



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2026 – 06:41 PM EDT

PDB ID : 3GCM / pdb_00003gcm
Title : Crystal Structure of E. coli polynucleotide phosphorylase bound to RNA and RNase E
Authors : Nurmohamed, S.; Luisi, B.L.
Deposited on : 2009-02-22
Resolution : 2.50 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

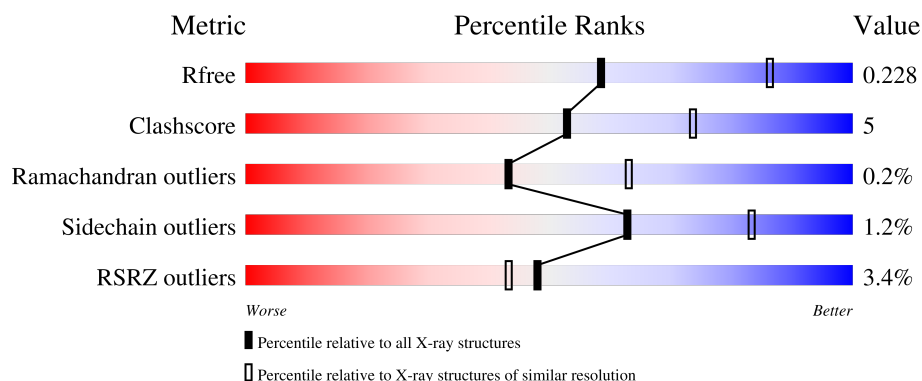
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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.



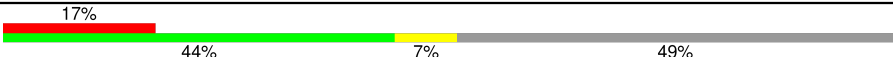
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	549	3% 89% 9% ..
1	B	549	2% 90% 8% ..
1	C	549	3% 89% 10% .
2	D	41	10% 39% 12% 49%
2	E	41	12% 46% . 51%

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Mol	Chain	Length	Quality of chain
2	F	41	

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
3	FLC	B	552	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13916 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 Polyribonucleotide nucleotidyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	544	Total	C	N	O	S	8	4	0
			4151	2604	724	802	21			
1	B	544	Total	C	N	O	S	0	5	0
			4170	2614	732	803	21			
1	C	544	Total	C	N	O	S	0	2	0
			4142	2598	721	802	21			

- Molecule 2 is a protein called Ribonuclease E.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	21	Total	C	N	O	0	0	0
			134	80	30	24			
2	E	20	Total	C	N	O	0	0	0
			130	78	29	23			
2	F	21	Total	C	N	O	0	0	0
			138	82	31	25			

- Molecule 3 is CITRATE ANION (CCD ID: FLC) (formula: $C_6H_5O_7$).

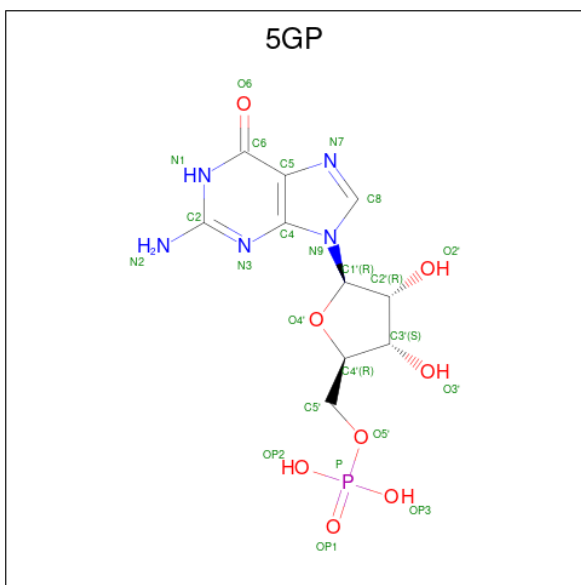


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		
3	A	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		
3	C	1	Total	C	O	0	0
			13	6	7		
3	C	1	Total	C	O	0	0
			13	6	7		

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

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

- Molecule 5 is GUANOSINE-5'-MONOPHOSPHATE (CCD ID: 5GP) (formula: C₁₀H₁₄N₅O₈P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			20	10	5	5		
5	B	1	Total	C	N	O	0	0
			20	10	5	5		
5	C	1	Total	C	N	O	0	0
			20	10	5	5		

