



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:14 AM UTC

PDB ID : 1RUO / pdb_00001ruo
Title : CATABOLITE GENE ACTIVATOR PROTEIN (CAP) MUTANT/DNA
COMPLEX + ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE
Authors : Parkinson, G.N.; Ebright, R.H.; Berman, H.M.
Deposited on : 1996-05-26
Resolution : 2.70 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

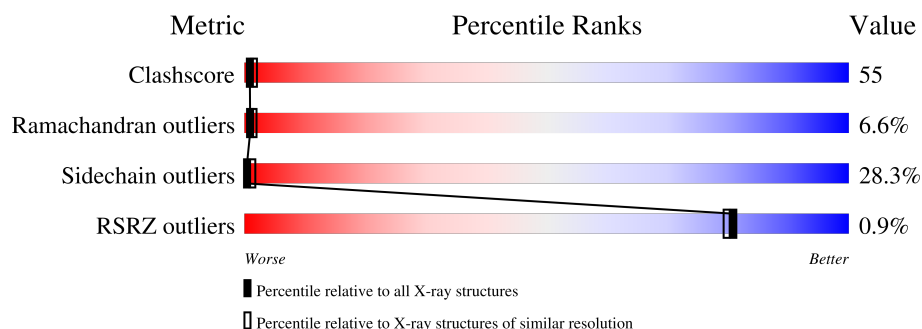
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	C	14	14% 86%
1	E	14	7% 64% 36%
2	D	17	76% 24%
2	F	17	71% 29%
3	A	209	25% 44% 21% 5% 5%
3	B	209	% 28% 44% 21% • •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	CMP	A	210	-	-	X	-
4	CMP	B	210	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4511 atoms, of which 13 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*AP*TP*GP*TP*GP*AP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	14	Total	C	N	O	P	0	0	0
			290	139	59	79	13			
1	E	14	Total	C	N	O	P	0	0	0
			290	139	59	79	13			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	17	Total	C	N	O	P	0	0	0
			341	166	56	103	16			
2	F	17	Total	C	N	O	P	0	0	0
			341	166	56	103	16			

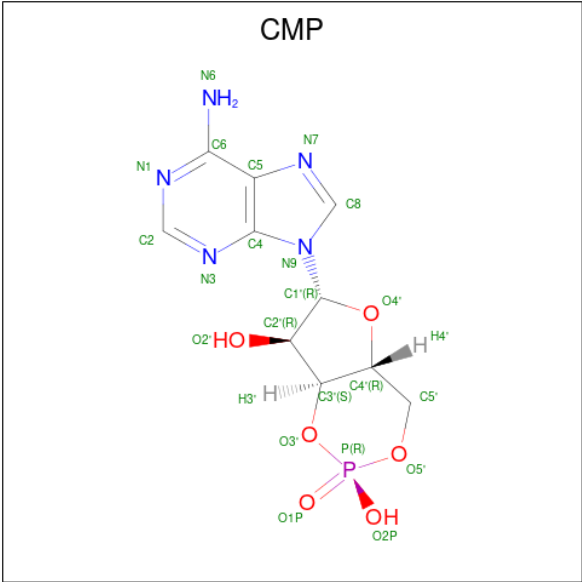
- Molecule 3 is a protein called PROTEIN (CATABOLITE GENE ACTIVATOR PROTEIN (CAP)).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	A	198	Total	C	H	N	O	S	0	0	0
			1574	999	4	274	288	9			
3	B	201	Total	C	H	N	O	S	0	0	0
			1596	1008	9	277	293	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	PHE	GLU	engineered mutation	UNP P0ACJ8
B	181	PHE	GLU	engineered mutation	UNP P0ACJ8

- 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		

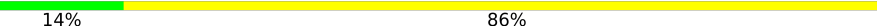
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	O	0	0
			1	1		
5	D	10	Total	O	0	0
			10	10		
5	E	2	Total	O	0	0
			2	2		
5	F	3	Total	O	0	0
			3	3		
5	A	13	Total	O	0	0
			13	13		
5	B	6	Total	O	0	0
			6	6		

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

Chain C: 



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

Chain E: 



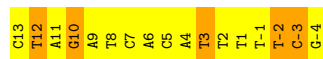
- Molecule 2: DNA (5'-D(*CP*TP*AP*GP*AP*TP*CP*AP*CP*AP*TP*TP*TP*TP*TP*CP*G)-3')

Chain D: 

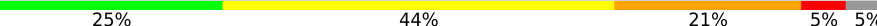


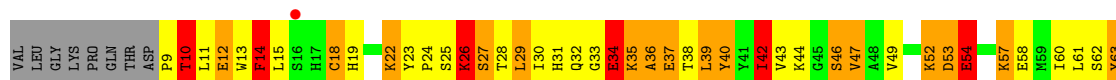
- Molecule 2: DNA (5'-D(*CP*TP*AP*GP*AP*TP*CP*AP*CP*AP*TP*TP*TP*TP*TP*CP*G)-3')

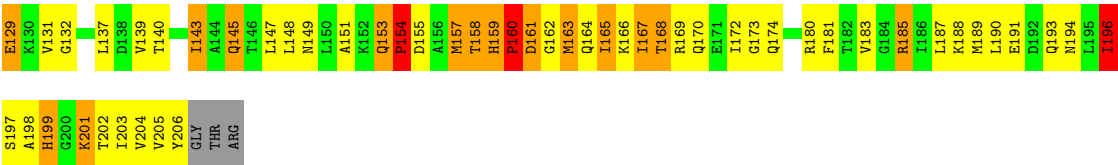
Chain F: 



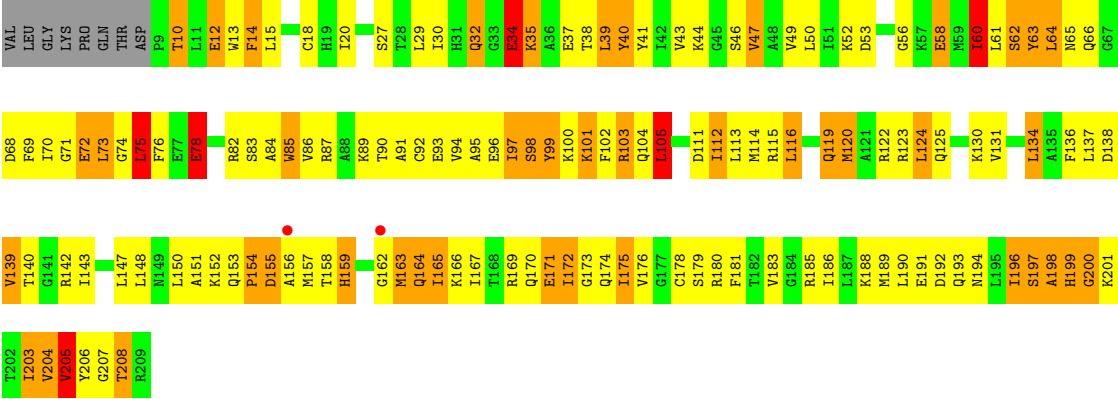
- Molecule 3: PROTEIN (CATABOLITE GENE ACTIVATOR PROTEIN (CAP))

Chain A: 





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



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	137.29Å 153.10Å 76.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.70 10.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	78.9 (10.00-2.70) 79.4 (10.00-2.70)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.91 (at 2.72Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.213 , (Not available) 0.218 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.482	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.21 , 225.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4511	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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	C	0.77	0/327	1.12	0/504
1	E	0.86	0/327	1.41	7/504 (1.4%)
2	D	0.91	0/380	1.32	5/584 (0.9%)
2	F	1.00	0/380	1.49	6/584 (1.0%)
3	A	1.10	3/1596 (0.2%)	1.47	26/2149 (1.2%)
3	B	1.09	3/1613 (0.2%)	1.47	28/2171 (1.3%)
All	All	1.04	6/4623 (0.1%)	1.43	72/6496 (1.1%)

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
3	B	0	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	163	MET	SD-CE	8.27	2.00	1.79
3	A	172	ILE	CA-CB	7.32	1.63	1.54
3	B	205	VAL	CA-CB	6.54	1.63	1.54
3	A	28	THR	CA-CB	5.86	1.61	1.52
3	B	172	ILE	CA-CB	-5.64	1.48	1.54
3	B	91	ALA	CA-CB	-5.19	1.45	1.53

