



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 14, 2026 – 11:25 AM UTC

PDB ID : 3E8E / pdb_00003e8e
Title : Crystal structures of the kinase domain of PKA in complex with ATP-competitive inhibitors
Authors : Concha, N.O.; Elkins, P.A.; Smallwood, A.; Ward, P.
Deposited on : 2008-08-19
Resolution : 2.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

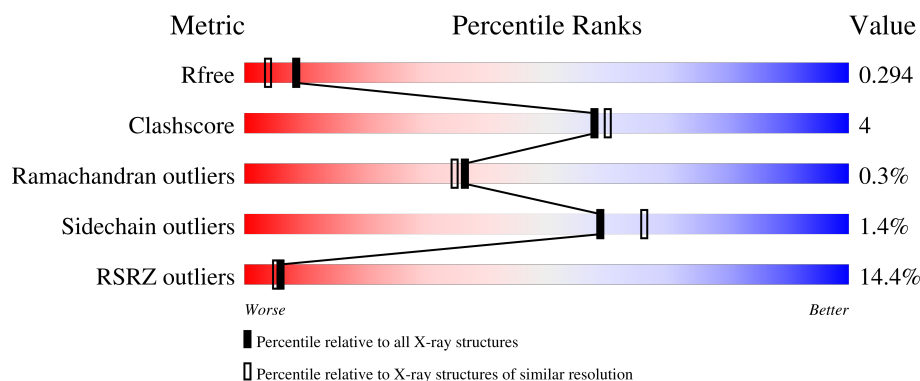
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	350	
1	B	350	
1	E	350	
1	I	350	
1	L	350	

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Mol	Chain	Length	Quality of chain
1	P	350	12% 88% 9% ..
2	C	20	40% 70% 30%
2	F	20	25% 70% 30%
2	G	20	35% 80% 15% 5%
2	J	20	40% 90% 10%
2	N	20	35% 90% 10%
2	Q	20	45% 80% 20%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18474 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.

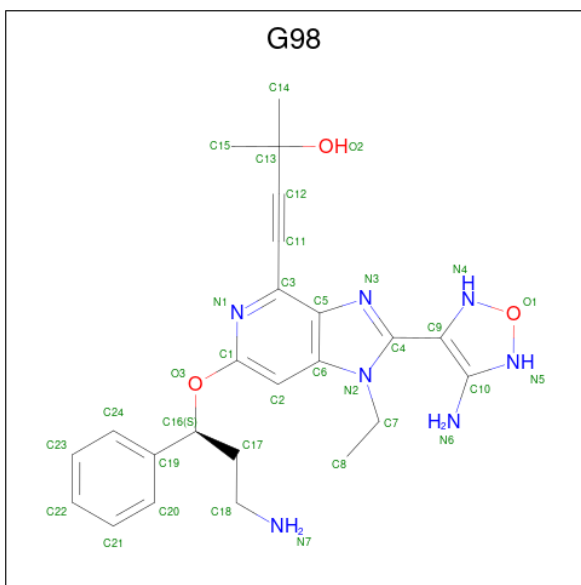
- Molecule 1 is a protein called cAMP-dependent protein kinase catalytic subunit alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	P	S	0	1	0
			2848	1840	477	519	3	9			
1	B	339	Total	C	N	O	P	S	0	1	0
			2804	1816	469	508	2	9			
1	E	344	Total	C	N	O	P	S	0	1	0
			2840	1834	475	519	3	9			
1	I	348	Total	C	N	O	P	S	0	1	0
			2858	1845	478	523	3	9			
1	L	338	Total	C	N	O	P	S	0	1	0
			2798	1813	468	506	2	9			
1	P	344	Total	C	N	O	P	S	0	2	0
			2844	1837	475	520	3	9			

- Molecule 2 is a protein called PKI inhibitor peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	20	Total	C	N	O	0	0	0
			157	94	32	31			
2	C	20	Total	C	N	O	0	0	0
			157	94	32	31			
2	F	20	Total	C	N	O	0	0	0
			157	94	32	31			
2	J	20	Total	C	N	O	0	0	0
			157	94	32	31			
2	N	20	Total	C	N	O	0	0	0
			157	94	32	31			
2	Q	20	Total	C	N	O	0	0	0
			157	94	32	31			