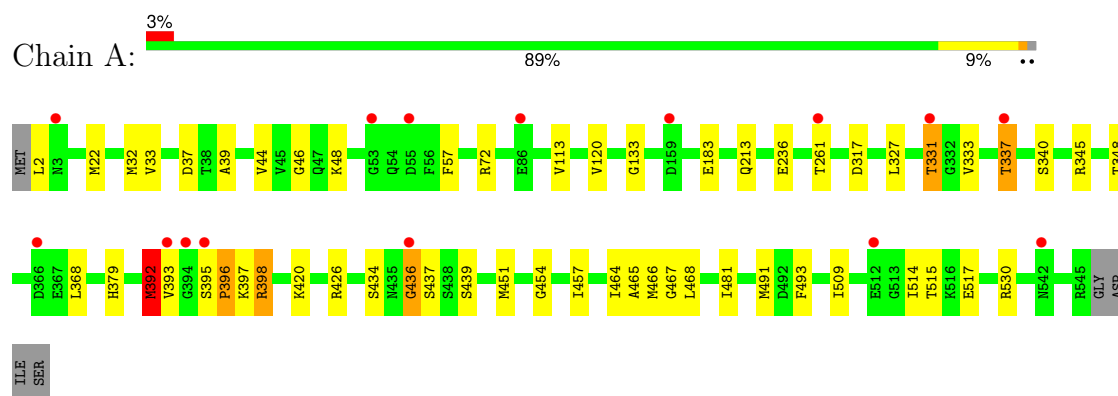
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	334	Total	O	0	0
			334	334		
6	D	8	Total	O	0	0
			8	8		
6	B	280	Total	O	0	0
			280	280		
6	E	9	Total	O	0	0
			9	9		
6	C	247	Total	O	0	0
			247	247		
6	F	6	Total	O	0	0
			6	6		

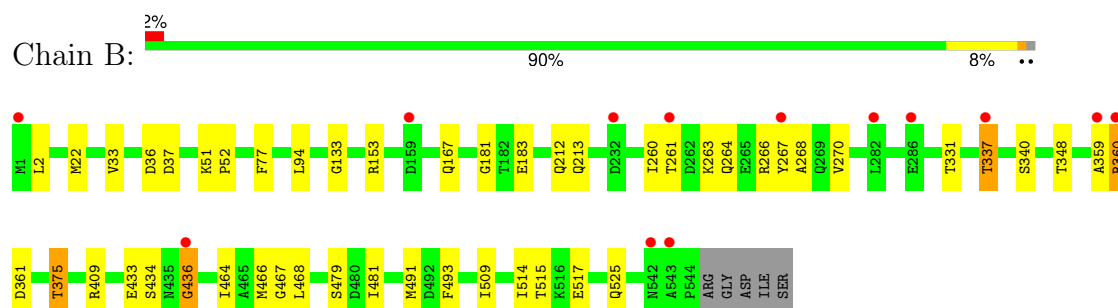
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

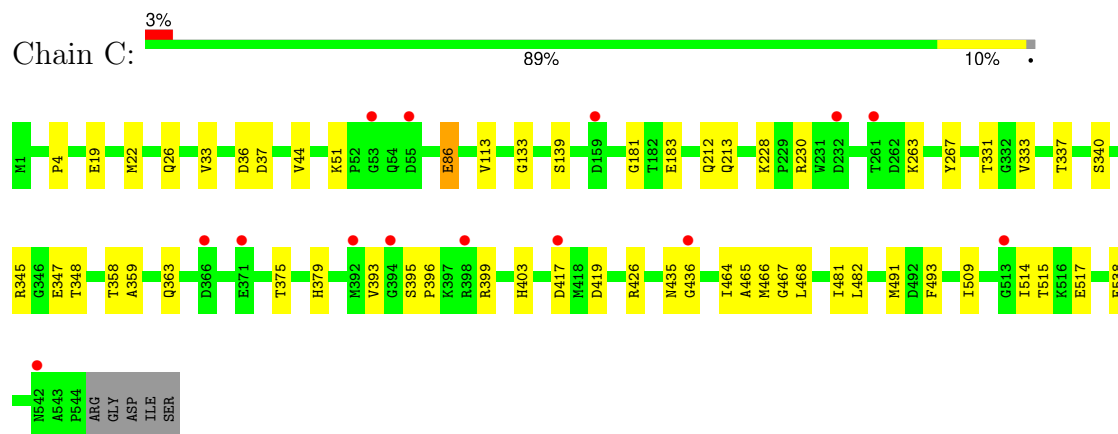
- Molecule 1: Polyribonucleotide nucleotidyltransferase



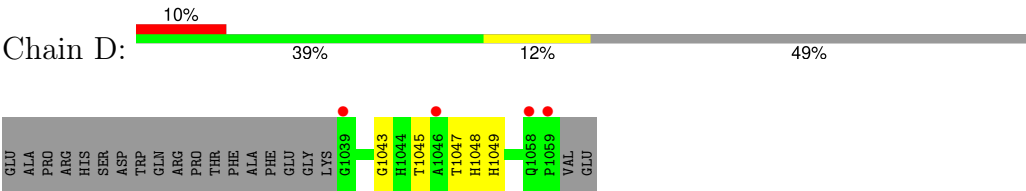
- Molecule 1: Polyribonucleotide nucleotidyltransferase



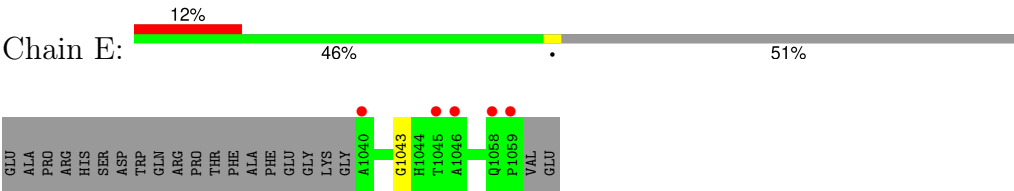
- Molecule 1: Polyribonucleotide nucleotidyltransferase