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	113	LEU	N-CA-C	-9.90	100.57	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	83	SER	N-CA-C	9.39	123.24	111.69
3	A	63	TYR	N-CA-C	-9.06	95.48	109.52
3	A	57	LYS	N-CA-C	8.64	122.76	110.23
3	A	75	LEU	N-CA-C	8.59	129.09	110.80
3	B	198	ALA	N-CA-C	8.20	120.05	108.74
1	E	19	DT	O5'-C5'-C4'	8.06	122.89	110.80
3	B	20	ILE	N-CA-C	7.98	118.71	107.37
3	B	10	THR	N-CA-C	-7.94	101.59	111.11
3	B	205	VAL	CB-CA-C	-7.92	98.31	111.29
3	B	96	GLU	N-CA-C	7.88	121.22	108.76
2	D	13	DG	P-O3'-C3'	7.51	131.47	120.20
3	B	154	PRO	N-CA-C	7.31	121.84	110.80
3	A	76	PHE	N-CA-C	7.15	122.10	112.88
3	A	154	PRO	N-CA-C	7.14	123.28	113.65
2	F	10	DG	N9-C1'-C2'	6.84	123.76	113.50
3	B	105	LEU	N-CA-C	-6.82	103.07	111.33
2	F	3	DT	N1-C1'-C2'	6.81	123.71	113.50
3	A	108	VAL	N-CA-C	-6.79	105.00	112.80
3	A	14	PHE	N-CA-C	-6.74	103.93	111.28
3	B	78	GLU	N-CA-C	6.65	120.50	109.46
1	E	27	DG	C1'-O4'-C4'	-6.44	100.03	109.70
2	D	16	DC	N1-C1'-C2'	-6.37	103.94	113.50
2	D	14	DA	P-O5'-C5'	-6.35	110.48	120.00
3	B	131	VAL	CB-CA-C	-6.19	104.04	111.97
3	B	134	LEU	N-CA-C	-6.17	106.30	113.88
3	B	89	LYS	N-CA-C	-6.16	104.47	112.23
3	B	60	ILE	CB-CA-C	-6.11	103.56	110.96
3	B	32	GLN	N-CA-C	-6.08	101.41	110.23
3	A	89	LYS	N-CA-C	-6.08	106.02	113.50
2	F	-2	DT	C6-N1-C1'	-6.05	110.27	119.35
2	D	17	DA	N9-C1'-C2'	-6.04	104.44	113.50
2	F	-2	DT	C2-N1-C1'	6.00	128.35	119.35
3	A	26	LYS	CA-C-N	-5.97	112.24	120.71
3	A	26	LYS	C-N-CA	-5.97	112.24	120.71
3	B	200	GLY	N-CA-C	-5.95	99.09	113.18
3	B	198	ALA	CA-C-N	5.92	132.84	121.54
3	B	198	ALA	C-N-CA	5.92	132.84	121.54
3	A	172	ILE	N-CA-C	-5.91	104.97	110.53
3	A	35	LYS	N-CA-C	5.90	118.02	108.34
3	B	175	ILE	CB-CA-C	-5.89	104.52	111.94
3	A	72	GLU	N-CA-C	-5.87	106.67	113.88
2	F	-3	DC	C2'-C3'-O3'	5.85	120.28	111.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	99	TYR	N-CA-C	-5.84	104.83	111.14
3	A	94	VAL	N-CA-C	5.79	116.49	108.84
3	B	75	LEU	N-CA-C	5.77	123.09	110.80
3	B	98	SER	N-CA-C	-5.77	102.76	110.55
3	B	47	VAL	N-CA-CB	-5.69	101.85	111.23
3	B	185	ARG	N-CA-C	5.64	117.43	111.28
3	A	159	HIS	CA-C-N	5.60	126.84	119.84
3	A	159	HIS	C-N-CA	5.60	126.84	119.84
3	A	19	HIS	N-CA-C	5.56	118.18	107.44
3	A	10	THR	N-CA-C	5.55	117.77	111.11
3	B	63	TYR	N-CA-C	-5.51	101.15	109.85
1	E	25	DG	C4'-C3'-O3'	5.50	118.25	110.00
3	A	172	ILE	N-CA-CB	5.45	117.55	110.57
3	A	196	ILE	N-CA-C	5.44	115.39	107.88
3	B	119	GLN	N-CA-C	-5.44	105.04	110.97
3	B	172	ILE	CB-CA-C	-5.39	104.98	111.88
3	A	77	GLU	N-CA-C	-5.37	99.37	110.80
1	E	18	DG	N9-C1'-C2'	5.35	121.53	113.50
1	E	19	DT	C5'-C4'-O4'	5.30	117.35	109.40
3	A	161	ASP	N-CA-C	-5.27	100.10	109.06
3	A	149	ASN	N-CA-C	5.18	116.61	111.07
1	E	14	DT	C2'-C3'-O3'	-5.17	103.74	111.50
3	A	64	LEU	N-CA-C	-5.17	99.18	108.48
3	B	172	ILE	N-CA-C	-5.16	104.94	110.36
3	B	163	MET	N-CA-C	5.15	117.33	110.53
2	F	12	DT	C4'-C3'-O3'	5.14	117.71	110.00
2	D	13	DG	O3'-P-O5'	5.07	111.60	104.00
3	B	85	TRP	N-CA-C	-5.04	102.54	110.36
1	E	25	DG	N9-C1'-C2'	-5.02	105.97	113.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	B	159	HIS	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	C	290	0	159	52	0
1	E	290	0	159	75	0
2	D	341	0	196	51	0
2	F	341	0	195	55	0
3	A	1570	4	1612	137	0
3	B	1587	9	1624	140	0
4	A	22	0	11	9	0
4	B	22	0	11	11	0
5	A	13	0	0	1	0
5	B	6	0	0	4	0
5	C	1	0	0	0	0
5	D	10	0	0	0	0
5	E	2	0	0	0	0
5	F	3	0	0	0	0
All	All	4498	13	3967	460	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 55.

