8	28
3	B	198/209 (95%)	178 (90%)	19 (10%)	1 (0%)	24	55
All	All	396/418 (95%)	361 (91%)	31 (8%)	4 (1%)	12	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	109	ASN
3	B	9	PRO
3	A	177	GLY
3	A	10	THR

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	172/180 (96%)	159 (92%)	13 (8%)	12	36
3	B	172/180 (96%)	158 (92%)	14 (8%)	11	33
All	All	344/360 (96%)	317 (92%)	27 (8%)	11	35

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	11	LEU
3	A	51	ILE
3	A	59	MET
3	A	72	GLU
3	A	80	GLN
3	A	92	CYS
3	A	113	LEU
3	A	130	LYS
3	A	140	THR
3	A	169	ARG
3	A	170	GLN
3	A	182	THR
3	A	187	LEU
3	B	12	GLU
3	B	20	ILE
3	B	22	LYS
3	B	28	THR
3	B	54	GLU
3	B	72	GLU
3	B	81	GLU
3	B	122	ARG
3	B	127	THR
3	B	138	ASP
3	B	148	LEU
3	B	155	ASP
3	B	171	GLU
3	B	199	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	80	GLN
3	A	107	GLN
3	A	149	ASN
3	A	159	HIS
3	A	170	GLN
3	B	17	HIS
3	B	65	ASN
3	B	104	GLN
3	B	107	GLN
3	B	109	ASN
3	B	125	GLN
3	B	133	ASN
3	B	145	GLN
3	B	149	ASN
3	B	159	HIS
3	B	170	GLN
3	B	174	GLN
3	B	193	GLN
3	B	199	HIS

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	210	-	25,25,25	0.90	2 (8%)	37,39,39	1.74	5 (13%)
4	CMP	A	210	-	25,25,25	0.95	2 (8%)	37,39,39	1.54	3 (8%)

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	210	-	-	0/4/31/31	0/4/4/4
4	CMP	A	210	-	-	0/4/31/31	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	210	CMP	P-O3'	2.57	1.62	1.57
4	B	210	CMP	O3'-C3'	-2.41	1.41	1.44
4	A	210	CMP	P-O5'	2.15	1.60	1.57
4	B	210	CMP	P-O5'	2.11	1.60	1.57

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	210	CMP	O3'-C3'-C4'	-7.59	104.98	110.71
4	A	210	CMP	O3'-C3'-C4'	-7.08	105.37	110.71
4	B	210	CMP	O3'-C3'-C2'	4.74	120.25	115.61
4	A	210	CMP	O2P-P-O1P	3.56	119.51	108.56
4	B	210	CMP	O2P-P-O1P	3.48	119.26	108.56
4	A	210	CMP	O3'-C3'-C2'	3.21	118.75	115.61
4	B	210	CMP	O5'-P-O3'	-2.30	102.63	105.70
4	B	210	CMP	O5'-P-O1P	-2.10	105.59	110.44

There are no chirality outliers.