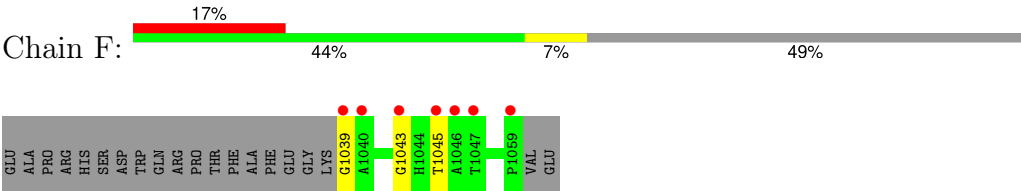
● Molecule 2: Ribonuclease E



● Molecule 2: Ribonuclease E



● Molecule 2: Ribonuclease E



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	176.34Å 176.34Å 189.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.4 (20.00-2.50) 99.2 (20.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.166 , 0.221 0.201 , 0.228	Depositor DCC
R_{free} test set	5134 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 33.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13916	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.42% 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: FLC, MG, 5GP

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	0.76	7/4226 (0.2%)	0.80	3/5728 (0.1%)
1	B	0.48	1/4245 (0.0%)	0.78	3/5751 (0.1%)
1	C	0.49	4/4211 (0.1%)	0.77	2/5708 (0.0%)
2	D	0.51	0/139	0.81	0/191
2	E	0.47	0/135	0.78	0/186
2	F	0.51	0/143	0.77	0/196
All	All	0.59	12/13099 (0.1%)	0.78	8/17760 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	1
1	C	0	1
All	All	0	5

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	392[A]	MET	CA-C	-20.48	1.25	1.52
1	A	392[B]	MET	CA-C	-20.48	1.25	1.52
1	A	398[A]	ARG	N-CA	-14.76	1.27	1.46
1	A	398[B]	ARG	N-CA	-14.76	1.27	1.46
1	A	454	GLY	C-O	-5.99	1.16	1.24
1	A	439	SER	C-O	-5.90	1.16	1.24
1	C	345	ARG	C-O	-5.57	1.17	1.23
1	A	57	PHE	C-O	-5.37	1.17	1.24
1	C	348	THR	C-O	-5.37	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	347	GLU	C-O	-5.25	1.17	1.24
1	B	433	GLU	C-O	-5.10	1.17	1.23
1	C	86	GLU	C-O	-5.07	1.18	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	392[A]	MET	CA-C-O	-7.24	110.16	120.51
1	A	392[B]	MET	CA-C-O	-7.24	110.16	120.51
1	B	434	SER	N-CA-C	6.18	119.20	108.76
1	B	51	LYS	CA-C-N	5.31	126.48	119.84
1	B	51	LYS	C-N-CA	5.31	126.48	119.84
1	C	51	LYS	CA-C-N	5.29	126.45	119.84
1	C	51	LYS	C-N-CA	5.29	126.45	119.84
1	A	120	VAL	N-CA-C	5.12	115.42	107.78

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	392[A]	MET	Mainchain
1	A	392[B]	MET	Mainchain
1	A	436	GLY	Peptide
1	B	436	GLY	Peptide
1	C	436	GLY	Peptide

5.2 Too-close contacts [i](#)

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	4151	0	4170	39	0
1	B	4170	0	4203	38	0
1	C	4142	0	4160	47	0
2	D	134	0	119	4	0
2	E	130	0	116	3	0
2	F	138	0	125	6	0
3	A	26	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	52	0	20	6	0
3	C	26	0	10	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	B	40	0	24	1	0
5	C	20	0	12	0	0
6	A	334	0	0	12	0
6	B	280	0	0	10	0
6	C	247	0	0	11	0
6	D	8	0	0	1	0
6	E	9	0	0	0	0
6	F	6	0	0	0	0
All	All	13916	0	12969	129	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 5.