- Molecule 3 is 4-[2-(4-amino-2,5-dihydro-1,2,5-oxadiazol-3-yl)-6-{{[(1S)-3-amino-1-phenylpropyl]oxy}}-1-ethyl-1H-imidazo[4,5-c]pyridin-4-yl]-2-methylbut-3-yn-2-ol (CCD ID: G98) (formula: C₂₄H₂₉N₇O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			34	24	7	3		
3	B	1	Total	C	N	O	0	0
			34	24	7	3		
3	E	1	Total	C	N	O	0	0
			34	24	7	3		
3	I	1	Total	C	N	O	0	0
			34	24	7	3		
3	L	1	Total	C	N	O	0	0
			34	24	7	3		
3	P	1	Total	C	N	O	0	0
			34	24	7	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	74	Total	O	0	0
			74	74		
4	G	3	Total	O	0	0
			3	3		
4	B	65	Total	O	0	0
			65	65		
4	C	4	Total	O	0	0
			4	4		
4	E	76	Total	O	0	0
			76	76		
4	F	2	Total	O	0	0
			2	2		

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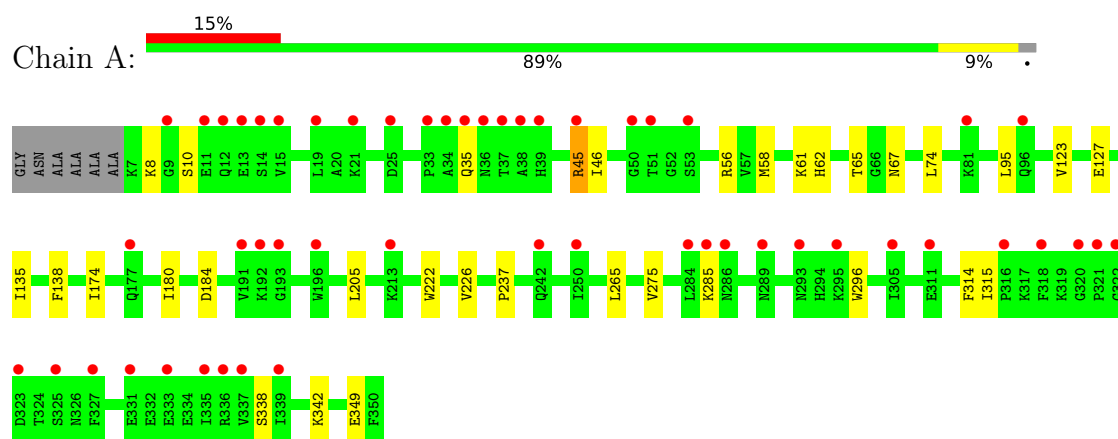
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	I	38	Total 38	O 38	0	0
4	J	2	Total 2	O 2	0	0
4	L	41	Total 41	O 41	0	0
4	N	4	Total 4	O 4	0	0
4	P	26	Total 26	O 26	0	0
4	Q	1	Total 1	O 1	0	0

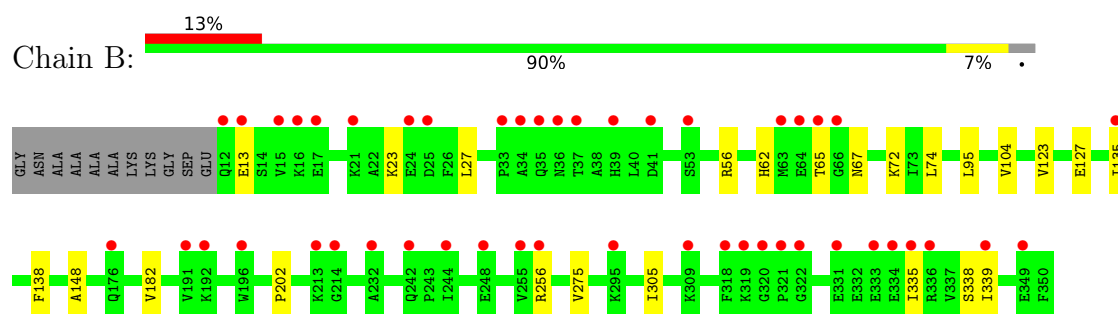
3 Residue-property plots [i](#)

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.