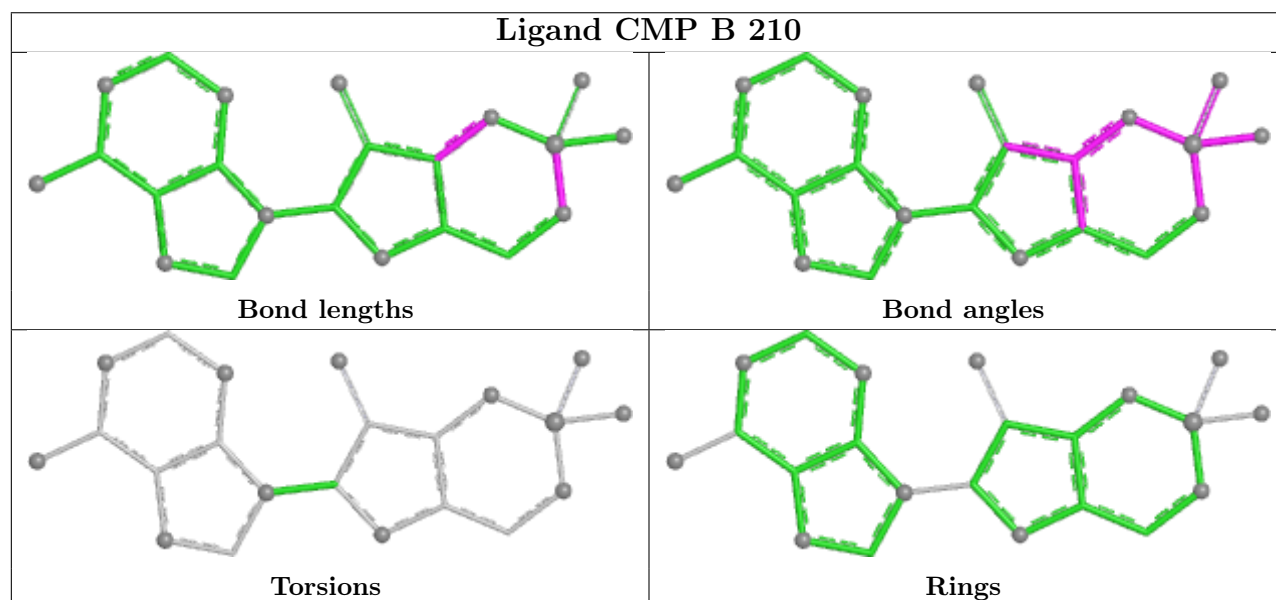
There are no torsion outliers.

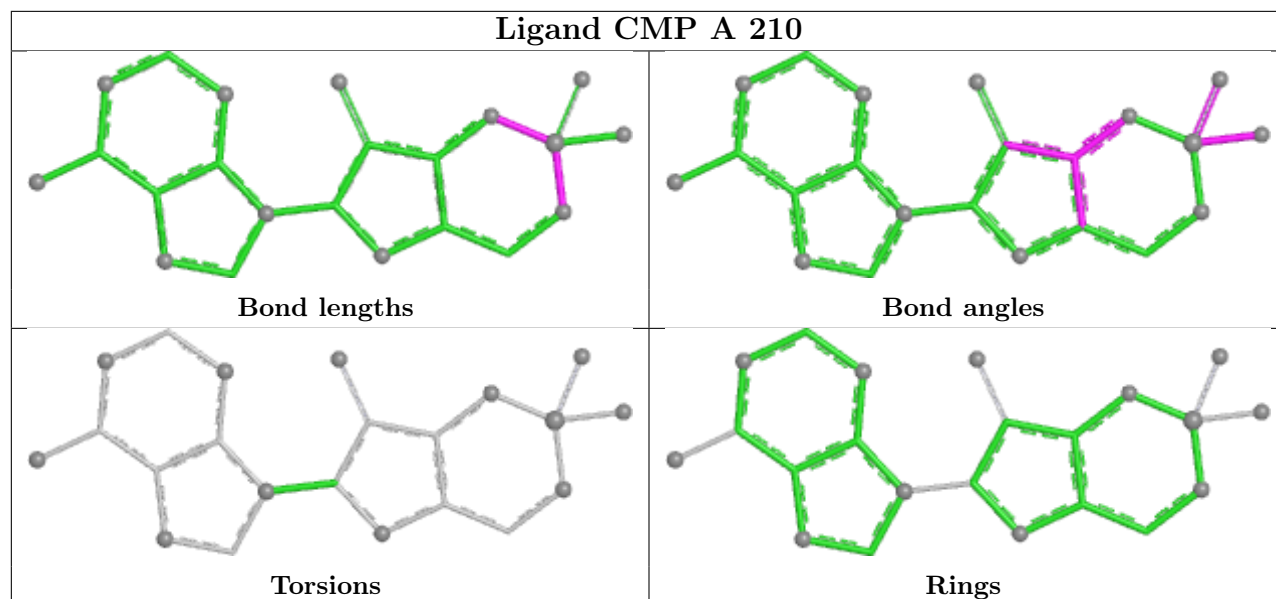
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	210	CMP	3	0
4	A	210	CMP	4	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	W	17/17 (100%)	0.50	0	100 100	62, 107, 140, 142	0
1	Y	17/17 (100%)	0.35	0	100 100	71, 107, 145, 147	0
2	X	21/21 (100%)	0.43	0	100 100	59, 88, 150, 151	0
2	Z	21/21 (100%)	0.40	1 (4%)	35 28	61, 95, 140, 145	0
3	A	200/209 (95%)	-0.32	1 (0%)	87 82	36, 50, 73, 92	0
3	B	200/209 (95%)	-0.04	1 (0%)	87 82	38, 65, 104, 130	0
All	All	476/494 (96%)	-0.08	3 (0%)	85 80	36, 60, 122, 151	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	207	GLY	2.1
3	A	207	GLY	2.1
2	Z	-8	DT	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.