All (129) 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:B:264:GLN:NE2	1:C:26:GLN:HA	1.74	1.02
1:B:77:PHE:HB2	5:B:550:5GP:H2'	1.50	0.92
2:E:1043:GLY:HA3	1:C:22:MET:HE1	1.51	0.91
1:B:264:GLN:HE21	1:C:26:GLN:HA	1.37	0.85
1:B:515:THR:CG2	6:B:608:HOH:O	2.30	0.80
1:C:337:THR:HG21	1:C:340:SER:HB3	1.64	0.79
1:A:236:GLU:HG3	6:A:646:HOH:O	1.84	0.78
1:A:22:MET:HE1	2:F:1043:GLY:HA3	1.66	0.78
3:B:552:FLC:HG1	6:B:1067:HOH:O	1.82	0.78
1:A:337:THR:HG21	1:A:340:SER:HB3	1.68	0.76
1:A:72[A]:ARG:HD3	6:A:689:HOH:O	1.87	0.74
1:B:337:THR:HG21	1:B:340:SER:HB3	1.71	0.73
3:B:552:FLC:CG	6:B:1067:HOH:O	2.37	0.70
6:C:1044:HOH:O	2:F:1045:THR:HG21	1.91	0.70
1:C:393:VAL:HG12	1:C:393:VAL:O	1.91	0.69
2:E:1043:GLY:CA	1:C:22:MET:HE1	2.23	0.68
6:C:1044:HOH:O	2:F:1045:THR:CG2	2.41	0.67
1:A:2:LEU:HG	1:A:22:MET:HG3	1.75	0.67
2:E:1043:GLY:HA3	1:C:22:MET:CE	2.22	0.67
1:C:358:THR:HG22	1:C:359:ALA:H	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:515:THR:HG23	6:B:608:HOH:O	1.92	0.66
1:C:538:GLU:HG3	6:C:627:HOH:O	1.95	0.65
1:B:263:LYS:O	1:B:267:TYR:HD2	1.79	0.65
1:B:479:SER:OG	1:B:525:GLN:NE2	2.29	0.65
1:A:395:SER:O	1:A:396:PRO:C	2.36	0.64
1:B:36:ASP:O	1:B:37:ASP:HB2	1.97	0.63
1:C:263:LYS:O	1:C:267:TYR:HD2	1.80	0.63
1:B:261:THR:HG21	6:B:702:HOH:O	1.99	0.62
1:B:466:MET:HB2	1:B:493:PHE:HB3	1.82	0.61
1:A:515:THR:HG22	1:A:517:GLU:H	1.66	0.60
1:A:331:THR:HG22	2:D:1048:HIS:HB2	1.84	0.60
1:B:153:ARG:HH21	3:B:552:FLC:CBC	2.14	0.59
1:C:435:ASN:HD21	1:C:482:LEU:HG	1.67	0.59
1:A:466:MET:HB2	1:A:493:PHE:HB3	1.84	0.59
1:B:515:THR:HG22	1:B:517:GLU:H	1.67	0.59
1:C:358:THR:HG22	1:C:359:ALA:N	2.17	0.58
1:A:379:HIS:HE1	6:A:602:HOH:O	1.87	0.57
1:B:33:VAL:HG21	1:B:133:GLY:HA2	1.85	0.57
1:B:375:THR:HG22	6:B:697:HOH:O	2.04	0.57
1:A:22:MET:HE1	2:F:1043:GLY:CA	2.34	0.56
1:C:139:SER:O	1:C:230:ARG:NH2	2.36	0.55
1:A:397:LYS:HE3	6:A:694:HOH:O	2.07	0.55
1:C:4:PRO:HB2	1:C:19:GLU:HG3	1.86	0.55
1:A:33:VAL:HG21	1:A:133:GLY:HA2	1.89	0.55
1:A:72[A]:ARG:HH22	1:A:368:LEU:HA	1.72	0.54
1:C:228:LYS:O	1:C:230:ARG:NH1	2.41	0.54
1:C:395:SER:HB2	1:C:396:PRO:HD2	1.88	0.54
1:B:94:LEU:HD11	3:B:552:FLC:HG2	1.90	0.53
1:A:327:LEU:CD2	1:A:345:ARG:HG3	2.38	0.53
1:A:420:LYS:O	1:A:420:LYS:HG3	2.08	0.53
1:C:515:THR:HG22	1:C:517:GLU:H	1.74	0.53
1:C:33:VAL:HG21	1:C:133:GLY:HA2	1.90	0.53
1:C:426:ARG:HD3	6:C:612:HOH:O	2.09	0.53
1:B:359:ALA:O	1:B:360:ARG:C	2.51	0.52
1:B:263:LYS:O	1:B:267:TYR:CD2	2.63	0.52
1:A:426:ARG:HD3	6:A:565:HOH:O	2.09	0.52
1:A:261:THR:HG22	1:A:317:ASP:HA	1.91	0.52
1:B:266:ARG:O	1:B:270:VAL:HG23	2.09	0.52
1:A:468:LEU:HB2	1:A:491:MET:HE2	1.92	0.51
1:B:264:GLN:HE22	1:C:26:GLN:HA	1.66	0.51
1:C:466:MET:HB2	1:C:493:PHE:HB3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:379:HIS:CD2	1:C:426:ARG:HE	2.30	0.50
1:A:32:MET:HE2	1:A:39:ALA:HB1	1.94	0.49
1:A:348:THR:OG1	1:A:436:GLY:O	2.25	0.49
1:B:181:GLY:H	1:B:212:GLN:HE22	1.60	0.49
1:B:183:GLU:HA	1:B:213:GLN:HE22	1.78	0.49
1:A:327:LEU:HD23	1:A:345:ARG:HG3	1.93	0.49
2:D:1049:HIS:CE1	6:D:1112:HOH:O	2.64	0.49
