



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:08 AM UTC

PDB ID : 1U0H / pdb_00001u0h
Title : STRUCTURAL BASIS FOR THE INHIBITION OF MAMMALIAN ADENYLYL CYCLASE BY MANT-GTP
Authors : Mou, T.C.; Gille, A.; Seifert, R.J.; Sprang, S.R.
Deposited on : 2004-07-13
Resolution : 2.90 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

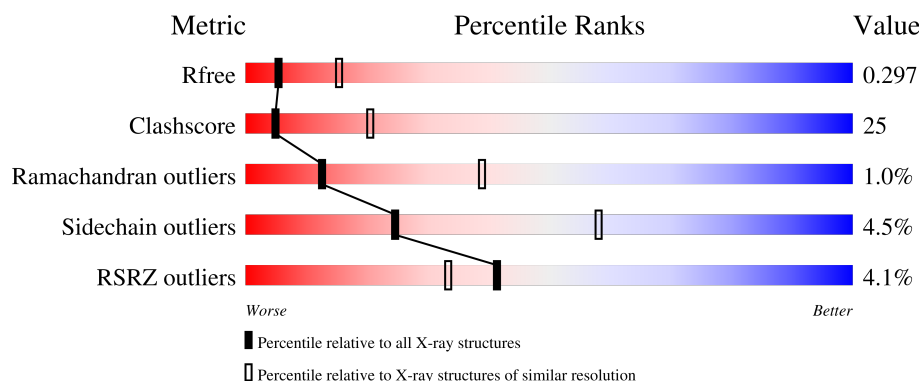
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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

Mol	Chain	Length	Quality of chain
1	A	225	4% 38% 43% • 16%
2	B	212	8% 49% 38% • 11%
3	C	394	% 47% 34% • 16%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 5769 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Adenylate cyclase, type V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	189	Total	C	N	O	S	0	0	0
			1476	929	259	271	17			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	356	MET	-	initiating methionine	UNP P30803
A	357	HIS	-	expression tag	UNP P30803
A	358	HIS	-	expression tag	UNP P30803
A	359	HIS	-	expression tag	UNP P30803
A	360	HIS	-	expression tag	UNP P30803
A	362	HIS	VAL	conflict	UNP P30803
A	476	MET	VAL	conflict	UNP P30803

- Molecule 2 is a protein called Adenylate cyclase, type II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	189	Total	C	N	O	S	0	0	0
			1467	936	242	279	10			

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(s), alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	330	Total	C	N	O	S	0	0	0
			2702	1714	470	505	13			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

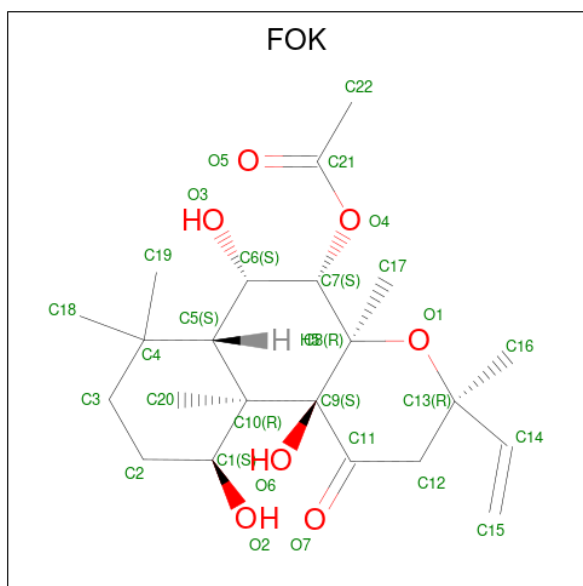
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		

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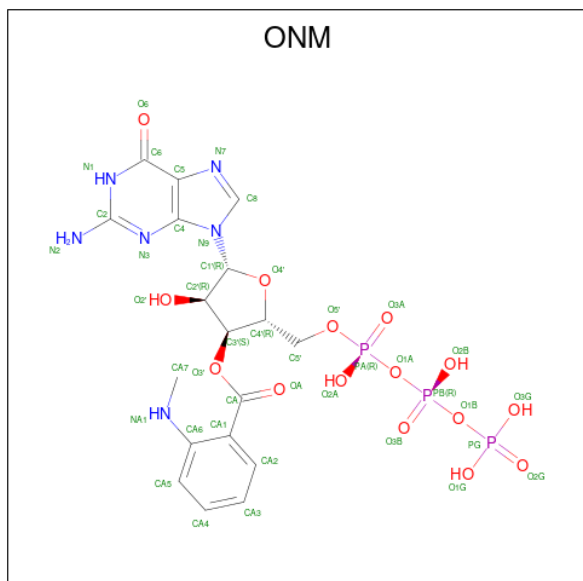
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	Mg	0	0
			1	1		

- Molecule 5 is FORSKOLIN (CCD ID: FOK) (formula: $C_{22}H_{34}O_7$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	C	O	
			29	22	7	

- Molecule 6 is 3'-O-(N-METHYLANTHRANILOYL)-GUANOSINE-5'-TRIPHOSPHATE (CCD ID: ONM) (formula: $C_{18}H_{23}N_6O_{15}P_3$).

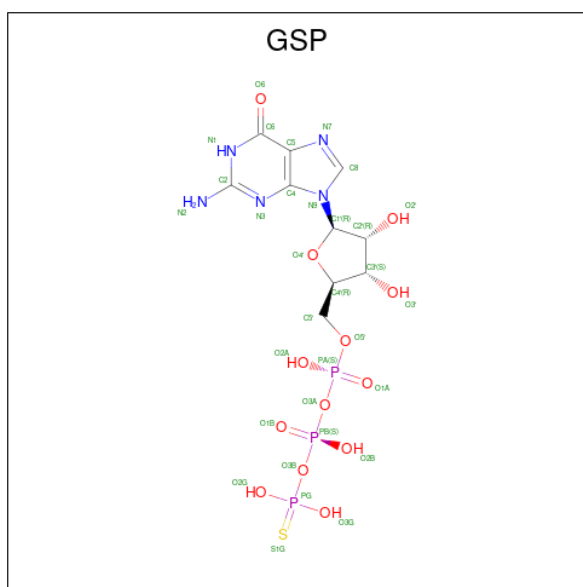


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			42	18	6	15	3		

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	1	Total	Cl	0	0
			1	1		

- Molecule 8 is 5'-GUANOSINE-DIPHOSPHATE-MONOTHIOPHOSPHATE (CCD ID: GSP) (formula: C₁₀H₁₆N₅O₁₃P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
8	C	1	Total	C	N	O	P	S	0	0
			32	10	5	13	3	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	4	Total	O	0	0
			4	4		
9	B	7	Total	O	0	0
			7	7		
9	C	6	Total	O	0	0
			6	6		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	118.40Å 133.00Å 70.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.90 15.00 – 2.91	Depositor EDS
% Data completeness (in resolution range)	88.1 (15.00-2.90) 87.3 (15.00-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.91Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.245 , 0.280 0.265 , 0.297	Depositor DCC
R_{free} test set	1064 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	40.9	Xtriage
Anisotropy	1.226	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 62.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	5769	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ONM, FOK, GSP, CL, MG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.44	0/1504	0.91	4/2027 (0.2%)
2	B	0.50	0/1492	0.95	3/2014 (0.1%)
3	C	0.46	0/2759	0.95	15/3733 (0.4%)
All	All	0.47	0/5755	0.94	22/7774 (0.3%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	137	VAL	CA-C-N	7.53	129.25	119.84
3	C	137	VAL	C-N-CA	7.53	129.25	119.84
2	B	1045	THR	N-CA-C	-5.88	104.87	111.28
3	C	259	GLU	N-CA-C	5.56	117.34	111.28
3	C	320	THR	CA-C-N	5.55	126.77	119.84
3	C	320	THR	C-N-CA	5.55	126.77	119.84
2	B	985	ALA	N-CA-C	-5.54	105.15	111.14
1	A	447	GLY	N-CA-C	-5.52	106.91	114.64
1	A	421	ALA	N-CA-C	-5.43	105.27	111.14
3	C	310	ASP	N-CA-C	-5.34	107.14	113.97
1	A	530	LYS	N-CA-C	-5.30	106.07	112.54
3	C	120	ALA	N-CA-C	-5.29	106.82	113.28
3	C	311	TYR	N-CA-C	-5.17	107.52	113.88
2	B	998	VAL	N-CA-C	5.16	116.79	108.85
3	C	92	VAL	N-CA-C	-5.16	105.68	110.53
3	C	63	LEU	N-CA-C	5.15	117.68	111.40
3	C	340	PHE	N-CA-C	-5.15	105.36	110.97
1	A	418	GLU	N-CA-C	-5.15	105.36	110.97
3	C	209	GLU	N-CA-C	5.10	117.01	107.99
3	C	378	ASP	N-CA-C	-5.05	106.95	113.01
3	C	312	PHE	CA-C-N	5.04	125.31	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	312	PHE	C-N-CA	5.04	125.31	119.47