All (460) 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:B:210:CMP:C2	4:B:210:CMP:H2	0.97	1.48
4:A:210:CMP:H2	4:A:210:CMP:C2	0.97	1.46
1:E:16:DG:H1'	1:E:15:DA:C8	1.82	1.14
2:F:4:DA:C8	2:F:3:DT:H72	1.82	1.13
2:D:15:DT:H72	3:B:181:PHE:CD1	1.86	1.10
1:E:17:DT:O3'	1:E:16:DG:H2'	1.54	1.07
1:E:17:DT:H2''	1:E:16:DG:H5'	1.09	1.06
1:E:20:DA:C8	1:E:19:DT:H72	1.93	1.03
1:E:27:DG:H4'	1:E:26:DC:H5'	1.41	1.02
1:E:17:DT:C2'	1:E:16:DG:H5'	1.88	1.02
1:E:14:DT:H2'	2:F:13:DC:H5	1.26	1.01
1:E:17:DT:H2''	1:E:16:DG:C5'	1.93	0.99
2:F:10:DG:OP1	3:A:139:VAL:HG12	1.64	0.97
3:B:178:CYS:SG	3:B:183:VAL:HG23	2.05	0.96
3:A:29:LEU:HD11	3:A:94:VAL:HG21	1.48	0.95
1:C:-4:DC:H2''	1:C:-3:DG:C8	2.01	0.95
3:B:163:MET:HB2	3:B:204:VAL:HG12	1.45	0.95
2:D:17:DA:H2''	2:D:18:DC:OP2	1.64	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:15:DT:C7	3:B:181:PHE:CD1	2.54	0.91
1:E:14:DT:H2'	2:F:13:DC:C5	2.06	0.90
1:E:16:DG:H1'	1:E:15:DA:N7	1.85	0.90
1:E:24:DA:C2	1:E:23:DA:C6	2.60	0.89
1:E:19:DT:H2''	1:E:18:DG:O5'	1.73	0.88
1:E:27:DG:C4	1:E:26:DC:C5	2.62	0.87
1:E:26:DC:H2'	1:E:25:DG:C8	2.12	0.85
3:B:169:ARG:HA	3:B:172:ILE:HD12	1.57	0.84
2:F:3:DT:H1'	2:F:2:DT:H5'	1.60	0.84
2:F:6:DA:H2''	2:F:5:DC:O5'	1.76	0.83
1:C:-3:DG:H2''	1:C:-2:DA:N7	1.94	0.83
2:D:19:DA:C2'	2:D:20:DT:H71	2.09	0.82
2:F:10:DG:P	3:A:139:VAL:HG12	2.19	0.82
3:A:71:GLY:HA2	4:A:210:CMP:O3'	1.79	0.82
1:C:7:DG:H5'	1:C:7:DG:C8	2.15	0.81
1:C:-1:DA:H2''	1:C:1:DA:OP2	1.80	0.81
1:C:7:DG:H2''	1:C:8:DA:C8	2.15	0.81
4:B:210:CMP:H3'	4:B:210:CMP:H8	1.63	0.80
2:F:12:DT:H4'	2:F:11:DA:OP1	1.80	0.80
1:C:7:DG:H1'	1:C:8:DA:C5	2.18	0.78
1:E:16:DG:OP2	1:E:16:DG:H3'	1.85	0.77
1:E:24:DA:C2	1:E:23:DA:C5	2.71	0.77
1:C:6:DT:C2'	1:C:7:DG:H5''	2.16	0.75
3:A:69:PHE:H	3:A:123:ARG:HH22	1.32	0.75
3:A:23:TYR:CD2	3:A:29:LEU:HD22	2.21	0.75
2:D:21:DT:C2	1:E:20:DA:H2	2.04	0.75
3:A:9:PRO:O	3:A:12:GLU:HG2	1.86	0.75
3:A:38:THR:HG23	3:A:97:ILE:O	1.86	0.74
3:B:188:LYS:O	3:B:191:GLU:HG3	1.87	0.74
3:B:49:VAL:HG22	3:B:86:VAL:HG22	1.70	0.74
3:B:150:LEU:HB3	3:B:165:ILE:HD13	1.69	0.74
2:F:6:DA:H2'	2:F:5:DC:C6	2.23	0.74
3:A:124:LEU:HD21	4:B:210:CMP:N6	2.03	0.74
3:A:204:VAL:HG12	3:A:206:TYR:HB2	1.68	0.74
1:E:23:DA:C2	1:E:22:DA:C4	2.75	0.73
3:A:113:LEU:HD11	3:B:114:MET:SD	2.28	0.73
1:C:4:DT:OP1	3:A:169:ARG:HD3	1.88	0.73
2:D:15:DT:H72	3:B:181:PHE:CG	2.24	0.73
2:D:10:DC:O2	2:F:9:DA:H2	1.72	0.72
2:F:-4:DG:OP2	2:F:-4:DG:H3'	1.89	0.72
2:D:20:DT:H2'	2:D:21:DT:H6	1.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:32:GLN:HG2	3:A:33:GLY:H	1.54	0.72
2:D:20:DT:H2''	2:D:21:DT:O4'	1.90	0.72
3:B:151:ALA:HB1	3:B:163:MET:HG3	1.71	0.72
1:E:23:DA:C2	1:E:22:DA:N3	2.58	0.72
3:A:83:SER:HB2	4:A:210:CMP:H5'1	1.71	0.71
2:F:13:DC:H1'	2:F:12:DT:O5'	1.89	0.71
3:A:145:GLN:O	3:A:148:LEU:HB3	1.90	0.71
3:A:201:LYS:NZ	3:A:203:ILE:HD11	2.06	0.71
3:A:73:LEU:HD21	3:B:124:LEU:HB3	1.73	0.70
3:B:134:LEU:HD23	3:B:142:ARG:HH11	1.55	0.70
1:C:7:DG:H2''	1:C:8:DA:N7	2.06	0.70
2:D:19:DA:H2''	2:D:20:DT:OP2	1.92	0.69
3:A:32:GLN:HG2	3:A:33:GLY:N	2.06	0.69
1:C:6:DT:H2'	1:C:7:DG:H5''	1.75	0.69
2:D:21:DT:H2'	2:D:22:DT:C6	2.28	0.69
2:D:24:DT:O2	1:E:23:DA:C2	2.46	0.69
3:A:73:LEU:HD13	4:A:210:CMP:H3'	1.74	0.69
1:C:7:DG:C4	1:C:8:DA:C6	2.80	0.69
1:E:24:DA:N3	1:E:23:DA:C5	2.61	0.68
2:F:8:DT:H72	3:A:181:PHE:CD2	2.27	0.68
3:A:34:GLU:O	3:A:81:GLU:HA	1.92	0.68
3:B:63:TYR:O	3:B:64:LEU:HD13	1.92	0.68
1:E:16:DG:H1'	1:E:15:DA:C5	2.27	0.68
2:D:24:DT:H2''	2:D:25:DC:H5'	1.74	0.68
3:B:30:ILE:HG23	3:B:82:ARG:HG3	1.76	0.68
3:A:18:CYS:SG	3:A:97:ILE:HG23	2.34	0.68
3:B:61:LEU:O	3:B:61:LEU:HD23	1.94	0.67
2:F:4:DA:N9	2:F:3:DT:H72	2.08	0.67
3:A:35:LYS:O	3:A:37:GLU:HG3	1.94	0.67
1:E:23:DA:N1	1:E:22:DA:C2	2.62	0.67
1:C:6:DT:C3'	1:C:7:DG:H5''	2.24	0.67
1:E:27:DG:H4'	1:E:26:DC:C5'	2.20	0.67
3:A:46:SER:HA	3:A:65:ASN:HA	1.77	0.67
3:B:100:LYS:H	3:B:100:LYS:HD3	1.60	0.67
3:A:111:ASP:O	3:A:114:MET:HB2	1.93	0.66
3:A:109:ASN:HD22	3:A:112:ILE:HG12	1.59	0.66
1:C:7:DG:C2'	1:C:8:DA:N7	2.59	0.66
3:A:29:LEU:HD23	3:A:88:ALA:HB2	1.78	0.66
3:B:32:GLN:HG3	3:B:85:TRP:CE3	2.30	0.66
3:B:72:GLU:HB3	3:B:116:LEU:HD21	1.76	0.66
3:A:124:LEU:CD2	4:B:210:CMP:N6	2.59	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:98:SER:OG	3:A:100:LYS:HG3	1.96	0.65
3:A:162:GLY:HA3	3:A:204:VAL:O	1.95	0.65
3:B:18:CYS:HB3	3:B:95:ALA:HB1	1.77	0.65
3:B:70:ILE:HG22	3:B:71:GLY:H	1.60	0.65
3:A:23:TYR:HD2	3:A:29:LEU:HD22	1.62	0.65
3:A:161:ASP:OD2	3:A:205:VAL:HG22	1.96	0.65
1:C:-1:DA:C2	1:C:1:DA:C4	2.84	0.65
1:C:5:DG:H2''	1:C:6:DT:C6	2.32	0.64
2:F:4:DA:C6	2:F:3:DT:O4	2.49	0.64
3:A:75:LEU:HD21	3:A:102:PHE:CD2	2.32	0.64
3:B:115:ARG:O	3:B:119:GLN:HG2	1.98	0.64
1:E:25:DG:H4'	1:E:24:DA:OP1	1.97	0.64
2:D:20:DT:H2'	2:D:21:DT:C6	2.32	0.64
3:A:36:ALA:HA	3:A:82:ARG:HH11	1.62	0.64
3:B:150:LEU:O	3:B:165:ILE:HG21	1.98	0.64
2:D:23:DT:C2	1:E:22:DA:H2	2.16	0.64
3:B:178:CYS:SG	3:B:183:VAL:CG2	2.84	0.64
1:E:20:DA:N7	1:E:19:DT:H72	2.13	0.63
3:B:120:MET:HA	3:B:123:ARG:HB2	1.80	0.63
1:E:26:DC:C2'	1:E:25:DG:C8	2.81	0.63
3:A:69:PHE:H	3:A:123:ARG:NH2	1.96	0.63
1:C:-3:DG:C2	1:C:-2:DA:C2	2.87	0.63
2:D:21:DT:O2	1:E:20:DA:H2	1.80	0.63
1:E:26:DC:H3'	1:E:25:DG:C8	2.33	0.63
1:C:7:DG:C2	1:C:8:DA:C2	2.86	0.63
2:D:16:DC:H3'	2:D:16:DC:C6	2.35	0.62
3:A:122:ARG:HH22	3:B:76:PHE:HB2	1.64	0.62
3:A:18:CYS:SG	3:A:96:GLU:C	2.82	0.62
3:B:83:SER:OG	4:B:210:CMP:O2P	2.15	0.62
3:B:87:ARG:HG2	3:B:87:ARG:HH11	1.65	0.62
3:B:18:CYS:SG	3:B:97:ILE:CG1	2.88	0.62
3:B:170:GLN:CG	5:B:212:HOH:O	2.47	0.62
3:A:196:ILE:HG12	3:A:197:SER:N	2.13	0.62
2:F:7:DC:H2''	2:F:6:DA:OP2	1.98	0.61
2:F:-2:DT:H2''	2:F:-3:DC:C6	2.36	0.61
3:A:153:GLN:HG3	3:A:154:PRO:N	2.15	0.61
1:E:26:DC:C3'	1:E:25:DG:C8	2.84	0.61
2:F:7:DC:H2'	2:F:7:DC:O5'	2.01	0.61
2:F:6:DA:C2'	2:F:5:DC:O5'	2.48	0.61
3:A:155:ASP:O	3:A:157:MET:HE3	2.00	0.61
2:F:4:DA:C4	2:F:3:DT:C5	2.88	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:170:GLN:HG2	5:B:212:HOH:O	2.01	0.61
1:E:27:DG:C4'	1:E:26:DC:H5'	2.26	0.61
1:C:8:DA:C4	1:C:9:DT:N3	2.69	0.60
3:B:61:LEU:O	3:B:62:SER:HB2	2.01	0.60
2:D:23:DT:N3	1:E:22:DA:C2	2.70	0.60
3:B:52:LYS:HD3	3:B:85:TRP:CH2	2.37	0.60
3:A:47:VAL:HG13	3:A:86:VAL:HG13	1.84	0.60
3:B:111:ASP:O	3:B:114:MET:HB2	2.01	0.60
3:B:204:VAL:O	3:B:205:VAL:HB	2.00	0.60
2:D:11:DT:H2''	2:D:12:DA:C8	2.36	0.59
1:C:3:DA:H1'	1:C:4:DT:H5'	1.85	0.59
3:B:18:CYS:SG	3:B:97:ILE:HD11	2.42	0.59
2:D:18:DC:H2''	2:D:19:DA:OP2	2.00	0.59