1:B:2:LEU:HD23	1:B:22:MET:HE3	1.94	0.49
1:C:375:THR:HG23	6:C:614:HOH:O	2.11	0.49
1:A:46:GLY:O	6:A:561:HOH:O	2.20	0.48
1:A:183:GLU:HA	1:A:213:GLN:HE22	1.79	0.48
1:B:509:ILE:HD13	1:B:514:ILE:HG21	1.94	0.48
1:A:437:SER:HB2	1:A:465:ALA:HB2	1.95	0.48
1:B:359:ALA:O	1:B:361:ASP:N	2.46	0.48
1:A:48:LYS:CD	6:A:653:HOH:O	2.62	0.47
1:A:379:HIS:CD2	1:A:426:ARG:HE	2.32	0.47
1:A:37:ASP:OD1	2:F:1039:GLY:HA3	2.15	0.47
1:A:48:LYS:HD2	6:A:653:HOH:O	2.14	0.47
1:A:509:ILE:HD13	1:A:514:ILE:HG21	1.96	0.47
1:A:72[A]:ARG:HD2	6:A:775:HOH:O	2.13	0.47
1:B:348:THR:OG1	1:B:436:GLY:O	2.29	0.47
1:B:267:TYR:O	1:B:270:VAL:N	2.48	0.47
1:C:465:ALA:HB3	6:C:573:HOH:O	2.15	0.47
1:C:379:HIS:HD2	1:C:426:ARG:HE	1.63	0.47
1:C:263:LYS:HG2	1:C:267:TYR:CE2	2.50	0.47
1:C:509:ILE:HD13	1:C:514:ILE:HG21	1.97	0.46
1:B:167:GLN:NE2	6:B:657:HOH:O	2.48	0.46
1:C:393:VAL:O	1:C:393:VAL:CG1	2.61	0.46
1:C:515:THR:HG22	6:C:1025:HOH:O	2.15	0.46
1:C:183:GLU:HA	1:C:213:GLN:HE22	1.80	0.46
1:A:530:ARG:HD3	6:A:607:HOH:O	2.16	0.46
1:A:379:HIS:HD2	1:A:426:ARG:HE	1.64	0.45
1:C:33:VAL:HG21	1:C:133:GLY:CA	2.46	0.45
1:B:468:LEU:HB2	1:B:491:MET:HE2	1.98	0.45
1:C:435:ASN:ND2	1:C:482:LEU:HG	2.31	0.45
1:C:181:GLY:H	1:C:212:GLN:HE22	1.64	0.45
6:A:685:HOH:O	2:D:1047:THR:HG21	2.17	0.45
1:C:395:SER:HB2	1:C:396:PRO:CD	2.47	0.45
1:C:467:GLY:HA3	1:C:481:ILE:HG21	1.98	0.45
1:B:261:THR:CG2	6:B:702:HOH:O	2.61	0.44
1:C:44:VAL:HG22	1:C:113:VAL:HG22	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:337:THR:HG22	6:B:596:HOH:O	2.16	0.44
6:C:1044:HOH:O	2:F:1045:THR:HG22	2.13	0.44
1:A:467:GLY:HA3	1:A:481:ILE:HG21	1.99	0.44
1:C:515:THR:CG2	6:C:1025:HOH:O	2.66	0.44
1:C:468:LEU:HB2	1:C:491:MET:HE2	1.99	0.44
1:C:263:LYS:CG	1:C:267:TYR:CE2	3.01	0.44
1:C:363:GLN:NE2	6:C:703:HOH:O	2.51	0.43
1:B:409:ARG:NH1	3:B:552:FLC:HA2	2.33	0.43
1:C:263:LYS:HG3	1:C:267:TYR:HE2	1.83	0.42
1:A:379:HIS:HD2	1:A:426:ARG:HH21	1.67	0.42
1:C:399[A]:ARG:HG2	1:C:403:HIS:CE1	2.54	0.42
1:C:417:ASP:OD2	1:C:419:ASP:HB2	2.20	0.42
1:A:44:VAL:HG22	1:A:113:VAL:HG22	2.01	0.42
1:A:22:MET:HE2	1:A:22:MET:HB3	1.90	0.42
1:B:359:ALA:C	1:B:361:ASP:N	2.77	0.41
1:C:515:THR:HG21	6:C:700:HOH:O	2.20	0.41
1:B:267:TYR:O	1:B:268:ALA:C	2.64	0.41
1:C:36:ASP:O	1:C:37:ASP:HB2	2.20	0.41
1:B:467:GLY:HA3	1:B:481:ILE:HG21	2.02	0.41
1:B:260:ILE:O	1:B:266:ARG:HD3	2.20	0.41
1:B:515:THR:HG22	1:B:517:GLU:N	2.34	0.41
1:C:358:THR:CG2	1:C:359:ALA:H	2.31	0.41
1:C:399[A]:ARG:HD3	1:C:403:HIS:CE1	2.55	0.41
3:B:552:FLC:HG2	6:B:1067:HOH:O	2.12	0.41
1:A:72[B]:ARG:CD	6:A:689:HOH:O	2.69	0.40
1:A:451:MET:HE3	1:A:457:ILE:HD11	2.03	0.40
2:D:1043:GLY:HA3	1:B:22:MET:SD	2.61	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	546/549 (100%)	532 (97%)	11 (2%)	3 (0%)	24	43
1	B	547/549 (100%)	530 (97%)	15 (3%)	2 (0%)	30	49
1	C	544/549 (99%)	532 (98%)	12 (2%)	0	100	100
2	D	19/41 (46%)	18 (95%)	1 (5%)	0	100	100
2	E	18/41 (44%)	18 (100%)	0	0	100	100
2	F	19/41 (46%)	19 (100%)	0	0	100	100
All	All	1693/1770 (96%)	1649 (97%)	39 (2%)	5 (0%)	43	55