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1476	0	1450	103	0
2	B	1467	0	1470	74	0
3	C	2702	0	2651	124	0
4	A	2	0	0	0	0
4	C	1	0	0	0	0
5	A	29	0	34	2	0
6	A	42	0	19	2	0
7	C	1	0	0	0	0
8	C	32	0	12	0	0
9	A	4	0	0	0	0
9	B	7	0	0	2	0
9	C	6	0	0	5	0
All	All	5769	0	5636	287	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 25.

All (287) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:411:GLU:HA	1:A:414:MET:HE2	1.43	1.00
3:C:119:LEU:H	3:C:119:LEU:HD12	1.40	0.87
1:A:378:MET:HE1	3:C:283:ARG:N	1.91	0.85
1:A:554:LEU:HA	1:A:559:ILE:HD12	1.59	0.84
3:C:207:ILE:HG12	3:C:224:VAL:HG12	1.61	0.82
1:A:530:LYS:HD2	1:A:531:ALA:N	1.93	0.81
2:B:1030:MET:HE2	2:B:1042:THR:HG23	1.63	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:456:ALA:HB2	1:A:535:TYR:HB3	1.61	0.81
1:A:508:SER:OG	1:A:511:VAL:HG23	1.81	0.80
2:B:1064:VAL:HG21	2:B:1070:LEU:HD12	1.61	0.79
1:A:456:ALA:HB1	1:A:536:LEU:HD12	1.63	0.79
1:A:530:LYS:HD2	1:A:531:ALA:H	1.45	0.78
3:C:99:LEU:HD11	3:C:182:ILE:HD13	1.65	0.77
3:C:127:ARG:HH11	3:C:149:HIS:HA	1.48	0.77
3:C:100:LYS:O	3:C:104:GLU:HG2	1.84	0.77
3:C:364:THR:HG22	3:C:375:VAL:HG11	1.66	0.76
3:C:257:ILE:HD12	3:C:259:GLU:HB2	1.67	0.75
2:B:953:ILE:HD12	2:B:953:ILE:H	1.51	0.74
3:C:255:MET:HB2	3:C:265:ARG:HD2	1.69	0.74
2:B:1043:GLU:O	2:B:1046:SER:HB3	1.88	0.74
1:A:404:ALA:HA	1:A:412:LEU:HD11	1.72	0.72
1:A:424:ASP:HB3	2:B:1013:GLN:HG2	1.70	0.72
5:A:101:FOK:H202	5:A:101:FOK:H193	1.71	0.71
1:A:526:ILE:N	1:A:526:ILE:HD12	2.05	0.71
1:A:391:SER:O	1:A:445:VAL:HG23	1.90	0.70
1:A:401:THR:HA	1:A:404:ALA:HB3	1.73	0.70
1:A:528:ILE:HD11	1:A:562:PHE:HB2	1.73	0.70
2:B:1007:ALA:HA	2:B:1019:ILE:HG22	1.74	0.69
1:A:529:THR:HG22	1:A:530:LYS:HD2	1.73	0.69
1:A:470:ILE:HD13	1:A:470:ILE:O	1.93	0.69
3:C:87:GLU:O	3:C:90:THR:HG22	1.92	0.69
1:A:546:CYS:O	1:A:549:GLU:HG2	1.93	0.69
1:A:395:ALA:HB1	1:A:483:MET:HE3	1.75	0.67
1:A:473:VAL:O	1:A:477:THR:HG23	1.94	0.67
3:C:107:VAL:HA	3:C:110:MET:HG3	1.75	0.67
5:A:101:FOK:H201	5:A:101:FOK:H173	1.77	0.66
1:A:452:ARG:HG2	1:A:454:ASP:H	1.60	0.66
3:C:43:LEU:HD11	3:C:219:PHE:HD2	1.61	0.66
2:B:940:ILE:H	2:B:940:ILE:HD12	1.61	0.65
1:A:397:ILE:HD13	1:A:483:MET:SD	2.36	0.65
2:B:1044:GLU:HA	2:B:1047:LEU:HD12	1.78	0.64
3:C:380:ARG:O	3:C:384:GLN:HG2	1.98	0.64
3:C:58:LYS:HA	3:C:61:ARG:HH12	1.60	0.64
2:B:1009:VAL:HG11	2:B:1015:PRO:HB3	1.80	0.63
2:B:953:ILE:HD12	2:B:953:ILE:N	2.13	0.63
3:C:365:CYS:H	3:C:371:ASN:ND2	1.97	0.62
2:B:915:LEU:HA	2:B:918:ILE:HD12	1.79	0.62
1:A:500:ARG:HG3	2:B:923:ASP:HB3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:121:ASN:HB2	3:C:124:ASN:HD22	1.65	0.62
1:A:426:LEU:HB2	1:A:462:MET:HE1	1.82	0.62
2:B:1030:MET:HE1	2:B:1041:VAL:C	2.25	0.62
2:B:1050:GLN:C	2:B:1052:LEU:H	2.06	0.62
3:C:63:LEU:HD21	3:C:369:THR:HG22	1.83	0.60
3:C:273:PHE:HE1	3:C:287:VAL:HG11	1.66	0.60
3:C:266:LEU:HD23	3:C:312:PHE:CZ	2.36	0.60
1:A:403:LEU:HD22	1:A:481:VAL:CG1	2.31	0.60
2:B:891:SER:HB2	9:B:8:HOH:O	2.01	0.60
1:A:463:GLY:O	1:A:467:ILE:HG12	2.00	0.60
3:C:166:SER:HA	3:C:169:TYR:CE2	2.37	0.60
1:A:384:ILE:HG22	1:A:385:GLN:N	2.17	0.59
1:A:378:MET:HE1	3:C:283:ARG:H	1.67	0.59
3:C:124:ASN:OD1	3:C:127:ARG:HD2	2.03	0.59
3:C:119:LEU:HD12	3:C:119:LEU:N	2.14	0.59
1:A:552:ALA:O	1:A:556:GLU:HG3	2.03	0.58
2:B:1009:VAL:HG12	2:B:1010:ILE:N	2.17	0.58
1:A:426:LEU:CB	1:A:462:MET:HE1	2.34	0.58
3:C:170:GLN:HG3	9:C:400:HOH:O	2.02	0.58
1:A:387:HIS:CG	1:A:448:LEU:HD21	2.39	0.57
1:A:452:ARG:HD3	1:A:454:ASP:HB3	1.85	0.57
3:C:121:ASN:HD22	3:C:124:ASN:ND2	2.03	0.57
3:C:61:ARG:HH11	3:C:61:ARG:HB2	1.70	0.57
2:B:888:MET:HG2	2:B:889:PHE:N	2.18	0.57
2:B:906:LYS:HB2	2:B:909:LEU:HB3	1.86	0.57
1:A:474:ARG:HG3	1:A:481:VAL:HG22	1.85	0.56
3:C:307:LYS:HB2	3:C:310:ASP:OD2	2.05	0.56
1:A:528:ILE:CD1	1:A:562:PHE:HB2	2.35	0.56
1:A:460:VAL:O	1:A:464:MET:HG2	2.05	0.56
3:C:203:LEU:HD23	3:C:204:THR:N	2.20	0.56
3:C:302:LEU:HA	3:C:333:ARG:HH21	1.69	0.56
2:B:1009:VAL:HG12	2:B:1010:ILE:H	1.70	0.56
1:A:423:PHE:HA	1:A:426:LEU:HD12	1.88	0.55
1:A:435:ILE:HD11	1:A:445:VAL:HG12	1.88	0.55
1:A:532:THR:O	1:A:536:LEU:HD13	2.07	0.55
3:C:45:LEU:HD11	3:C:221:MET:HG3	1.89	0.55
2:B:948:THR:HG23	2:B:972:MET:SD	2.46	0.55
3:C:99:LEU:HD21	3:C:182:ILE:HG23	1.88	0.55
3:C:219:PHE:O	3:C:220:HIS:HD2	1.89	0.55
1:A:423:PHE:CE2	1:A:483:MET:HE2	2.41	0.55
3:C:151:LYS:O	3:C:155:GLU:HG2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:371:ASN:O	3:C:375:VAL:HG22	2.07	0.55