3:A:188:LYS:O	3:A:191:GLU:HB3	2.02	0.59
3:B:157:MET:HB3	3:B:164:GLN:NE2	2.18	0.59
3:B:186:ILE:O	3:B:190:LEU:HB2	2.03	0.59
3:A:73:LEU:CD1	4:A:210:CMP:H3'	2.33	0.59
2:D:13:DG:OP1	3:B:138:ASP:HB2	2.01	0.59
2:D:24:DT:H2'	2:D:25:DC:O4'	2.02	0.59
2:D:18:DC:H2''	2:D:19:DA:C8	2.37	0.58
1:E:15:DA:C8	1:E:14:DT:H73	2.38	0.58
3:B:40:TYR:HB2	3:B:70:ILE:HB	1.84	0.58
3:B:170:GLN:HA	5:B:212:HOH:O	2.03	0.58
1:E:24:DA:C6	1:E:23:DA:N6	2.72	0.58
3:A:38:THR:HG21	3:A:96:GLU:HB3	1.84	0.58
3:B:32:GLN:HE21	3:B:85:TRP:HZ3	1.51	0.58
1:C:8:DA:C4	1:C:9:DT:C4	2.91	0.57
1:C:6:DT:H71	3:A:181:PHE:CD1	2.40	0.57
1:C:7:DG:C1'	1:C:8:DA:C5	2.88	0.57
3:A:109:ASN:ND2	3:A:112:ILE:HG12	2.19	0.57
3:A:25:SER:O	3:A:26:LYS:HB2	2.02	0.57
1:C:7:DG:H1'	1:C:8:DA:C8	2.38	0.57
3:A:18:CYS:SG	3:A:97:ILE:N	2.78	0.57
1:C:-3:DG:N2	1:C:-2:DA:C2	2.73	0.56
2:F:6:DA:C8	2:F:6:DA:O5'	2.58	0.56
3:A:44:LYS:HB3	3:A:93:GLU:HB2	1.87	0.56
3:B:105:LEU:HB3	3:B:112:ILE:HD11	1.87	0.56
1:E:21:DA:N6	1:E:20:DA:N1	2.54	0.56
3:A:43:VAL:HB	3:A:93:GLU:O	2.05	0.56
1:E:19:DT:H2''	1:E:18:DG:C5'	2.35	0.56
3:A:72:GLU:HG3	3:A:120:MET:HE2	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:58:GLU:O	5:B:213:HOH:O	2.17	0.56
1:C:7:DG:N3	1:C:8:DA:C2	2.74	0.56
3:A:69:PHE:HD2	3:A:119:GLN:OE1	1.88	0.56
1:E:14:DT:C2'	2:F:13:DC:C5	2.84	0.56
3:A:73:LEU:CD1	4:A:210:CMP:H2'	2.35	0.56
4:B:210:CMP:H3'	4:B:210:CMP:C8	2.35	0.56
1:C:7:DG:H1'	1:C:8:DA:C4	2.40	0.56
3:A:113:LEU:CD1	3:B:114:MET:SD	2.93	0.56
3:A:167:ILE:HG12	3:A:168:THR:H	1.71	0.56
1:C:5:DG:C8	1:C:6:DT:H72	2.42	0.55
3:B:72:GLU:OE2	3:B:73:LEU:HD13	2.06	0.55
1:C:6:DT:C3'	1:C:7:DG:C5'	2.84	0.55
3:A:124:LEU:HD21	4:B:210:CMP:C6	2.41	0.55
3:A:173:GLY:N	3:A:183:VAL:HG21	2.21	0.55
3:A:98:SER:HB3	3:A:101:LYS:HB3	1.88	0.55
3:B:159:HIS:ND1	3:B:203:ILE:HG23	2.22	0.55
3:A:90:THR:O	3:A:92:CYS:N	2.39	0.55
2:D:15:DT:H71	3:B:181:PHE:HD1	1.72	0.55
1:E:19:DT:C2'	1:E:18:DG:O5'	2.52	0.55
2:D:19:DA:H2''	2:D:20:DT:H71	1.87	0.55
3:A:40:TYR:HB2	3:A:70:ILE:HB	1.88	0.55
3:B:154:PRO:O	3:B:155:ASP:HB2	2.06	0.55
2:D:11:DT:H2''	2:D:12:DA:OP2	2.06	0.54
2:F:4:DA:C6	2:F:3:DT:C4	2.94	0.54
3:A:24:PRO:HD2	3:A:27:SER:OG	2.07	0.54
3:B:50:LEU:O	3:B:84:ALA:HB1	2.07	0.54
2:D:18:DC:C4	2:D:19:DA:N6	2.75	0.54
3:A:71:GLY:HA2	4:A:210:CMP:P	2.47	0.54
3:B:52:LYS:HB2	3:B:85:TRP:CZ3	2.41	0.54
2:D:21:DT:C2	1:E:20:DA:C2	2.92	0.54
2:F:4:DA:C5	2:F:3:DT:C4	2.95	0.54
3:A:39:LEU:HD23	3:A:71:GLY:O	2.08	0.54
3:B:99:TYR:O	3:B:103:ARG:HG3	2.08	0.54
3:B:52:LYS:HD3	3:B:85:TRP:CZ3	2.42	0.54
3:B:159:HIS:N	3:B:162:GLY:O	2.40	0.54
2:D:24:DT:C2'	2:D:25:DC:O4'	2.56	0.54
1:E:15:DA:C8	1:E:14:DT:C7	2.90	0.54
3:B:43:VAL:HB	3:B:93:GLU:HB3	1.90	0.54
3:B:90:THR:O	3:B:92:CYS:SG	2.63	0.53
2:D:15:DT:C7	3:B:181:PHE:HD1	2.15	0.53
1:E:27:DG:H1'	1:E:26:DC:C6	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:16:DC:H3'	2:D:16:DC:H6	1.71	0.53
2:D:23:DT:N3	1:E:22:DA:H2	2.06	0.53
1:E:27:DG:N3	1:E:26:DC:C5	2.76	0.53
3:A:198:ALA:O	3:A:199:HIS:HB2	2.08	0.53
3:B:140:THR:HG22	3:B:186:ILE:HG23	1.91	0.53
3:A:54:GLU:CD	3:A:54:GLU:H	2.17	0.53
3:A:53:ASP:HB3	3:A:57:LYS:O	2.09	0.53
2:D:17:DA:O5'	2:D:17:DA:C8	2.61	0.53
2:F:7:DC:C5	2:F:6:DA:N6	2.77	0.53
2:F:-1:DT:H2''	2:F:-2:DT:OP2	2.08	0.53
1:C:6:DT:H3'	1:C:7:DG:C5'	2.39	0.53
2:D:20:DT:C2	2:D:21:DT:C6	2.96	0.53
3:A:10:THR:HA	3:A:13:TRP:HB3	1.91	0.53
1:E:23:DA:C2	1:E:22:DA:C2	2.96	0.52
3:B:164:GLN:HA	3:B:203:ILE:HA	1.91	0.52
3:B:169:ARG:HA	3:B:172:ILE:CD1	2.33	0.52
3:A:53:ASP:HB2	3:B:136:PHE:CE1	2.45	0.52
3:A:201:LYS:HZ2	3:A:203:ILE:HD11	1.74	0.52
3:B:137:LEU:HD12	3:B:137:LEU:H	1.73	0.52
1:E:19:DT:H73	3:B:180:ARG:NH1	2.25	0.52
1:C:8:DA:C5	1:C:9:DT:C4	2.97	0.52
1:C:7:DG:H1'	1:C:8:DA:N7	2.24	0.52
3:A:158:THR:HA	3:A:163:MET:SD	2.50	0.52
3:A:75:LEU:HD23	3:A:116:LEU:HD21	1.91	0.52
1:E:24:DA:N3	1:E:23:DA:N7	2.58	0.51
2:D:26:DG:O6	1:E:26:DC:N3	2.43	0.51
3:B:159:HIS:CG	3:B:159:HIS:O	2.62	0.51
2:D:16:DC:C6	2:D:16:DC:C3'	2.93	0.51
3:A:60:ILE:HD13	3:A:63:TYR:HE2	1.75	0.51
3:A:109:ASN:HD21	3:A:112:ILE:H	1.58	0.51
3:B:205:VAL:HG12	3:B:206:TYR:N	2.24	0.51
3:B:65:ASN:O	3:B:68:ASP:HB2	2.11	0.51
3:A:60:ILE:HD13	3:A:63:TYR:CE2	2.44	0.51
3:B:192:ASP:C	3:B:194:ASN:H	2.19	0.51
2:F:3:DT:H1'	2:F:2:DT:C5'	2.37	0.51
3:A:22:LYS:NZ	3:A:91:ALA:HB1	2.26	0.51
1:E:24:DA:N1	1:E:23:DA:C6	2.80	0.50
3:A:191:GLU:HA	3:A:196:ILE:O	2.11	0.50
3:B:97:ILE:HG22	3:B:98:SER:O	2.11	0.50
3:B:163:MET:O	3:B:203:ILE:HA	2.11	0.50
2:D:20:DT:C2'	2:D:21:DT:H6	2.22	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:201:LYS:CE	3:A:203:ILE:HD11	2.42	0.50
3:B:75:LEU:HD21	3:B:102:PHE:CD2	2.47	0.50
1:E:21:DA:C6	1:E:20:DA:C2	3.00	0.50
3:A:75:LEU:HD21	3:A:102:PHE:CE2	2.46	0.50
3:B:18:CYS:SG	3:B:97:ILE:HG13	2.52	0.50
3:B:137:LEU:HD12	3:B:137:LEU:N	2.27	0.50
1:E:15:DA:OP2	1:E:15:DA:H8	1.94	0.50
3:A:108:VAL:HG12	3:A:108:VAL:O	2.12	0.50
3:A:109:ASN:OD1	3:A:111:ASP:HB2	2.12	0.50
3:B:53:ASP:O	3:B:56:GLY:N	2.45	0.50
1:E:21:DA:N6	1:E:20:DA:C6	2.79	0.49
1:E:16:DG:O3'	1:E:15:DA:C8	2.64	0.49
3:B:60:ILE:HD11	3:B:171:GLU:OE1	2.13	0.49
1:C:3:DA:H2''	1:C:4:DT:O5'	2.11	0.49
3:A:71:GLY:HA2	4:A:210:CMP:O1P	2.12	0.49
1:C:7:DG:C1'	1:C:8:DA:N7	2.75	0.49
3:A:185:ARG:HD2	5:A:212:HOH:O	2.12	0.49
3:B:41:TYR:HD2	3:B:69:PHE:CD1	2.31	0.49
3:B:205:VAL:HG12	3:B:206:TYR:CD2	2.48	0.49
2:D:19:DA:H2'	2:D:20:DT:H71	1.91	0.49
3:A:118:ALA:O	3:A:122:ARG:HG2	2.12	0.49
2:F:7:DC:N4	3:A:181:PHE:CE2	2.76	0.49
1:C:8:DA:N6	2:F:9:DA:N6	2.61	0.49
3:A:131:VAL:HG11	3:B:130:LYS:HD3	1.93	0.49
3:A:173:GLY:CA	3:A:183:VAL:HG21	2.43	0.49
3:B:156:ALA:HB1	3:B:163:MET:SD	2.52	0.49
1:C:7:DG:C2'	1:C:8:DA:C8	2.92	0.48
3:A:30:ILE:HG22	3:A:31:HIS:N	2.27	0.48
3:A:39:LEU:HB2	3:A:99:TYR:CE1	2.48	0.48
3:B:37:GLU:OE2	3:B:100:LYS:HD2	2.12	0.48
3:B:140:THR:HG22	3:B:186:ILE:CG2	2.43	0.48
1:C:-1:DA:C2	2:F:-2:DT:O2	2.66	0.48
2:F:8:DT:H73	3:A:185:ARG:HH22	1.78	0.48
2:F:6:DA:O5'	2:F:6:DA:H8	1.96	0.48
3:A:10:THR:HG21	3:A:112:ILE:HG23	1.95	0.48
3:A:102:PHE:CE1	3:A:112:ILE:HD12	2.49	0.48
1:E:23:DA:H2''	1:E:22:DA:O5'	2.14	0.48
1:C:5:DG:N7	3:A:180:ARG:NH1	2.58	0.48
3:A:109:ASN:ND2	3:A:112:ILE:H	2.11	0.48
3:B:190:LEU:HA	3:B:190:LEU:HD12	1.63	0.48
3:B:10:THR:O	3:B:13:TRP:HB3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:203:ILE:O	3:B:203:ILE:HG22	2.14	0.47
3:A:37:GLU:H	3:A:82:ARG:NH1	2.12	0.47
3:B:76:PHE:CE2	3:B:120:MET:HE1	2.49	0.47
2:F:4:DA:N7	2:F:3:DT:H72	2.25	0.47