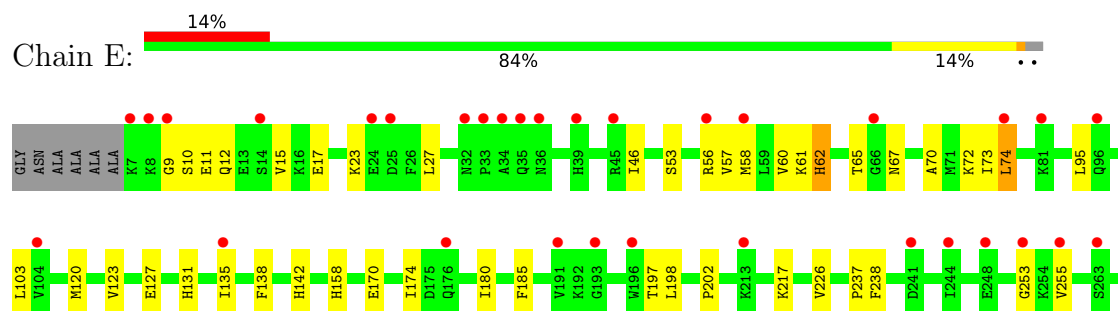
- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha

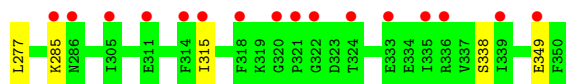


- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha

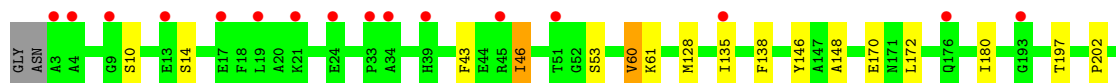
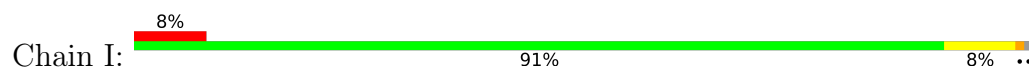


- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha

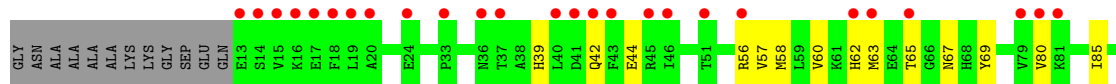
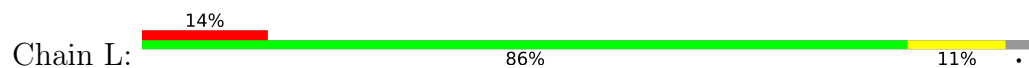




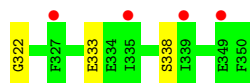
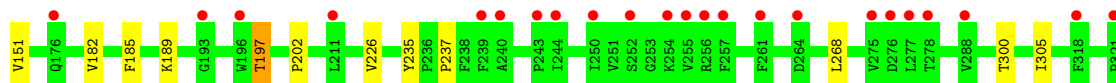
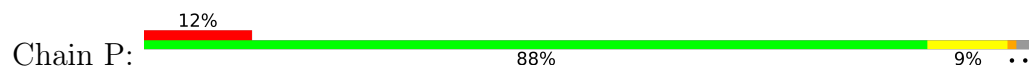
- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha



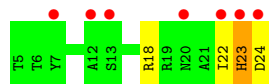
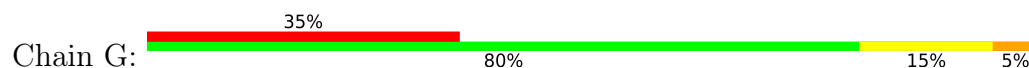
- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha



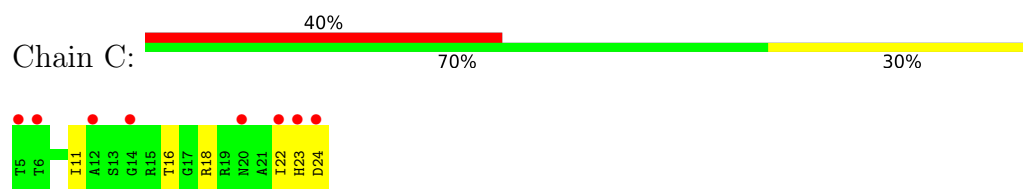
- Molecule 1: cAMP-dependent protein kinase catalytic subunit alpha