1:A:497:LEU:HD22	2:B:916:ASN:HA	1.89	0.54
2:B:886:CYS:HB2	2:B:969:ILE:HD13	1.89	0.54
2:B:1030:MET:HE3	2:B:1040:GLN:O	2.07	0.54
3:C:61:ARG:HA	3:C:65:VAL:HB	1.89	0.54
3:C:143:PRO:HB2	3:C:145:GLU:OE2	2.08	0.54
1:A:384:ILE:HG22	1:A:385:GLN:H	1.73	0.54
2:B:975:PHE:O	2:B:978:ALA:HB3	2.07	0.54
3:C:337:ALA:O	3:C:340:PHE:HB3	2.08	0.54
3:C:166:SER:HA	3:C:169:TYR:CZ	2.43	0.54
3:C:360:TYR:OH	3:C:386:MET:HG3	2.07	0.54
1:A:420:PHE:HA	1:A:423:PHE:HB2	1.90	0.53
1:A:476:MET:SD	1:A:476:MET:N	2.82	0.53
3:C:59:GLN:OE1	3:C:59:GLN:HA	2.09	0.53
3:C:308:ILE:C	3:C:310:ASP:H	2.16	0.53
3:C:124:ASN:O	3:C:128:VAL:HG23	2.08	0.53
3:C:279:ASN:O	3:C:283:ARG:HG3	2.08	0.53
1:A:529:THR:HG22	1:A:530:LYS:N	2.24	0.52
6:A:100:ONM:H3'	2:B:1025:ASN:HD22	1.74	0.52
2:B:953:ILE:H	2:B:953:ILE:CD1	2.21	0.52
1:A:425:LYS:O	1:A:429:GLU:HG3	2.09	0.52
1:A:499:LEU:HD12	1:A:500:ARG:HG2	1.92	0.52
2:B:1059:ARG:NH1	2:B:1062:ILE:HG13	2.25	0.52
1:A:378:MET:HE2	3:C:284:THR:OG1	2.10	0.52
1:A:435:ILE:HD11	1:A:445:VAL:CG1	2.40	0.52
1:A:527:HIS:NE2	1:A:561:THR:HB	2.25	0.52
3:C:127:ARG:O	3:C:131:ILE:HG12	2.10	0.52
3:C:207:ILE:HD13	3:C:237:CYS:SG	2.50	0.52
3:C:176:GLN:HB3	9:C:403:HOH:O	2.09	0.52
1:A:461:GLU:O	1:A:464:MET:HB2	2.09	0.51
1:A:493:HIS:HB2	1:A:507:TRP:O	2.10	0.51
1:A:525:ARG:NH2	1:A:565:LEU:HD23	2.25	0.51
2:B:1064:VAL:CG2	2:B:1070:LEU:HD12	2.36	0.51
6:A:100:ONM:HA3	2:B:1020:TRP:HB3	1.92	0.51
1:A:505:ASP:HB3	1:A:507:TRP:CZ2	2.45	0.51
3:C:104:GLU:HA	3:C:107:VAL:HG22	1.93	0.51
3:C:386:MET:HB2	3:C:388:LEU:HG	1.91	0.51
1:A:513:LEU:HD13	1:A:559:ILE:HD13	1.91	0.51
3:C:44:LEU:HD12	3:C:222:PHE:HB3	1.91	0.51
2:B:888:MET:HG3	2:B:1000:ILE:HG12	1.93	0.51
3:C:284:THR:C	3:C:357:HIS:HB3	2.36	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1026:VAL:HG22	2:B:1064:VAL:HG11	1.93	0.50
3:C:58:LYS:HA	3:C:61:ARG:NH1	2.26	0.50
2:B:1030:MET:CE	2:B:1041:VAL:C	2.85	0.50
2:B:1048:ILE:O	2:B:1052:LEU:HG	2.12	0.50
1:A:470:ILE:HD13	1:A:470:ILE:C	2.36	0.50
3:C:165:ARG:HB3	3:C:168:GLU:OE2	2.11	0.50
3:C:294:GLN:O	3:C:297:LEU:HB3	2.12	0.50
1:A:525:ARG:HH21	1:A:565:LEU:HD23	1.77	0.49
1:A:529:THR:HG22	1:A:530:LYS:CD	2.40	0.49
3:C:241:VAL:CG1	3:C:285:ILE:HD13	2.43	0.49
1:A:431:HIS:HB3	1:A:452:ARG:CZ	2.41	0.49
1:A:427:ALA:HA	1:A:462:MET:SD	2.52	0.49
2:B:949:GLY:HA3	2:B:968:HIS:HD2	1.77	0.49
3:C:173:ASP:C	3:C:175:ALA:H	2.19	0.49
1:A:382:ILE:HD12	1:A:382:ILE:O	2.13	0.49
2:B:923:ASP:C	2:B:925:LEU:H	2.19	0.49
1:A:403:LEU:HD22	1:A:481:VAL:HG12	1.93	0.49
2:B:1049:LEU:O	2:B:1054:TYR:HB2	2.13	0.49
3:C:104:GLU:HB2	3:C:135:MET:HE2	1.95	0.49
2:B:964:ARG:HA	2:B:964:ARG:NE	2.27	0.48
1:A:491:ARG:NH2	2:B:907:GLU:HG3	2.29	0.48
2:B:886:CYS:O	2:B:947:ALA:HA	2.13	0.48
2:B:906:LYS:O	2:B:907:GLU:HB2	2.13	0.48
1:A:403:LEU:HB2	1:A:479:VAL:HG11	1.95	0.48
2:B:918:ILE:HG12	2:B:986:ILE:HD13	1.96	0.48
3:C:99:LEU:HD11	3:C:182:ILE:CD1	2.42	0.48
3:C:45:LEU:N	3:C:45:LEU:HD12	2.29	0.48
3:C:43:LEU:HD12	3:C:221:MET:CE	2.44	0.48
3:C:174:CYS:HB2	3:C:178:PHE:CD2	2.49	0.48
1:A:474:ARG:HG3	1:A:481:VAL:CG2	2.44	0.48
2:B:998:VAL:HG12	2:B:999:GLY:N	2.29	0.47
2:B:1067:LYS:HG2	2:B:1070:LEU:HD21	1.95	0.47
2:B:890:ALA:HB1	2:B:996:LEU:HD21	1.95	0.47
3:C:191:VAL:O	3:C:191:VAL:HG13	2.14	0.47
2:B:984:ASP:HA	2:B:987:ASN:HD22	1.79	0.47
3:C:254:ASN:HA	3:C:311:TYR:CE2	2.48	0.47
1:A:541:GLU:CG	1:A:565:LEU:HD12	2.44	0.47
2:B:929:PRO:HD2	9:B:26:HOH:O	2.14	0.47
3:C:307:LYS:HB3	3:C:309:GLU:OE2	2.14	0.47
1:A:423:PHE:HE2	1:A:483:MET:HE2	1.80	0.47
2:B:1030:MET:HE2	2:B:1042:THR:CG2	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:43:LEU:HD11	3:C:219:PHE:CD2	2.45	0.47
3:C:44:LEU:HD12	3:C:222:PHE:CB	2.44	0.47
1:A:470:ILE:HD11	1:A:481:VAL:HG23	1.96	0.47
3:C:119:LEU:H	3:C:119:LEU:CD1	2.16	0.47
2:B:1007:ALA:CA	2:B:1019:ILE:HG22	2.44	0.47
3:C:166:SER:HB2	3:C:171:LEU:HD22	1.96	0.47
2:B:1043:GLU:O	2:B:1047:LEU:HG	2.15	0.47
2:B:1055:THR:OG1	2:B:1076:ASN:HB2	2.15	0.47
2:B:1049:LEU:HA	2:B:1052:LEU:HD12	1.97	0.47
3:C:166:SER:HB2	3:C:171:LEU:CD2	2.45	0.47
3:C:250:SER:HB2	3:C:292:ASN:O	2.14	0.47
3:C:292:ASN:CG	3:C:293:LYS:H	2.23	0.46
2:B:895:PHE:CZ	2:B:915:LEU:HB2	2.51	0.46
1:A:548:GLY:O	1:A:555:LYS:HB2	2.15	0.46
2:B:995:LYS:HD2	2:B:1037:ASP:OD2	2.16	0.46
3:C:121:ASN:HB2	3:C:124:ASN:ND2	2.28	0.46
1:A:393:LEU:C	1:A:393:LEU:HD23	2.40	0.46
3:C:203:LEU:HD23	3:C:204:THR:O	2.16	0.46
3:C:376:PHE:O	3:C:380:ARG:HG3	2.15	0.46
2:B:1007:ALA:HB1	2:B:1017:TYR:OH	2.16	0.46