1:C:8:DA:C5	1:C:9:DT:O4	2.68	0.47
1:C:8:DA:N1	2:F:8:DT:C4	2.83	0.47
3:A:47:VAL:CG1	3:A:86:VAL:HG13	2.44	0.47
3:B:205:VAL:HG12	3:B:206:TYR:CG	2.50	0.47
2:F:8:DT:H72	3:A:181:PHE:CG	2.49	0.47
3:B:164:GLN:HG3	3:B:203:ILE:HG13	1.97	0.47
2:D:20:DT:C2'	2:D:21:DT:C6	2.97	0.47
3:B:116:LEU:HA	3:B:119:GLN:HG3	1.97	0.47
2:F:-3:DC:H2''	2:F:-4:DG:O5'	2.15	0.47
3:A:153:GLN:HG3	3:A:154:PRO:CD	2.45	0.46
3:B:156:ALA:O	3:B:163:MET:SD	2.74	0.46
3:B:78:GLU:OE2	3:B:78:GLU:HA	2.16	0.46
3:B:74:GLY:HA3	3:B:99:TYR:HE2	1.81	0.46
1:E:27:DG:N9	1:E:26:DC:C5	2.84	0.46
3:B:120:MET:HA	3:B:123:ARG:HD2	1.98	0.46
3:A:113:LEU:HD23	3:B:113:LEU:HB3	1.97	0.46
3:B:87:ARG:HG2	3:B:87:ARG:NH1	2.26	0.46
3:A:75:LEU:HD21	3:A:102:PHE:HD2	1.78	0.46
1:E:20:DA:C5	1:E:19:DT:C4	3.04	0.45
3:B:62:SER:CB	4:B:210:CMP:H2	2.45	0.45
3:B:71:GLY:HA2	4:B:210:CMP:O1P	2.17	0.45
1:C:8:DA:C2	2:F:8:DT:N3	2.85	0.45
3:B:32:GLN:HG3	3:B:85:TRP:CZ3	2.52	0.45
3:A:43:VAL:HG21	3:A:95:ALA:HB2	1.98	0.45
2:F:9:DA:C8	2:F:8:DT:C7	3.00	0.45
2:D:19:DA:C1'	2:D:20:DT:H71	2.47	0.45
3:A:23:TYR:CE2	3:A:29:LEU:HD13	2.52	0.45
3:B:61:LEU:CD2	4:B:210:CMP:C6	3.00	0.45
1:C:9:DT:H2'	2:D:10:DC:C6	2.52	0.45
1:E:17:DT:H4'	1:E:16:DG:OP1	2.17	0.45
2:F:7:DC:O5'	2:F:7:DC:C2'	2.60	0.45
3:A:36:ALA:HA	3:A:82:ARG:NH1	2.29	0.45
3:B:70:ILE:HG22	3:B:71:GLY:N	2.29	0.45
3:B:104:GLN:O	3:B:105:LEU:C	2.60	0.45
1:C:-1:DA:C2'	1:C:1:DA:OP2	2.58	0.44
1:E:18:DG:N7	3:B:180:ARG:NH2	2.65	0.44
3:A:201:LYS:HE3	3:A:203:ILE:HD11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:38:THR:O	3:B:82:ARG:NH2	2.49	0.44
1:E:16:DG:OP2	1:E:16:DG:C3'	2.60	0.44
2:F:13:DC:H2''	2:F:12:DT:OP2	2.16	0.44
3:A:72:GLU:O	3:A:75:LEU:HB2	2.17	0.44
3:A:124:LEU:CD2	4:B:210:CMP:HN61	2.28	0.44
1:E:26:DC:H2'	1:E:25:DG:N9	2.30	0.44
2:F:8:DT:H73	3:A:185:ARG:NH2	2.33	0.44
3:A:30:ILE:HG22	3:A:31:HIS:H	1.82	0.44
2:D:21:DT:N3	1:E:20:DA:C2	2.86	0.44
1:E:27:DG:C2	1:E:26:DC:C4	3.06	0.44
3:A:72:GLU:O	3:A:74:GLY:N	2.51	0.44
3:A:92:CYS:O	3:A:92:CYS:SG	2.72	0.44
3:A:40:TYR:CD1	3:A:40:TYR:N	2.86	0.44
3:B:40:TYR:HD2	3:B:94:VAL:HG11	1.83	0.44
3:B:189:MET:O	3:B:192:ASP:HB2	2.18	0.44
2:D:19:DA:N9	2:D:20:DT:H71	2.33	0.44
3:A:151:ALA:HB1	3:A:163:MET:HB2	1.99	0.44
3:B:164:GLN:HB2	3:B:203:ILE:CG1	2.48	0.44
1:C:7:DG:C6	2:F:8:DT:O4	2.71	0.43
3:A:25:SER:O	3:A:26:LYS:CB	2.66	0.43
2:D:22:DT:H2'	2:D:23:DT:H5'	1.99	0.43
3:B:64:LEU:HD12	3:B:64:LEU:HA	1.90	0.43
3:B:34:GLU:HG3	3:B:35:LYS:NZ	2.33	0.43
3:B:139:VAL:O	3:B:143:ILE:HG13	2.18	0.43
1:C:5:DG:H2''	1:C:6:DT:H6	1.77	0.43
1:E:15:DA:C4	1:E:14:DT:C5	3.07	0.43
2:F:2:DT:O2	2:F:1:DT:C2	2.71	0.43
3:A:190:LEU:HB2	3:A:196:ILE:CG2	2.48	0.43
1:E:19:DT:C7	3:B:180:ARG:NH1	2.80	0.43
3:B:40:TYR:CD1	3:B:40:TYR:N	2.86	0.43
1:C:7:DG:C6	2:F:8:DT:C4	3.07	0.43
1:C:8:DA:N3	1:C:9:DT:C2	2.87	0.43
1:E:20:DA:N7	1:E:19:DT:C7	2.81	0.43
1:C:-3:DG:O3'	1:C:-2:DA:C8	2.72	0.43
1:C:8:DA:N1	2:F:8:DT:O4	2.52	0.43
1:E:20:DA:C8	1:E:19:DT:C7	2.82	0.43
1:E:16:DG:H1'	1:E:15:DA:C4	2.53	0.43
3:B:196:ILE:HG13	3:B:197:SER:H	1.84	0.43
1:C:8:DA:H2''	1:C:9:DT:C6	2.53	0.43
1:E:15:DA:H2'	1:E:14:DT:H71	2.00	0.43
3:B:37:GLU:HB3	3:B:100:LYS:NZ	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:3:DT:N1	2:F:2:DT:H72	2.34	0.42
2:F:-2:DT:H6	2:F:-2:DT:H2'	1.47	0.42
2:F:-3:DC:C2'	2:F:-4:DG:C8	3.02	0.42
2:F:10:DG:P	3:A:139:VAL:CG1	3.00	0.42
2:F:8:DT:C7	3:A:181:PHE:CD2	3.00	0.42
3:B:18:CYS:SG	3:B:97:ILE:CD1	3.07	0.42
1:E:21:DA:N1	1:E:20:DA:C2	2.88	0.42
1:E:15:DA:C5	1:E:14:DT:C4	3.07	0.42
3:A:62:SER:O	3:A:63:TYR:HD2	2.03	0.42
3:A:159:HIS:CD2	3:A:160:PRO:HD2	2.55	0.42
3:B:169:ARG:HH22	3:B:198:ALA:HB1	1.84	0.42
3:B:207:GLY:O	3:B:208:THR:HB	2.19	0.42
1:E:19:DT:OP2	3:B:170:GLN:HB2	2.19	0.42
3:A:64:LEU:HD23	3:A:123:ARG:NH2	2.34	0.42
3:A:70:ILE:HD13	3:A:70:ILE:HG21	1.85	0.42
1:C:6:DT:O4	3:A:181:PHE:CE1	2.73	0.42
3:B:14:PHE:CD1	3:B:14:PHE:C	2.97	0.42
3:B:159:HIS:ND1	3:B:162:GLY:O	2.53	0.42
3:A:131:VAL:HG11	3:B:130:LYS:CE	2.50	0.42
3:B:72:GLU:OE1	3:B:123:ARG:NH1	2.53	0.42
3:B:166:LYS:HA	3:B:201:LYS:HA	2.01	0.42
3:B:205:VAL:CG1	3:B:206:TYR:CD2	3.03	0.42
3:A:9:PRO:HB2	3:A:12:GLU:OE1	2.19	0.42
3:A:139:VAL:O	3:A:140:THR:C	2.62	0.42
3:A:14:PHE:CD2	3:A:14:PHE:C	2.97	0.42
3:A:42:ILE:HD11	3:A:70:ILE:HD11	2.02	0.42
1:E:27:DG:H1'	1:E:26:DC:OP2	2.20	0.41
1:E:24:DA:N1	1:E:23:DA:N6	2.68	0.41
2:F:4:DA:C4	2:F:3:DT:C4	3.08	0.41
3:B:98:SER:OG	3:B:101:LYS:HB2	2.20	0.41
2:D:16:DC:H2''	2:D:17:DA:OP1	2.19	0.41
2:D:23:DT:OP1	3:B:199:HIS:C	2.64	0.41
3:B:157:MET:HB3	3:B:164:GLN:HE21	1.84	0.41
2:F:1:DT:H2''	2:F:-1:DT:OP2	2.19	0.41
3:A:158:THR:HG23	3:A:163:MET:SD	2.60	0.41
3:B:52:LYS:HD3	3:B:85:TRP:CZ2	2.56	0.41
3:B:191:GLU:O	3:B:194:ASN:N	2.52	0.41
3:A:61:LEU:O	3:A:62:SER:HB2	2.19	0.41
3:B:114:MET:O	3:B:115:ARG:C	2.64	0.41
3:B:134:LEU:CD2	3:B:142:ARG:HH11	2.29	0.41
2:D:11:DT:H1'	2:D:12:DA:C8	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:21:DT:O2	2:D:22:DT:C2	2.74	0.41
2:F:2:DT:C2	2:F:1:DT:C5	3.09	0.41
3:A:97:ILE:HB	3:A:98:SER:H	1.54	0.41
1:E:15:DA:C6	1:E:14:DT:O4	2.74	0.41
2:F:-1:DT:C2'	2:F:-2:DT:OP2	2.69	0.41
3:A:52:LYS:HE3	3:A:85:TRP:CE2	2.56	0.41
3:A:85:TRP:HE3	3:A:85:TRP:H	1.69	0.41
3:A:129:GLU:O	3:A:132:GLY:N	2.53	0.41
3:B:38:THR:OG1	3:B:40:TYR:CE1	2.72	0.41
3:B:39:LEU:N	3:B:99:TYR:HE1	2.18	0.41
2:D:10:DC:C5	2:D:11:DT:O4	2.74	0.41
3:A:46:SER:O	3:A:47:VAL:HG23	2.21	0.41
3:A:73:LEU:HD12	4:A:210:CMP:H2'	2.03	0.41
3:A:158:THR:CG2	3:A:163:MET:SD	3.09	0.41
2:D:14:DA:OP2	3:B:179:SER:HB3	2.20	0.40
2:D:17:DA:C2'	2:D:18:DC:OP2	2.53	0.40
2:F:-3:DC:H2''	2:F:-4:DG:C8	2.55	0.40
3:A:143:ILE:O	3:A:147:LEU:HD12	2.21	0.40
3:B:44:LYS:HB3	3:B:93:GLU:HB2	2.02	0.40
3:B:173:GLY:HA2	3:B:183:VAL:HG21	2.03	0.40
1:E:25:DG:C4'	1:E:24:DA:OP1	2.69	0.40
3:A:49:VAL:HG23	3:A:64:LEU:CD1	2.51	0.40
3:B:102:PHE:C	3:B:104:GLN:N	2.79	0.40
3:B:152:LYS:O	3:B:153:GLN:HB2	2.21	0.40
3:B:188:LYS:HA	3:B:188:LYS:HD3	1.84	0.40
3:A:87:ARG:HG3	3:A:88:ALA:O	2.22	0.40
3:B:137:LEU:O	3:B:142:ARG:HG3	2.21	0.40
3:B:191:GLU:N	3:B:196:ILE:HG22	2.37	0.40
3:B:196:ILE:HD12	3:B:196:ILE:HA	1.71	0.40
1:C:-4:DC:H6	1:C:-4:DC:H2'	1.63	0.40
1:C:7:DG:C2	1:C:8:DA:N1	2.89	0.40
3:B:191:GLU:HB3	3:B:196:ILE:HG22	2.04	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	196/209 (94%)	152 (78%)	30 (15%)	14 (7%)	1	1
3	B	199/209 (95%)	148 (74%)	39 (20%)	12 (6%)	1	2
All	All	395/418 (94%)	300 (76%)	69 (18%)	26 (7%)	1	1