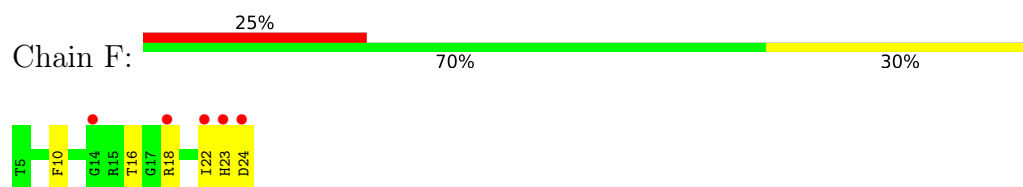
- Molecule 2: PKI inhibitor peptide



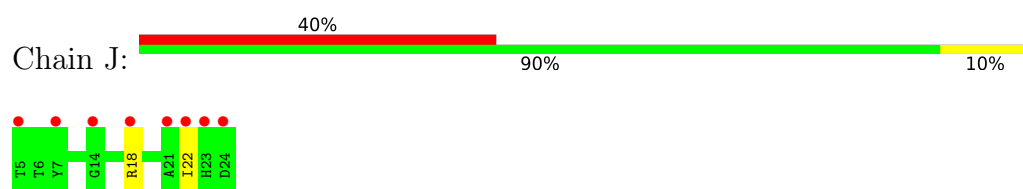
● Molecule 2: PKI inhibitor peptide



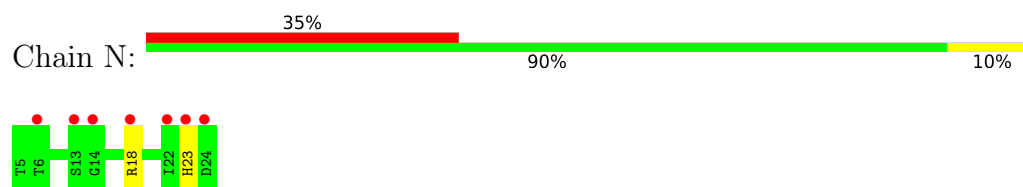
● Molecule 2: PKI inhibitor peptide



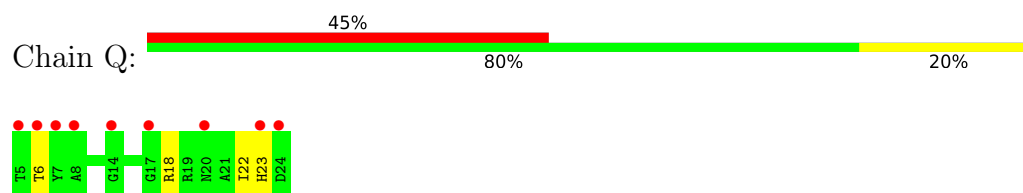
● Molecule 2: PKI inhibitor peptide



● Molecule 2: PKI inhibitor peptide



● Molecule 2: PKI inhibitor peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.58Å 95.85Å 179.10Å 90.00° 102.53° 90.00°	Depositor
Resolution (Å)	42.03 – 2.00 42.03 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.5 (42.03-2.00) 93.5 (42.03-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.243 , 0.279 0.259 , 0.294	Depositor DCC
R_{free} test set	8636 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.681	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.001 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	18474	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 25.30 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.2535e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: TPO, SEP, G98

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	0.50	0/2890	0.74	0/3888
1	B	0.50	0/2857	0.74	0/3849
1	E	0.52	0/2882	0.79	0/3880
1	I	0.47	0/2900	0.74	1/3906 (0.0%)
1	L	0.47	0/2851	0.74	0/3840
1	P	0.46	0/2889	0.73	0/3890
2	C	0.53	0/159	0.82	0/212
2	F	0.55	0/159	0.80	0/212
2	G	0.47	0/159	0.70	0/212
2	J	0.46	0/159	0.68	0/212
2	N	0.51	0/159	0.80	0/212
2	Q	0.55	0/159	0.76	0/212
All	All	0.49	0/18223	0.75	1/24525 (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	E	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	46	ILE	CB-CA-C	-5.17	107.31	111.71

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	62	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	A	2848	0	2824	25	0
1	B	2804	0	2776	21	0
1	E	2840	0	2802	38	0
1	I	2858	0	2815	20	0
1	L	2798	0	2777	21	0
1	P	2844	0	2803	20	0
2	C	157	0	146	4	0
2	F	157	0	146	6	0
2	G	157	0	146	3	0
2	J	157	0	146	2	0
2	N	157	0	146	1	0
2	Q	157	0	146	3	0
3	A	34	0	29	6	0
3	B	34	0	29	5	0
3	E	34	0	29	7	0
3	I	34	0	29	2	0
3	L	34	0	29	1	0
3	P	34	0	29	3	0
4	A	74	0	0	1	0
4	B	65	0	0	0	0
4	C	4	0	0	0	0
4	E	76	0	0	1	0
4	F	2	0	0	0	0
4	G	3	0	0	0	0
4	I	38	0	0	1	0
4	J	2	0	0	0	0
4	L	41	0	0	0	0
4	N	4	0	0	0	0
4	P	26	0	0	0	0
4	Q	1	0	0	0	0
All	All	18474	0	17847	162	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 4.