3:C:176:GLN:N	9:C:403:HOH:O	2.49	0.46
3:C:264:ASN:OD1	3:C:266:LEU:N	2.49	0.46
2:B:882:TYR:CD1	2:B:882:TYR:N	2.84	0.46
3:C:57:VAL:HG22	3:C:221:MET:HG2	1.98	0.46
2:B:1029:ARG:HH12	2:B:1033:THR:HG23	1.82	0.45
3:C:103:ILE:HG23	3:C:104:GLU:N	2.31	0.45
1:A:526:ILE:N	1:A:526:ILE:CD1	2.75	0.45
2:B:1050:GLN:C	2:B:1052:LEU:N	2.72	0.45
3:C:131:ILE:HG13	3:C:153:LEU:CD1	2.46	0.45
1:A:382:ILE:HG12	2:B:913:ARG:NH1	2.31	0.45
1:A:494:CYS:HB3	1:A:506:VAL:HG12	1.97	0.45
2:B:949:GLY:HA3	2:B:968:HIS:CD2	2.52	0.45
3:C:55:THR:HG22	3:C:366:ALA:HB1	1.99	0.45
3:C:330:GLU:OE2	3:C:338:LYS:HE2	2.16	0.45
2:B:888:MET:SD	2:B:972:MET:HE2	2.57	0.44
2:B:1000:ILE:HD12	2:B:1039:ILE:HG21	1.98	0.44
1:A:395:ALA:CB	1:A:483:MET:HE3	2.46	0.44
1:A:550:ARG:HB2	1:A:550:ARG:NH1	2.33	0.44
3:C:187:GLN:HB3	3:C:189:ASP:OD1	2.18	0.44
3:C:292:ASN:HA	3:C:364:THR:O	2.17	0.44
1:A:382:ILE:HG22	2:B:916:ASN:ND2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:ILE:HD11	1:A:492:VAL:CG1	2.48	0.44
1:A:397:ILE:HD11	1:A:419:LEU:CD2	2.48	0.44
2:B:934:VAL:HA	2:B:948:THR:HG22	2.00	0.44
2:B:936:LYS:HE3	2:B:939:THR:HG23	2.00	0.44
3:C:143:PRO:HB2	3:C:145:GLU:CD	2.43	0.43
3:C:195:GLN:HA	3:C:195:GLN:NE2	2.33	0.43
1:A:378:MET:CE	3:C:283:ARG:N	2.74	0.43
1:A:400:PHE:HA	1:A:403:LEU:HD23	1.99	0.43
3:C:193:SER:O	3:C:197:LEU:HG	2.17	0.43
3:C:255:MET:CB	3:C:265:ARG:HD2	2.45	0.43
1:A:442:TYR:CD1	1:A:442:TYR:C	2.96	0.43
3:C:59:GLN:HG3	3:C:372:ILE:HG13	2.00	0.43
2:B:904:VAL:HG11	3:C:207:ILE:HB	2.00	0.43
3:C:260:ASP:OD1	3:C:260:ASP:C	2.61	0.43
3:C:334:VAL:O	3:C:335:THR:C	2.61	0.43
2:B:989:HIS:CG	3:C:280:ARG:HG2	2.54	0.43
3:C:277:TRP:NE1	3:C:349:SER:HA	2.33	0.43
3:C:108:ALA:HB2	3:C:135:MET:HE1	2.00	0.43
3:C:210:THR:HB	3:C:221:MET:HB3	2.01	0.43
1:A:488:HIS:HB3	1:A:514:ALA:HB2	2.01	0.43
1:A:530:LYS:CD	1:A:531:ALA:N	2.75	0.43
3:C:214:VAL:HG11	3:C:380:ARG:CZ	2.49	0.42
1:A:422:ARG:O	1:A:426:LEU:HG	2.19	0.42
1:A:409:ALA:O	1:A:413:VAL:HG23	2.18	0.42
1:A:529:THR:HG22	1:A:530:LYS:H	1.84	0.42
2:B:1040:GLN:OE1	2:B:1059:ARG:HD2	2.20	0.42
1:A:422:ARG:HH21	1:A:425:LYS:CD	2.32	0.42
1:A:530:LYS:HD2	1:A:530:LYS:N	2.34	0.42
3:C:254:ASN:HB2	3:C:311:TYR:CZ	2.53	0.42
3:C:365:CYS:HB2	9:C:399:HOH:O	2.18	0.42
1:A:397:ILE:O	1:A:398:GLU:C	2.62	0.42
1:A:456:ALA:O	1:A:460:VAL:HG23	2.19	0.42
1:A:493:HIS:O	1:A:507:TRP:HE3	2.02	0.42
2:B:1033:THR:HB	2:B:1059:ARG:HH12	1.85	0.42
1:A:475:GLU:HB2	1:A:476:MET:SD	2.60	0.42
1:A:499:LEU:HA	1:A:502:TRP:HE1	1.84	0.42
3:C:253:TYR:HB3	9:C:401:HOH:O	2.20	0.42
3:C:331:ASP:OD1	3:C:333:ARG:HG3	2.19	0.42
1:A:432:CYS:SG	1:A:462:MET:HB2	2.60	0.42
1:A:512:THR:HG22	1:A:516:HIS:HD2	1.85	0.41
1:A:513:LEU:CD1	1:A:559:ILE:HD13	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:172:ILE:HD12	3:C:174:CYS:SG	2.60	0.41
3:C:235:ILE:O	3:C:235:ILE:HD13	2.20	0.41
1:A:378:MET:HE1	3:C:284:THR:H	1.86	0.41
1:A:426:LEU:HD13	1:A:465:ASP:HB3	2.01	0.41
1:A:456:ALA:HB1	1:A:536:LEU:CD1	2.43	0.41
3:C:50:GLU:HG2	3:C:201:ARG:NH2	2.36	0.41
3:C:260:ASP:OD1	3:C:262:GLN:N	2.53	0.41
3:C:147:TYR:HD1	3:C:182:ILE:HG22	1.86	0.41
2:B:998:VAL:CG1	2:B:999:GLY:N	2.84	0.41
3:C:228:ARG:C	3:C:230:GLU:N	2.78	0.41
3:C:292:ASN:CG	3:C:293:LYS:N	2.79	0.41
3:C:281:TRP:O	3:C:282:LEU:HD23	2.19	0.41
1:A:435:ILE:HG13	1:A:444:CYS:HA	2.03	0.41
2:B:905:ASN:HB3	3:C:236:GLN:OE1	2.20	0.41
3:C:90:THR:HG23	3:C:91:LYS:HG3	2.02	0.41
3:C:134:VAL:HA	3:C:137:VAL:HG23	2.02	0.41
3:C:134:VAL:HB	3:C:140:PHE:CE1	2.55	0.41
3:C:254:ASN:HA	3:C:311:TYR:CD2	2.56	0.41
1:A:410:GLN:O	1:A:414:MET:HG3	2.21	0.41
1:A:507:TRP:O	1:A:508:SER:HB3	2.21	0.41
3:C:201:ARG:C	3:C:201:ARG:HD3	2.46	0.41
2:B:972:MET:HE3	2:B:975:PHE:HB3	2.03	0.40
1:A:378:MET:CE	3:C:284:THR:H	2.34	0.40
1:A:379:PHE:CE1	3:C:281:TRP:HB3	2.56	0.40
3:C:54:SER:O	3:C:58:LYS:HG3	2.22	0.40
3:C:308:ILE:C	3:C:310:ASP:N	2.78	0.40
1:A:530:LYS:HA	1:A:562:PHE:HE1	1.85	0.40
2:B:937:ILE:O	2:B:938:LYS:HB3	2.21	0.40
3:C:119:LEU:HD21	3:C:128:VAL:HG21	2.03	0.40
3:C:121:ASN:HD22	3:C:124:ASN:HD22	1.68	0.40
3:C:353:GLY:HA2	3:C:357:HIS:NE2	2.36	0.40
2:B:889:PHE:O	2:B:998:VAL:HA	2.21	0.40
3:C:280:ARG:HA	3:C:283:ARG:NE	2.36	0.40
2:B:895:PHE:O	2:B:896:LYS:C	2.64	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	187/225 (83%)	171 (91%)	15 (8%)	1 (0%)	24	54
2	B	185/212 (87%)	171 (92%)	13 (7%)	1 (0%)	24	54
3	C	326/394 (83%)	292 (90%)	29 (9%)	5 (2%)	8	28
All	All	698/831 (84%)	634 (91%)	57 (8%)	7 (1%)	12	39