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	26	LYS
3	A	36	ALA
3	A	75	LEU
3	A	91	ALA
3	A	199	HIS
3	B	12	GLU
3	B	62	SER
3	B	75	LEU
3	B	155	ASP
3	B	199	HIS
3	B	205	VAL
3	B	34	GLU
3	A	34	GLU
3	A	97	ILE
3	A	160	PRO
3	A	194	ASN
3	B	200	GLY
3	B	208	THR
3	B	66	GLN
3	B	193	GLN
3	A	54	GLU
3	B	139	VAL
3	A	167	ILE
3	A	193	GLN
3	A	42	ILE
3	A	165	ILE

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	171/180 (95%)	116 (68%)	55 (32%)	0	1
3	B	172/180 (96%)	130 (76%)	42 (24%)	1	2
All	All	343/360 (95%)	246 (72%)	97 (28%)	0	1

All (97) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	10	THR
3	A	11	LEU
3	A	12	GLU
3	A	14	PHE
3	A	15	LEU
3	A	18	CYS
3	A	22	LYS
3	A	26	LYS
3	A	27	SER
3	A	29	LEU
3	A	34	GLU
3	A	37	GLU
3	A	39	LEU
3	A	40	TYR
3	A	42	ILE
3	A	46	SER
3	A	47	VAL
3	A	52	LYS
3	A	53	ASP
3	A	54	GLU
3	A	58	GLU
3	A	75	LEU
3	A	78	GLU
3	A	81	GLU
3	A	87	ARG
3	A	90	THR
3	A	97	ILE
3	A	100	LYS
3	A	105	LEU
3	A	106	ILE
3	A	111	ASP