The worst 5 of 162 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:135:ILE:HD11	1:B:138:PHE:CD1	2.00	0.97
1:A:135:ILE:HD11	1:A:138:PHE:CD1	2.03	0.93
1:E:135:ILE:HD11	1:E:138:PHE:CD1	2.05	0.92
1:B:135:ILE:HD11	1:B:138:PHE:HD1	1.43	0.84
1:E:135:ILE:HD11	1:E:138:PHE:HD1	1.41	0.83

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	340/350 (97%)	332 (98%)	8 (2%)	0	100	100
1	B	336/350 (96%)	327 (97%)	9 (3%)	0	100	100
1	E	340/350 (97%)	327 (96%)	13 (4%)	0	100	100
1	I	344/350 (98%)	333 (97%)	11 (3%)	0	100	100
1	L	335/350 (96%)	327 (98%)	8 (2%)	0	100	100
1	P	341/350 (97%)	333 (98%)	7 (2%)	1 (0%)	36	35
2	C	18/20 (90%)	17 (94%)	0	1 (6%)	1	0
2	F	18/20 (90%)	17 (94%)	0	1 (6%)	1	0
2	G	18/20 (90%)	17 (94%)	0	1 (6%)	1	0
2	J	18/20 (90%)	16 (89%)	2 (11%)	0	100	100
2	N	18/20 (90%)	16 (89%)	1 (6%)	1 (6%)	1	0
2	Q	18/20 (90%)	17 (94%)	0	1 (6%)	1	0
All	All	2144/2220 (97%)	2079 (97%)	59 (3%)	6 (0%)	36	35

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	N	23	HIS
2	Q	23	HIS
2	F	23	HIS
2	G	23	HIS
2	C	23	HIS

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	302/302 (100%)	298 (99%)	4 (1%)	61	68
1	B	298/302 (99%)	296 (99%)	2 (1%)	76	82
1	E	300/302 (99%)	293 (98%)	7 (2%)	44	49
1	I	299/302 (99%)	293 (98%)	6 (2%)	48	54
1	L	298/302 (99%)	293 (98%)	5 (2%)	53	60
1	P	300/302 (99%)	297 (99%)	3 (1%)	68	75
2	C	15/15 (100%)	15 (100%)	0	100	100
2	F	15/15 (100%)	15 (100%)	0	100	100
2	G	15/15 (100%)	15 (100%)	0	100	100
2	J	15/15 (100%)	15 (100%)	0	100	100
2	N	15/15 (100%)	15 (100%)	0	100	100
2	Q	15/15 (100%)	15 (100%)	0	100	100
All	All	1887/1902 (99%)	1860 (99%)	27 (1%)	59	66

5 of 27 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	53	SER
1	I	268	LEU
1	P	81	LYS
1	I	256	ARG

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Mol	Chain	Res	Type
1	I	333	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 20 such sidechains are listed below:

Mol	Chain	Res	Type
1	I	274	GLN
1	L	181	GLN
1	P	36	ASN
1	L	326	ASN
2	C	23	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

16 non-standard protein/DNA/RNA residues 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$
1	TPO	A	197	1	8,10,11	0.75	0	10,14,16	1.01	0
1	SEP	P	10	1	8,9,10	1.75	1 (12%)	7,12,14	1.64	1 (14%)
1	SEP	E	10	1	8,9,10	1.70	1 (12%)	7,12,14	2.39	1 (14%)
1	TPO	E	197	1	8,10,11	0.85	0	10,14,16	1.15	1 (10%)
1	SEP	I	338	1	8,9,10	1.62	1 (12%)	7,12,14	1.64	1 (14%)
1	SEP	P	338	1	8,9,10	1.55	1 (12%)	7,12,14	1.74	2 (28%)
1	SEP	I	10	1	8,9,10	1.68	1 (12%)	7,12,14	3.73	2 (28%)
1	TPO	P	197	1	8,10,11	0.77	0	10,14,16	1.11	1 (10%)
1	SEP	L	338	1	8,9,10	1.69	1 (12%)	7,12,14	1.91	2 (28%)
1	SEP	A	338	1	8,9,10	1.48	1 (12%)	7,12,14	1.59	2 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	TPO	I	197	1	8,10,11	0.76	0	10,14,16	1.16	1 (10%)
1	TPO	L	197	1	8,10,11	0.77	0	10,14,16	1.14	1 (10%)
1	SEP	B	338	1	8,9,10	1.54	1 (12%)	7,12,14	1.68	2 (28%)
1	SEP	E	338	1	8,9,10	1.60	1 (12%)	7,12,14	1.87	1 (14%)
1	SEP	A	10	1	8,9,10	1.73	1 (12%)	7,12,14	2.06	2 (28%)
1	TPO	B	197	1	8,10,11	0.77	0	10,14,16	0.91	0