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	360	ARG
1	A	392[A]	MET
1	A	392[B]	MET
1	A	396	PRO
1	B	52	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	441/446 (99%)	433 (98%)	8 (2%)	51	77
1	B	444/446 (100%)	440 (99%)	4 (1%)	70	87
1	C	440/446 (99%)	436 (99%)	4 (1%)	70	87
2	D	10/28 (36%)	9 (90%)	1 (10%)	7	15
2	E	10/28 (36%)	10 (100%)	0	100	100
2	F	11/28 (39%)	11 (100%)	0	100	100
All	All	1356/1422 (95%)	1339 (99%)	17 (1%)	63	82

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	331	THR
1	A	333	VAL
1	A	337	THR
1	A	393	VAL
1	A	398[A]	ARG
1	A	398[B]	ARG
1	A	434	SER
1	A	464	ILE
2	D	1045	THR
1	B	331	THR
1	B	337	THR
1	B	375	THR
1	B	464	ILE
1	C	86	GLU
1	C	331	THR
1	C	333	VAL
1	C	464	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	64	GLN
1	A	212	GLN
1	A	213	GLN
1	A	243	ASN
1	A	264	GLN
1	A	269	GLN
1	A	363	GLN
1	A	379	HIS
2	D	1049	HIS
1	B	3	ASN
1	B	47	GLN
1	B	62	ASN
1	B	64	GLN
1	B	167	GLN
1	B	212	GLN
1	B	213	GLN
1	B	234	GLN
1	B	264	GLN
1	B	363	GLN
1	B	525	GLN
2	E	1048	HIS

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Mol	Chain	Res	Type
1	C	3	ASN
1	C	47	GLN
1	C	62	ASN
1	C	212	GLN
1	C	213	GLN
1	C	363	GLN
1	C	379	HIS
2	F	1049	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 14 ligands modelled in this entry, 3 are monoatomic - leaving 11 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	$\# Z > 2$
3	FLC	B	555	-	12,12,12	1.05	0	17,17,17	1.24	1 (5%)
5	5GP	B	551	-	22,22,26	0.37	0	33,33,40	0.50	0
3	FLC	A	550	-	12,12,12	1.07	0	17,17,17	1.14	1 (5%)
3	FLC	A	551	-	12,12,12	1.06	0	17,17,17	1.20	1 (5%)
3	FLC	C	551	-	12,12,12	1.06	0	17,17,17	1.18	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	5GP	B	550	-	22,22,26	0.34	0	33,33,40	0.48	0
5	5GP	C	550	-	22,22,26	0.36	0	33,33,40	0.29	0
3	FLC	B	553	-	12,12,12	1.04	0	17,17,17	1.18	1 (5%)
3	FLC	B	554	-	12,12,12	1.08	0	17,17,17	1.19	1 (5%)
3	FLC	C	552	-	12,12,12	1.08	0	17,17,17	1.23	1 (5%)
3	FLC	B	552	-	12,12,12	1.08	0	17,17,17	1.31	1 (5%)

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
3	FLC	B	555	-	-	2/16/16/16	-
5	5GP	B	551	-	-	2/6/22/26	0/3/3/3
3	FLC	A	550	-	-	0/16/16/16	-
3	FLC	A	551	-	-	3/16/16/16	-
3	FLC	C	551	-	-	0/16/16/16	-
5	5GP	B	550	-	-	0/6/22/26	0/3/3/3
5	5GP	C	550	-	-	0/6/22/26	0/3/3/3
3	FLC	B	553	-	-	2/16/16/16	-
3	FLC	B	554	-	-	5/16/16/16	-
3	FLC	C	552	-	-	0/16/16/16	-
3	FLC	B	552	-	-	2/16/16/16	-

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	554	FLC	OB2-CBC-CB	3.20	119.28	113.14
3	C	551	FLC	OB2-CBC-CB	3.17	119.23	113.14
3	B	553	FLC	OB2-CBC-CB	3.17	119.22	113.14
3	B	555	FLC	OB2-CBC-CB	3.17	119.21	113.14
3	B	552	FLC	OB2-CBC-CB	3.15	119.17	113.14
3	A	551	FLC	OB2-CBC-CB	3.14	119.16	113.14
3	C	552	FLC	OB2-CBC-CB	3.11	119.10	113.14
3	A	550	FLC	OB2-CBC-CB	2.94	118.78	113.14

There are no chirality outliers.