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Mol	Chain	Res	Type
3	A	115	ARG
3	A	120	MET
3	A	126	VAL
3	A	129	GLU
3	A	137	LEU
3	A	143	ILE
3	A	145	GLN
3	A	153	GLN
3	A	154	PRO
3	A	157	MET
3	A	158	THR
3	A	160	PRO
3	A	164	GLN
3	A	165	ILE
3	A	166	LYS
3	A	168	THR
3	A	170	GLN
3	A	174	GLN
3	A	185	ARG
3	A	187	LEU
3	A	189	MET
3	A	196	ILE
3	A	201	LYS
3	A	202	THR
3	B	12	GLU
3	B	14	PHE
3	B	15	LEU
3	B	27	SER
3	B	29	LEU
3	B	34	GLU
3	B	35	LYS
3	B	39	LEU
3	B	40	TYR
3	B	46	SER
3	B	47	VAL
3	B	58	GLU
3	B	60	ILE
3	B	64	LEU
3	B	72	GLU
3	B	73	LEU
3	B	78	GLU
3	B	97	ILE

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Mol	Chain	Res	Type
3	B	101	LYS
3	B	103	ARG
3	B	105	LEU
3	B	112	ILE
3	B	116	LEU
3	B	120	MET
3	B	122	ARG
3	B	124	LEU
3	B	125	GLN
3	B	147	LEU
3	B	148	LEU
3	B	158	THR
3	B	164	GLN
3	B	165	ILE
3	B	167	ILE
3	B	171	GLU
3	B	174	GLN
3	B	175	ILE
3	B	176	VAL
3	B	196	ILE
3	B	197	SER
3	B	203	ILE
3	B	204	VAL
3	B	205	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	19	HIS
3	A	65	ASN
3	A	107	GLN
3	B	17	HIS
3	B	32	GLN
3	B	65	ASN
3	B	109	ASN
3	B	164	GLN
3	B	170	GLN

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](#)

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	1.58	3 (12%)	37,39,39	2.10	7 (18%)
4	CMP	A	210	-	25,25,25	1.05	3 (12%)	37,39,39	2.25	7 (18%)

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 (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	210	CMP	O3'-C3'	-5.28	1.36	1.44
4	B	210	CMP	P-O5'	3.35	1.61	1.57
4	A	210	CMP	P-O5'	2.53	1.60	1.57
4	A	210	CMP	P-O3'	2.38	1.61	1.57
4	B	210	CMP	C2-N3	-2.23	1.30	1.33
4	A	210	CMP	C2-N3	-2.11	1.30	1.33

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	210	CMP	O3'-C3'-C2'	10.00	125.42	115.61
4	B	210	CMP	O3'-C3'-C4'	-7.11	105.34	110.71
4	B	210	CMP	O3'-P-O1P	-6.23	97.04	110.39
4	B	210	CMP	C2'-C1'-N9	-4.43	102.29	113.30
4	A	210	CMP	O2P-P-O1P	4.38	122.01	108.56
4	A	210	CMP	C5'-C4'-C3'	-4.08	104.89	112.56
4	A	210	CMP	O2P-P-O3'	3.61	115.44	107.04
4	B	210	CMP	O2P-P-O1P	3.19	118.37	108.56
4	A	210	CMP	O3'-P-O1P	-3.05	103.87	110.39
4	B	210	CMP	O5'-P-O1P	2.87	117.07	110.44
4	A	210	CMP	O2P-P-O5'	-2.45	101.17	107.16
4	A	210	CMP	O3'-C3'-C4'	-2.40	108.90	110.71
4	B	210	CMP	O5'-P-O3'	-2.35	102.55	105.70
4	B	210	CMP	O4'-C4'-C3'	-2.25	100.16	104.92

There are no chirality outliers.