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	1013	GLN
3	C	138	PRO
3	C	321	PRO
3	C	175	ALA
1	A	430	ASN
3	C	48	ALA
3	C	309	GLU

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	158/189 (84%)	150 (95%)	8 (5%)	21	53
2	B	162/184 (88%)	154 (95%)	8 (5%)	22	54
3	C	297/351 (85%)	285 (96%)	12 (4%)	28	62
All	All	617/724 (85%)	589 (96%)	28 (4%)	24	58

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	388	ASP
1	A	403	LEU
1	A	470	ILE
1	A	499	LEU
1	A	526	ILE
1	A	528	ILE
1	A	530	LYS
1	A	541	GLU
2	B	879	HIS
2	B	894	ASP
2	B	940	ILE
2	B	965	GLN
2	B	992	ASN
2	B	1013	GLN
2	B	1044	GLU
2	B	1062	ILE
3	C	56	ILE
3	C	61	ARG
3	C	98	ASN
3	C	139	ASP
3	C	181	LYS
3	C	182	ILE
3	C	188	ASP
3	C	221	MET
3	C	229	ASP
3	C	235	ILE
3	C	262	GLN
3	C	370	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	380	HIS
1	A	385	GLN
1	A	387	HIS
1	A	431	HIS
1	A	455	HIS
1	A	493	HIS
1	A	515	ASN
2	B	968	HIS
2	B	989	HIS

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Mol	Chain	Res	Type
2	B	1001	ASN
2	B	1013	GLN
2	B	1016	GLN
2	B	1025	ASN
2	B	1050	GLN
2	B	1076	ASN
3	C	41	HIS
3	C	64	HIS
3	C	97	ASN
3	C	121	ASN
3	C	176	GLN
3	C	218	ASN
3	C	220	HIS
3	C	371	ASN
3	C	377	ASN
3	C	384	GLN
3	C	387	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	FOK	A	101	-	29,31,31	3.49	10 (34%)	38,54,54	1.59	7 (18%)
6	ONM	A	100	4	44,45,45	3.80	18 (40%)	65,69,69	2.76	24 (36%)
8	GSP	C	395	4	33,34,34	1.85	3 (9%)	47,54,54	1.55	7 (14%)

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
5	FOK	A	101	-	-	5/7/80/80	0/3/3/3
6	ONM	A	100	4	-	5/32/48/48	0/4/4/4
8	GSP	C	395	4	-	2/21/38/38	0/3/3/3

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	101	FOK	O7-C11	14.28	1.43	1.21
6	A	100	ONM	CA4-CA5	12.65	1.60	1.38
6	A	100	ONM	CA3-CA2	10.74	1.57	1.38
6	A	100	ONM	OA-CA	8.80	1.47	1.22
6	A	100	ONM	CA5-CA6	8.18	1.53	1.39
8	C	395	GSP	PB-O3B	7.84	1.68	1.59
5	A	101	FOK	C10-C5	6.53	1.67	1.56
5	A	101	FOK	C4-C5	-5.96	1.48	1.56
6	A	100	ONM	PB-O1B	-5.84	1.53	1.59
6	A	100	ONM	PB-O1A	5.60	1.65	1.59
6	A	100	ONM	PA-O1A	4.38	1.64	1.59
6	A	100	ONM	PB-O3B	4.12	1.65	1.50
6	A	100	ONM	PA-O3A	3.91	1.64	1.50
8	C	395	GSP	PB-O3A	3.76	1.63	1.59
6	A	100	ONM	PG-O2G	3.24	1.60	1.50
5	A	101	FOK	C9-C11	3.18	1.58	1.54
6	A	100	ONM	CA1-CA	2.91	1.56	1.50
6	A	100	ONM	CA3-CA4	2.87	1.44	1.38
5	A	101	FOK	C17-C8	2.57	1.56	1.51
6	A	100	ONM	CA7-NA1	2.53	1.49	1.45
6	A	100	ONM	O3'-C3'	2.50	1.48	1.44
6	A	100	ONM	CA2-CA1	2.45	1.43	1.39
5	A	101	FOK	C10-C1	2.42	1.59	1.55
5	A	101	FOK	C20-C10	2.40	1.58	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	101	FOK	C3-C4	2.38	1.58	1.54
6	A	100	ONM	C5-N7	-2.32	1.34	1.39
6	A	100	ONM	C6-N1	2.31	1.43	1.38
5	A	101	FOK	O1-C13	2.30	1.49	1.45
5	A	101	FOK	C9-C10	2.20	1.63	1.57
8	C	395	GSP	C2'-C3'	2.19	1.59	1.53
6	A	100	ONM	C4-N3	2.15	1.39	1.34