All (16) torsion outliers are listed below:

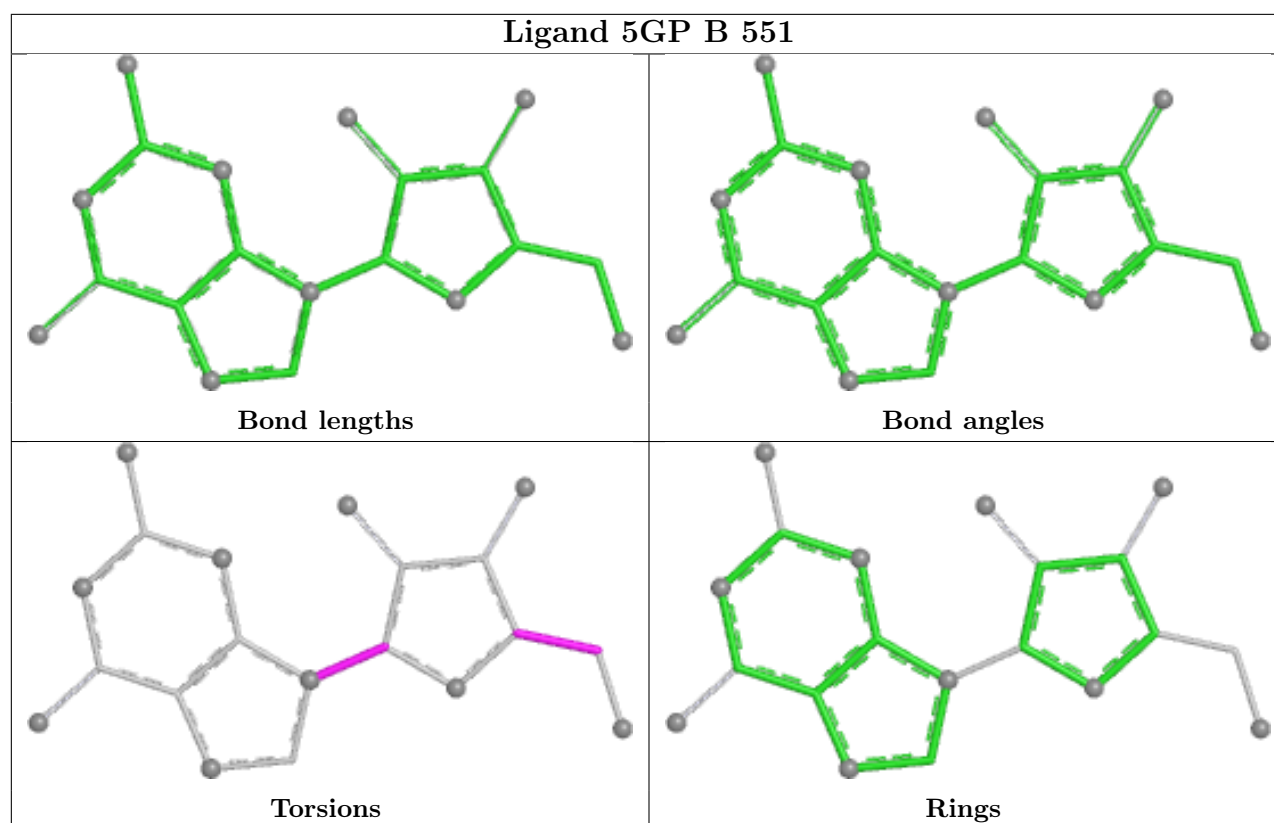
Mol	Chain	Res	Type	Atoms
3	B	554	FLC	CAC-CA-CB-OHB
3	B	554	FLC	CAC-CA-CB-CBC
3	B	554	FLC	CAC-CA-CB-CG
3	B	552	FLC	CA-CB-CG-CGC
3	B	553	FLC	CBC-CB-CG-CGC
3	B	554	FLC	CBC-CB-CG-CGC
3	B	553	FLC	OHB-CB-CG-CGC
3	B	554	FLC	OHB-CB-CG-CGC
3	A	551	FLC	CB-CA-CAC-OA1
5	B	551	5GP	O4'-C4'-C5'-O5'
3	B	552	FLC	CBC-CB-CG-CGC
5	B	551	5GP	C2'-C1'-N9-C4
3	A	551	FLC	CB-CA-CAC-OA2
3	B	555	FLC	CB-CA-CAC-OA1
3	A	551	FLC	OHB-CB-CG-CGC
3	B	555	FLC	CB-CA-CAC-OA2