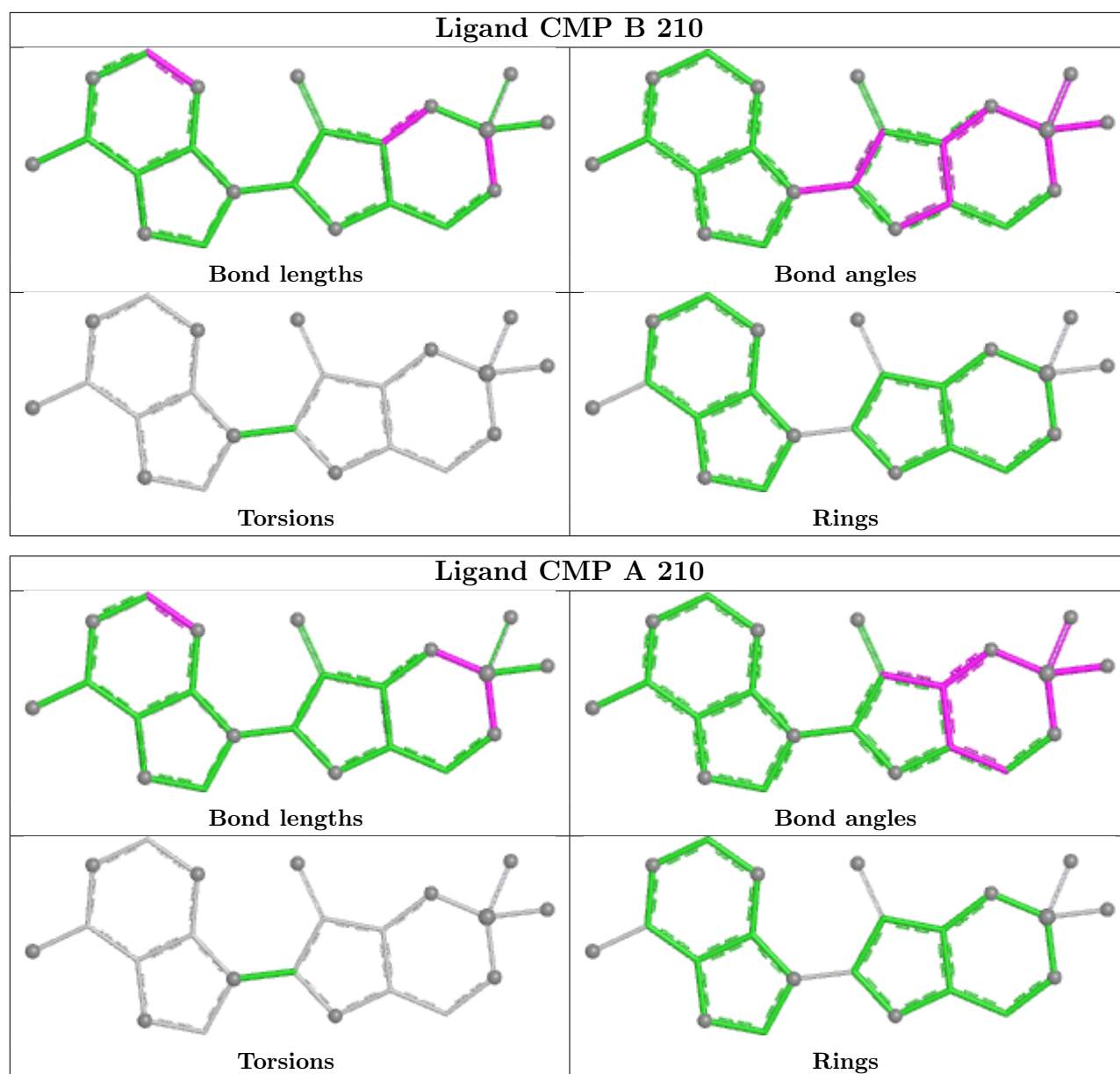
There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	210	CMP	11	0
4	A	210	CMP	9	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	C	14/14 (100%)	-0.35	0 100 100	26, 37, 49, 54	0
1	E	14/14 (100%)	0.07	1 (7%) 22 19	32, 44, 52, 53	0
2	D	17/17 (100%)	-0.02	0 100 100	31, 45, 59, 61	0
2	F	17/17 (100%)	-0.17	0 100 100	22, 35, 45, 52	0
3	A	198/209 (94%)	-0.58	1 (0%) 87 86	6, 20, 43, 57	0
3	B	201/209 (96%)	-0.52	2 (0%) 79 78	6, 22, 43, 58	1 (0%)
All	All	461/480 (96%)	-0.49	4 (0%) 81 80	6, 24, 48, 61	1 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	162	GLY	3.4
1	E	26	DC	2.2
3	A	16	SER	2.2
3	B	156	ALA	2.2

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.

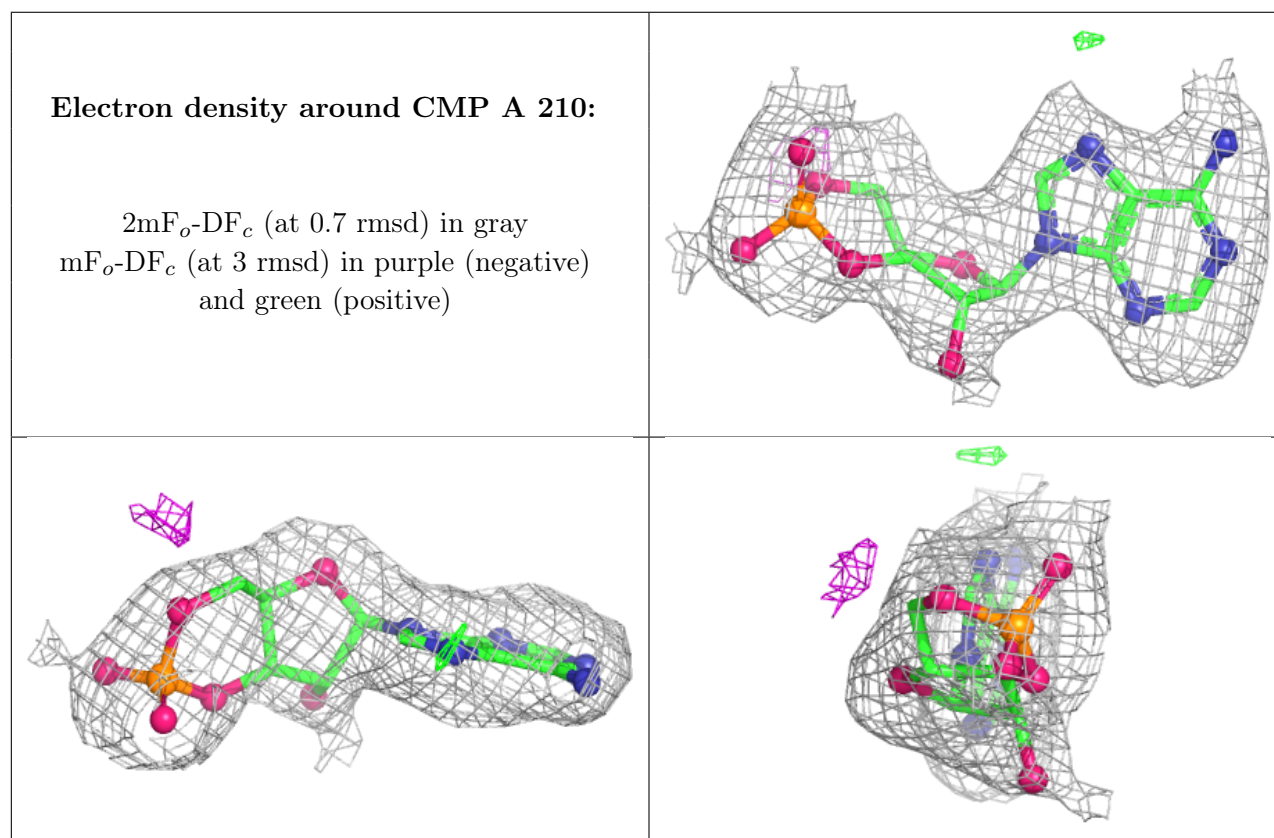
6.4 Ligands [i](#)

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.

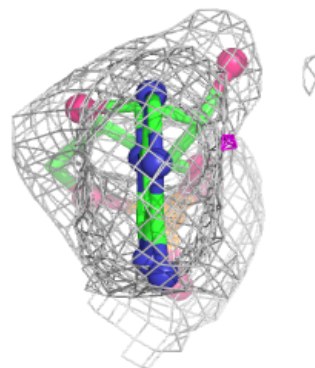
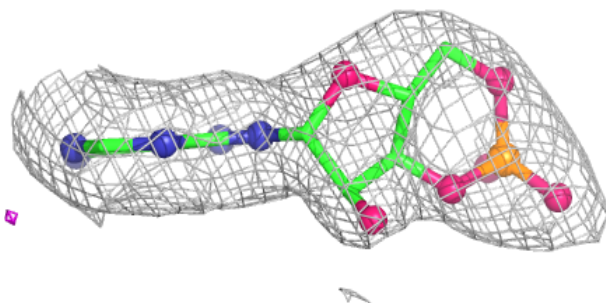
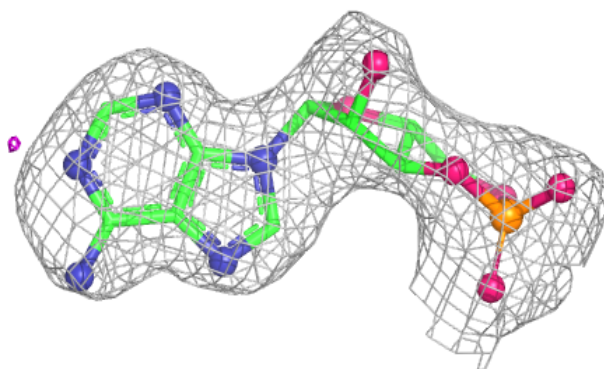
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	CMP	A	210	22/22	0.98	0.05	6,13,22,28	0
4	CMP	B	210	22/22	0.99	0.04	6,10,15,18	0

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 210:

$2mF_o - DF_c$ (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.