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	100	ONM	O1G-PG-O3G	-8.75	74.98	107.80
6	A	100	ONM	C2-N3-C4	5.99	122.63	112.30
6	A	100	ONM	O2A-PA-O1A	5.92	123.27	107.27
6	A	100	ONM	O3G-PG-O1B	5.51	123.12	104.64
6	A	100	ONM	C3'-O3'-CA	5.25	125.63	117.24
6	A	100	ONM	CA3-CA4-CA5	-5.14	113.90	120.24
6	A	100	ONM	O1G-PG-O1B	-5.03	87.78	104.64
6	A	100	ONM	O3'-CA-CA1	4.77	119.29	111.71
8	C	395	GSP	O2B-PB-O3B	4.32	118.94	107.27
6	A	100	ONM	C5-C4-N3	-4.21	121.69	128.39
6	A	100	ONM	O2B-PB-O1B	4.02	118.15	107.27
8	C	395	GSP	O6-C6-N1	3.92	127.49	120.11
5	A	101	FOK	C10-C5-C4	-3.84	113.01	116.32
6	A	100	ONM	C5'-C4'-C3'	3.82	127.09	114.38
5	A	101	FOK	C20-C10-C1	-3.66	101.21	107.40
6	A	100	ONM	CA7-NA1-CA6	3.61	128.06	122.37
5	A	101	FOK	C2-C3-C4	3.58	118.80	113.40
8	C	395	GSP	O3B-PB-O1B	-3.51	100.14	110.70
5	A	101	FOK	C13-O1-C8	3.45	125.27	119.86
6	A	100	ONM	O5'-C5'-C4'	3.41	120.61	108.99
6	A	100	ONM	C6-C5-N7	3.17	136.06	130.29
6	A	100	ONM	C2-N1-C6	-3.14	119.42	125.11
8	C	395	GSP	PB-O3B-PG	-3.01	122.14	133.17
6	A	100	ONM	C5-C6-N1	2.96	120.78	113.25
5	A	101	FOK	C17-C8-C9	2.85	118.34	115.19
6	A	100	ONM	O1B-PB-O3B	-2.71	102.56	110.70
6	A	100	ONM	O2B-PB-O3B	-2.60	100.36	112.44
5	A	101	FOK	C19-C4-C18	-2.48	104.26	107.87
6	A	100	ONM	N1-C2-N3	-2.47	118.80	123.32
8	C	395	GSP	C1'-N9-C8	2.37	133.45	126.73
6	A	100	ONM	N9-C4-N3	2.37	130.68	125.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	100	ONM	O3'-CA-OA	-2.35	119.73	123.55
6	A	100	ONM	O5'-PA-O3A	-2.24	100.08	108.94
8	C	395	GSP	O2A-PA-O1A	2.19	122.64	112.44
6	A	100	ONM	C4-C5-N7	-2.15	107.26	110.67
8	C	395	GSP	O3G-PG-O3B	2.14	111.79	104.64
6	A	100	ONM	CA4-CA5-CA6	2.09	122.94	118.69
5	A	101	FOK	C19-C4-C5	2.05	119.06	111.92

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	101	FOK	C16-C13-C14-C15
6	A	100	ONM	O4'-C4'-C5'-O5'
5	A	101	FOK	C22-C21-O4-C7
5	A	101	FOK	O5-C21-O4-C7
6	A	100	ONM	C3'-C4'-C5'-O5'
5	A	101	FOK	O1-C13-C14-C15
8	C	395	GSP	PA-O3A-PB-O1B
6	A	100	ONM	PB-O1B-PG-O2G
6	A	100	ONM	PB-O1B-PG-O1G
8	C	395	GSP	PA-O3A-PB-O2B
5	A	101	FOK	C12-C13-C14-C15
6	A	100	ONM	PA-O1A-PB-O2B

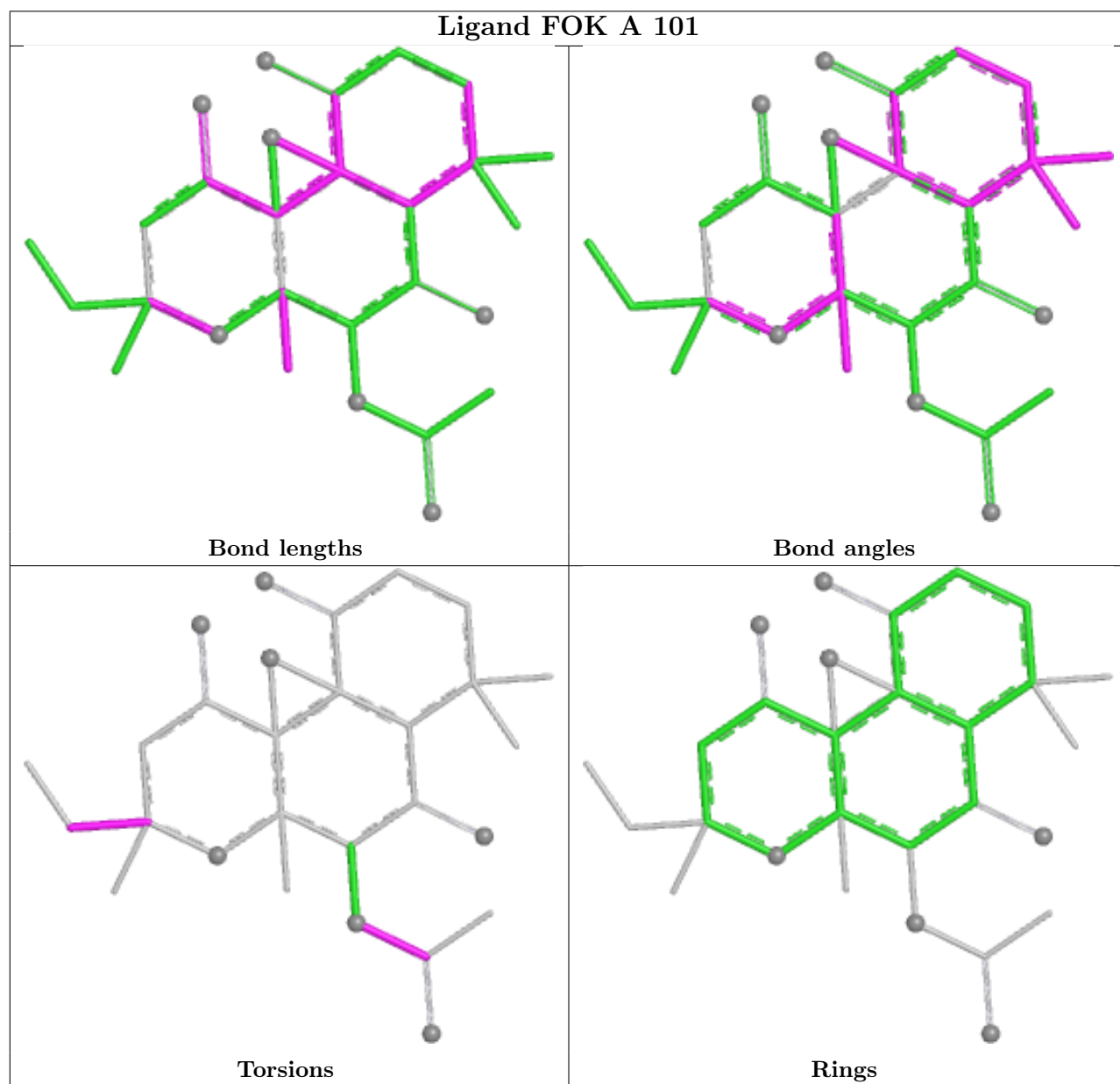
There are no ring outliers.