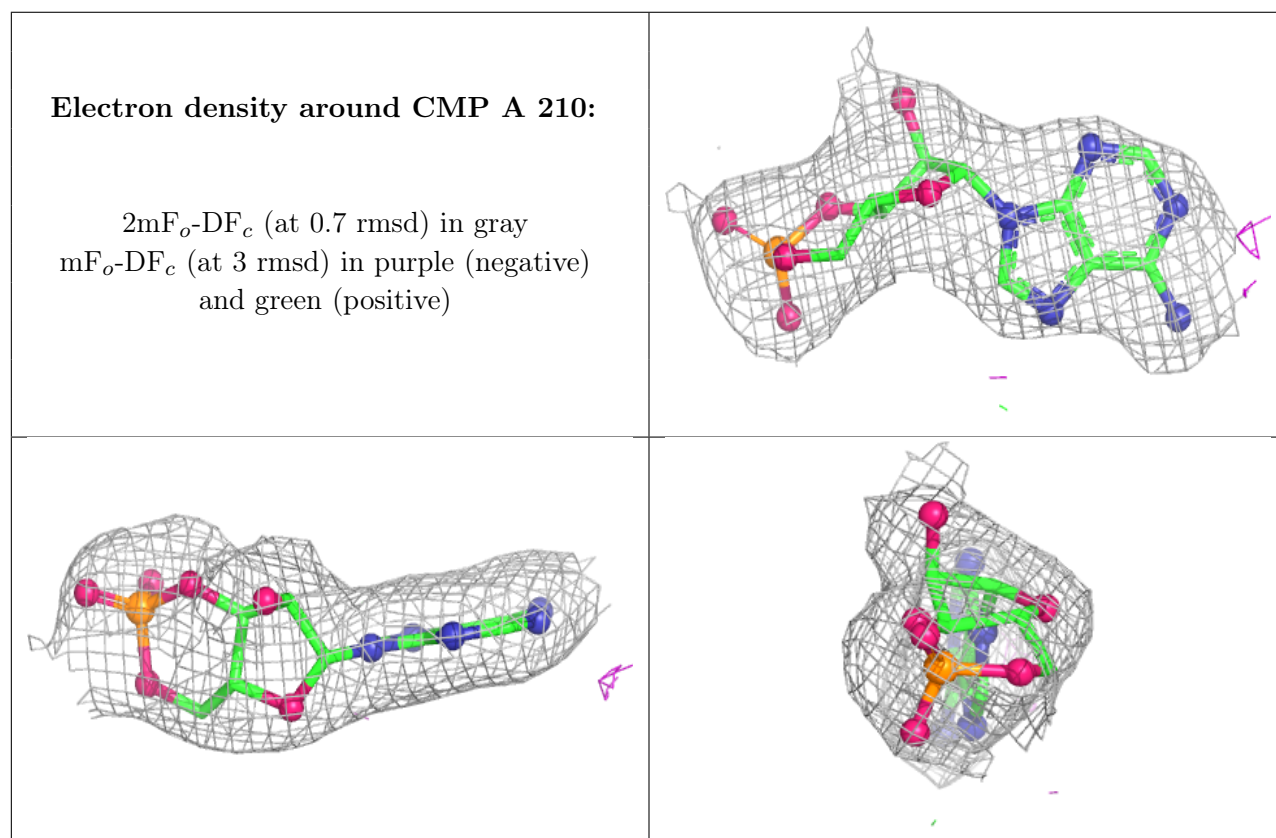
6.4 Ligands [i](#)