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	550	5GP	1	0
3	B	552	FLC	6	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	544/549 (99%)	-0.07	15 (2%)	55 50	10, 21, 44, 62	5 (0%)
1	B	544/549 (99%)	0.00	13 (2%)	59 55	11, 23, 46, 68	6 (1%)
1	C	544/549 (99%)	0.05	14 (2%)	57 52	12, 25, 47, 64	3 (0%)
2	D	21/41 (51%)	0.93	4 (19%)	3 2	17, 37, 47, 52	0
2	E	20/41 (48%)	1.30	5 (25%)	2 1	33, 50, 62, 65	0
2	F	21/41 (51%)	1.47	7 (33%)	1 1	24, 48, 61, 61	0
All	All	1694/1770 (95%)	0.04	58 (3%)	48 43	10, 23, 49, 68	14 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	159	ASP	7.5
1	C	159	ASP	7.3
1	A	159	ASP	6.3
1	C	436	GLY	4.6
2	F	1059	PRO	4.4
2	E	1059	PRO	4.3
2	F	1039	GLY	4.2
1	B	267	TYR	4.1
1	A	395	SER	4.0
2	D	1059	PRO	3.8
2	E	1058	GLN	3.4
1	A	394	GLY	3.4
1	A	261	THR	3.3
1	B	360	ARG	3.3
1	B	542	ASN	3.3
1	C	55	ASP	3.1
2	E	1045	THR	3.0
1	C	394	GLY	2.9
1	C	398	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
2	F	1046	ALA	2.9
1	B	1	MET	2.9
2	F	1045	THR	2.8
2	E	1040	ALA	2.8
1	B	232	ASP	2.7
1	A	53	GLY	2.7
1	B	359	ALA	2.7
1	A	86	GLU	2.6
1	C	261	THR	2.6
2	D	1058	GLN	2.6
1	A	436	GLY	2.6
1	C	366	ASP	2.6
1	C	53	GLY	2.5
1	A	55	ASP	2.4
1	B	282	LEU	2.4
1	B	261	THR	2.4
1	A	542	ASN	2.4
2	F	1043	GLY	2.3
1	A	3	ASN	2.3
2	F	1040	ALA	2.3
1	A	393	VAL	2.3
1	C	232	ASP	2.2
1	B	543	ALA	2.2
2	D	1046	ALA	2.2
1	A	512	GLU	2.2
1	B	286	GLU	2.2
1	C	392	MET	2.1
1	A	366	ASP	2.1
2	E	1046	ALA	2.1
1	C	542	ASN	2.1
1	A	331	THR	2.1
1	C	417	ASP	2.1
1	B	436	GLY	2.1
2	D	1039	GLY	2.1
1	B	337	THR	2.1
1	C	371	GLU	2.0
1	C	513	GLY	2.0
1	A	337	THR	2.0
2	F	1047	THR	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](#)

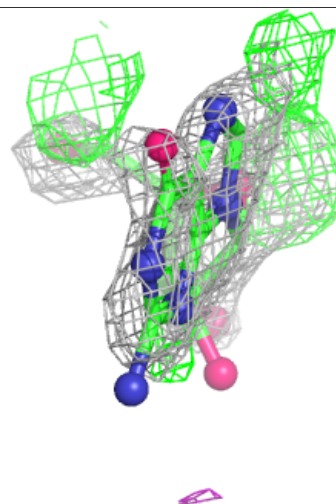
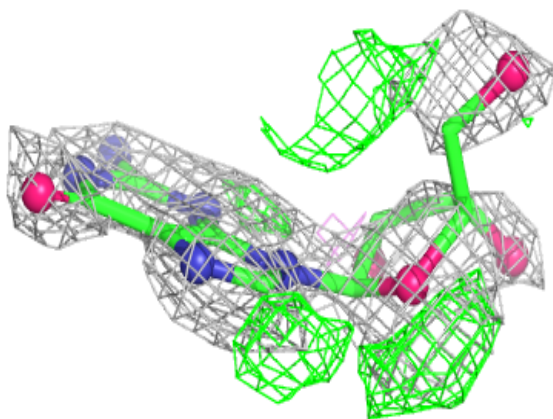
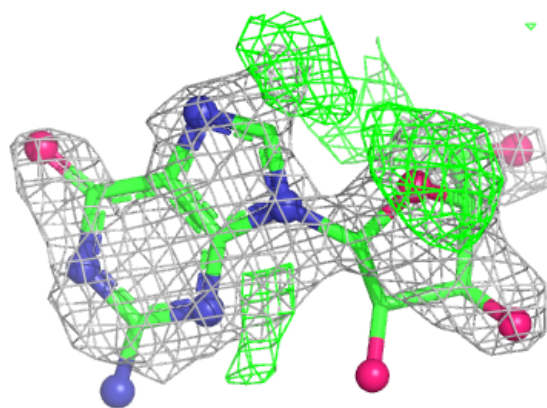
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	FLC	B	554	13/13	0.63	0.21	77,80,81,81	0
5	5GP	B	550	20/24	0.68	0.31	18,20,23,25	20
5	5GP	C	550	20/24	0.73	0.34	23,24,29,31	20
3	FLC	B	553	13/13	0.78	0.13	41,48,52,54	0
5	5GP	B	551	20/24	0.79	0.23	15,16,18,21	20
4	MG	C	553	1/1	0.80	0.17	40,40,40,40	0
3	FLC	B	552	13/13	0.80	0.21	60,62,64,64	0
3	FLC	C	551	13/13	0.83	0.12	45,48,51,53	0
3	FLC	A	551	13/13	0.84	0.13	30,42,51,54	0
4	MG	A	552	1/1	0.84	0.15	33,33,33,33	0
3	FLC	A	550	13/13	0.85	0.12	46,49,50,50	0
3	FLC	B	555	13/13	0.87	0.13	40,47,54,57	0
4	MG	B	556	1/1	0.88	0.12	36,36,36,36	0
3	FLC	C	552	13/13	0.90	0.11	30,43,51,55	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 5GP B 551:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.