2 monomers are involved in 4 short contacts:

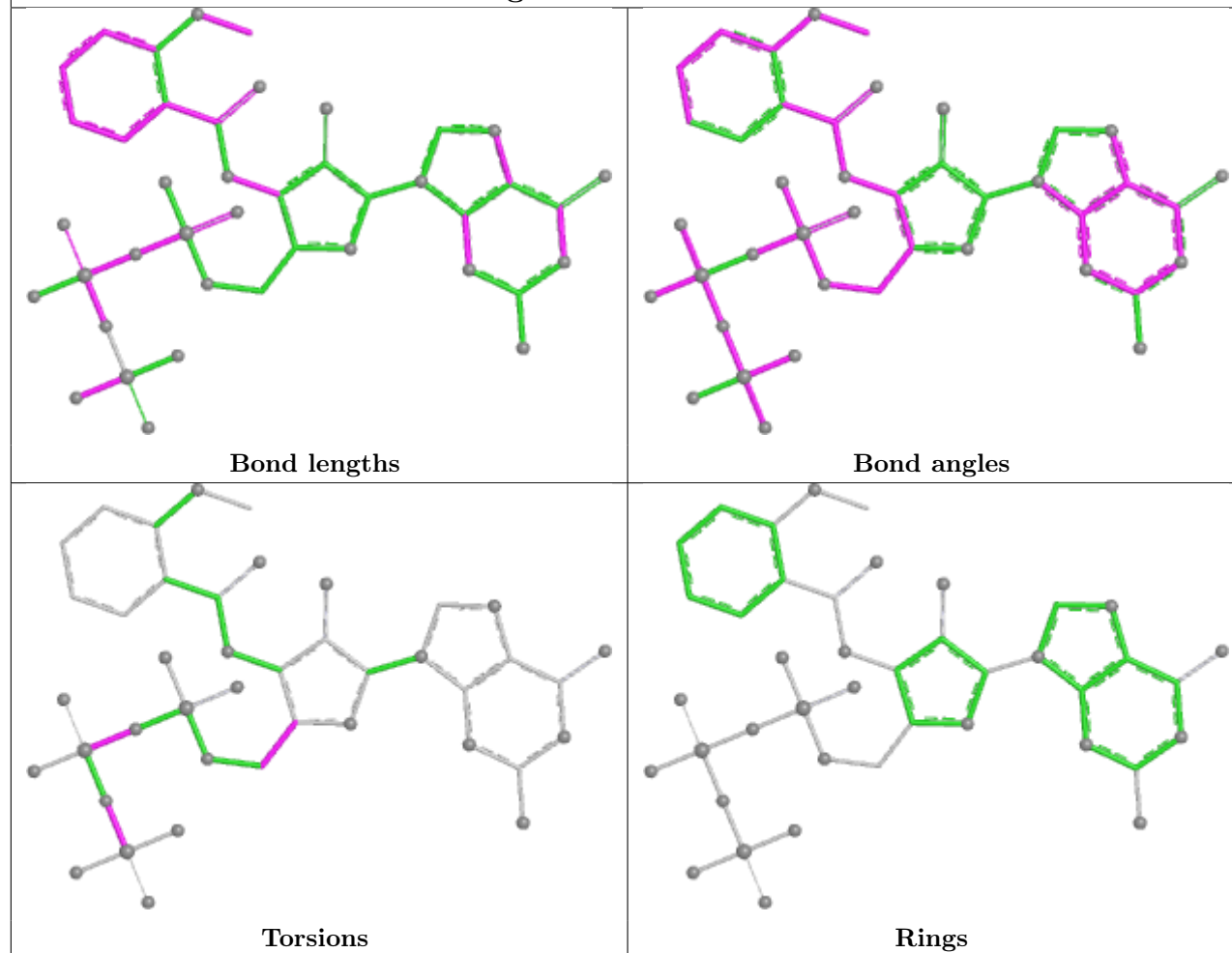
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	101	FOK	2	0
6	A	100	ONM	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.

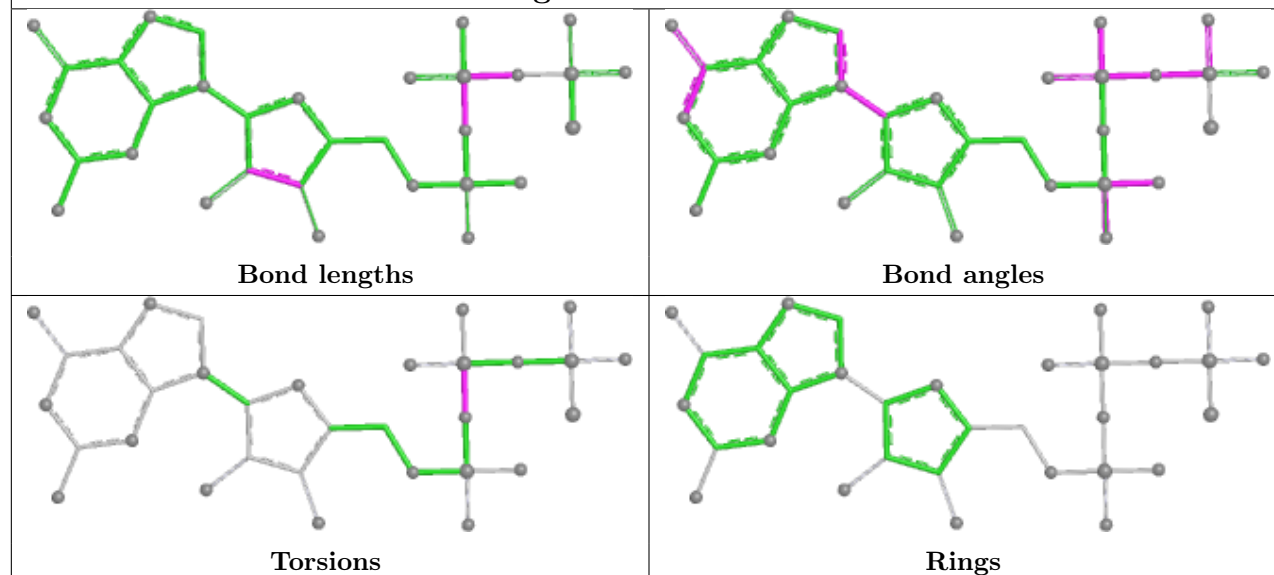
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



Ligand ONM A 100



Ligand GSP C 395



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	189/225 (84%)	0.81	10 (5%) 32 25	27, 64, 97, 116	0
2	B	189/212 (89%)	0.96	16 (8%) 16 14	20, 46, 80, 103	0
3	C	330/394 (83%)	0.25	3 (0%) 81 75	26, 49, 75, 103	0
All	All	708/831 (85%)	0.59	29 (4%) 41 33	20, 52, 84, 116	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	498	GLY	3.3
2	B	1046	SER	2.8
2	B	1031	ASP	2.8
2	B	1055	THR	2.8
1	A	456	ALA	2.6
2	B	1003	GLY	2.6
3	C	135	MET	2.6
2	B	1056	CYS	2.6
1	A	480	ASN	2.5
1	A	430	ASN	2.5
2	B	909	LEU	2.5
3	C	39	ALA	2.5
2	B	886	CYS	2.5
2	B	966	TYR	2.4
1	A	526	ILE	2.4
2	B	1048	ILE	2.3
1	A	400	PHE	2.3
2	B	984	ASP	2.3
1	A	407	CYS	2.3
2	B	1012	ALA	2.3
2	B	986	ILE	2.3
1	A	520	GLY	2.3
2	B	1034	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
3	C	353	GLY	2.2
2	B	1059	ARG	2.2
1	A	479	VAL	2.2
2	B	885	VAL	2.1
1	A	439	GLY	2.1
2	B	931	PHE	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](#)