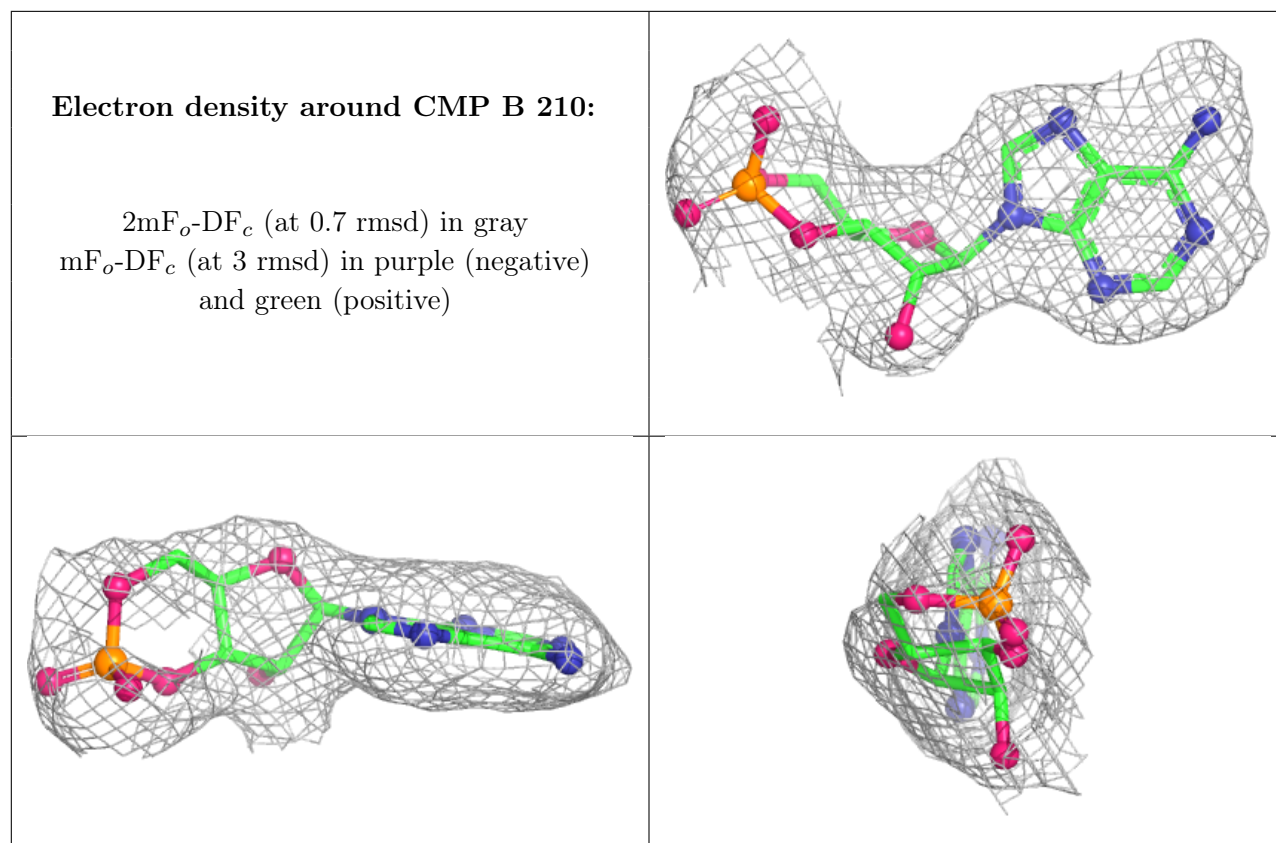
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

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
4	CMP	A	210	22/22	0.97	0.07	41,44,48,51	0
4	CMP	B	210	22/22	0.97	0.06	36,42,43,45	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.