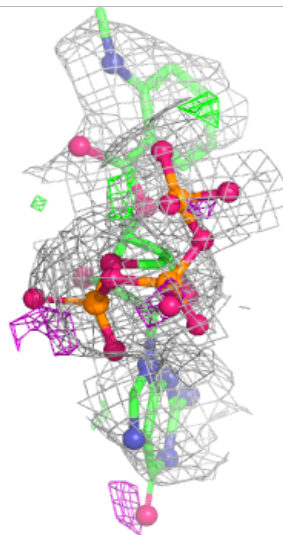
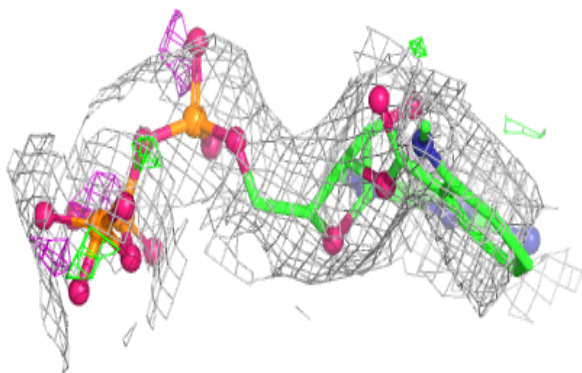
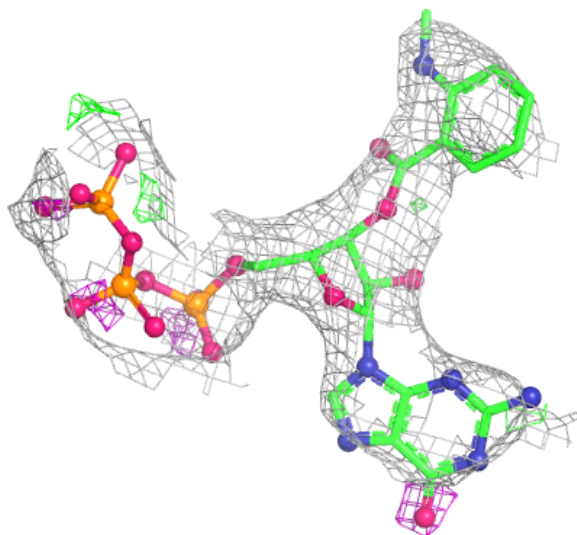
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MG	A	581	1/1	0.75	0.10	38,38,38,38	0
6	ONM	A	100	42/42	0.80	0.14	54,74,82,89	0
7	CL	C	397	1/1	0.87	0.10	49,49,49,49	0
4	MG	A	582	1/1	0.88	0.07	38,38,38,38	0
5	FOK	A	101	29/29	0.89	0.13	30,36,44,44	0
8	GSP	C	395	32/32	0.90	0.12	23,38,66,67	0
4	MG	C	396	1/1	0.93	0.08	21,21,21,21	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.

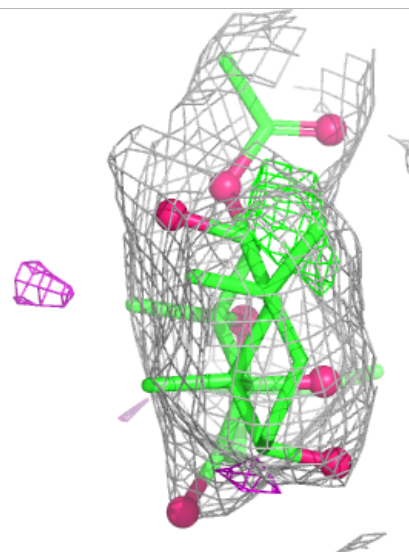
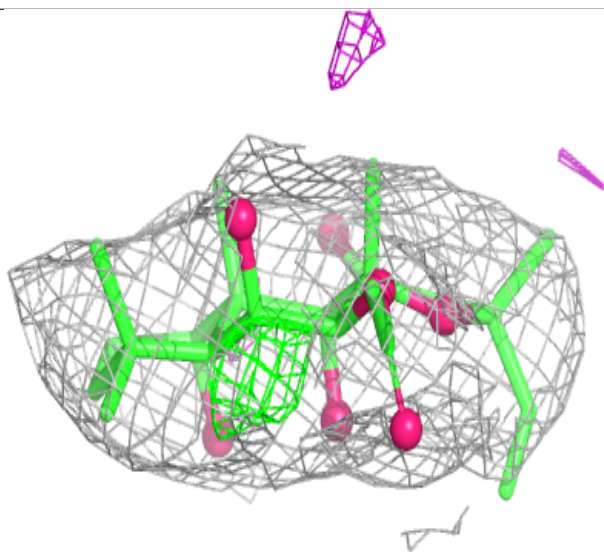
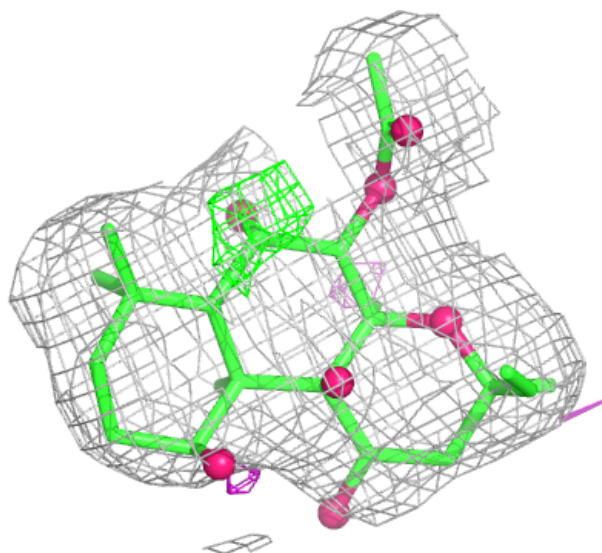
Electron density around ONM A 100:

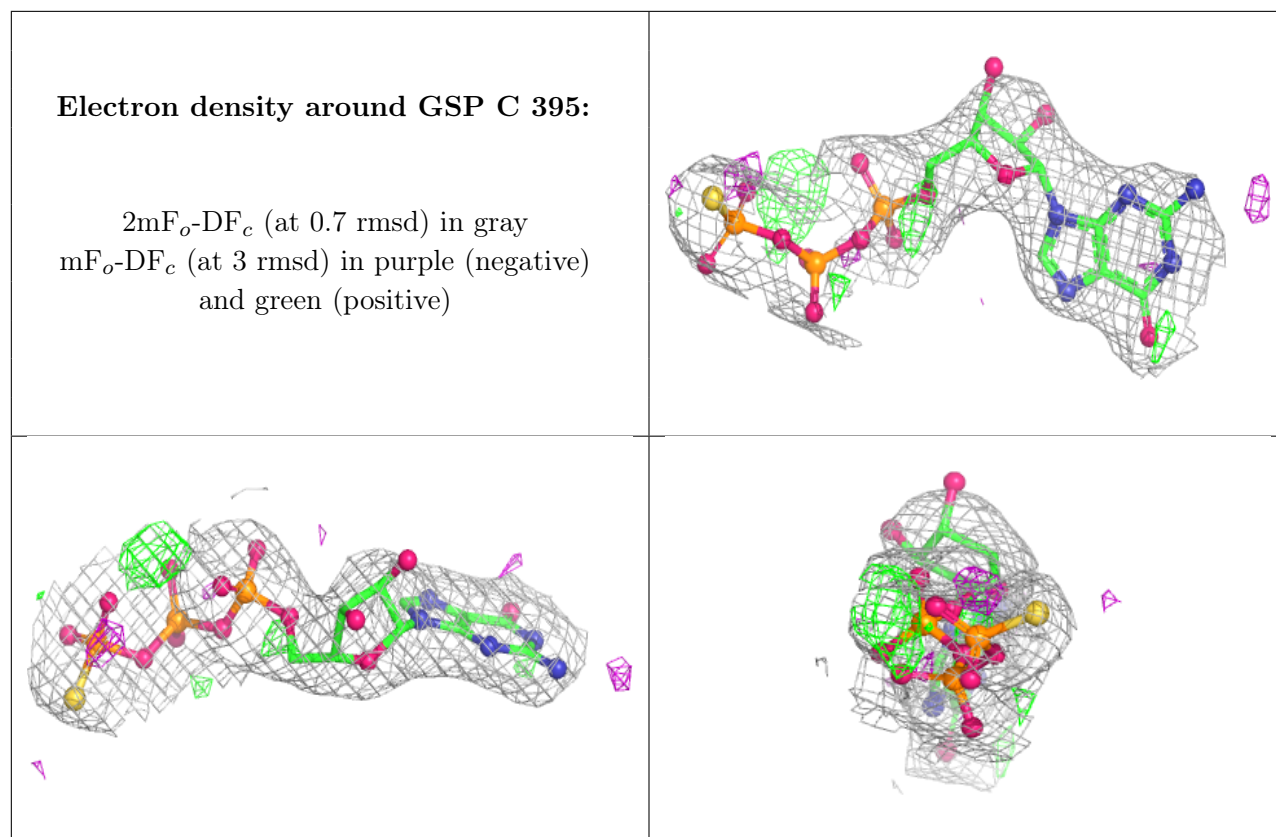
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FOK A 101:

$2mF_o-DF_c$ (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.