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
1	TPO	A	197	1	-	2/9/11/13	-
1	SEP	P	10	1	-	0/6/8/10	-
1	SEP	E	10	1	-	4/6/8/10	-
1	TPO	E	197	1	-	1/9/11/13	-
1	SEP	I	338	1	-	3/6/8/10	-
1	SEP	P	338	1	-	4/6/8/10	-
1	SEP	I	10	1	-	2/6/8/10	-
1	TPO	P	197	1	-	0/9/11/13	-
1	SEP	L	338	1	-	3/6/8/10	-
1	SEP	A	338	1	-	4/6/8/10	-
1	TPO	I	197	1	-	1/9/11/13	-
1	TPO	L	197	1	-	1/9/11/13	-
1	SEP	B	338	1	-	5/6/8/10	-
1	SEP	E	338	1	-	3/6/8/10	-
1	SEP	A	10	1	-	3/6/8/10	-
1	TPO	B	197	1	-	1/9/11/13	-

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	P	10	SEP	P-O1P	3.81	1.62	1.50
1	A	10	SEP	P-O1P	3.71	1.62	1.50
1	L	338	SEP	P-O1P	3.62	1.61	1.50
1	I	338	SEP	P-O1P	3.57	1.61	1.50
1	E	338	SEP	P-O1P	3.48	1.61	1.50

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	10	SEP	OG-CB-CA	9.35	117.24	108.14
1	E	10	SEP	OG-CB-CA	5.79	113.78	108.14
1	A	10	SEP	OG-CB-CA	4.74	112.76	108.14
1	E	338	SEP	OG-CB-CA	4.47	112.50	108.14
1	L	338	SEP	OG-CB-CA	4.38	112.41	108.14

There are no chirality outliers.

5 of 37 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	10	SEP	CB-OG-P-O1P
1	A	10	SEP	CB-OG-P-O2P
1	A	10	SEP	CB-OG-P-O3P
1	A	197	TPO	O-C-CA-CB
1	A	338	SEP	C-CA-CB-OG

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	P	10	SEP	2	0
1	P	197	TPO	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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
3	G98	E	351	-	32,37,37	2.59	4 (12%)	39,53,53	1.80	8 (20%)
3	G98	I	351	-	32,37,37	2.62	4 (12%)	39,53,53	1.72	8 (20%)
3	G98	L	351	-	32,37,37	2.66	4 (12%)	39,53,53	1.70	9 (23%)
3	G98	B	351	-	32,37,37	2.64	3 (9%)	39,53,53	1.78	9 (23%)
3	G98	P	351	-	32,37,37	2.63	4 (12%)	39,53,53	1.60	7 (17%)
3	G98	A	351	-	32,37,37	2.76	3 (9%)	39,53,53	1.70	8 (20%)

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	G98	E	351	-	-	8/21/33/33	0/4/4/4
3	G98	I	351	-	-	6/21/33/33	0/4/4/4
3	G98	L	351	-	-	6/21/33/33	0/4/4/4
3	G98	B	351	-	-	10/21/33/33	0/4/4/4
3	G98	P	351	-	-	8/21/33/33	0/4/4/4
3	G98	A	351	-	-	7/21/33/33	0/4/4/4

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	351	G98	C3-C11	-14.42	1.28	1.43
3	L	351	G98	C3-C11	-13.83	1.28	1.43
3	B	351	G98	C3-C11	-13.77	1.28	1.43
3	P	351	G98	C3-C11	-13.72	1.29	1.43
3	I	351	G98	C3-C11	-13.53	1.29	1.43

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	351	G98	C5-C6-N2	5.91	109.54	105.22
3	E	351	G98	C5-C6-N2	5.82	109.47	105.22
3	B	351	G98	C5-C6-N2	5.70	109.38	105.22
3	L	351	G98	C5-C6-N2	5.53	109.26	105.22
3	P	351	G98	C5-C6-N2	5.27	109.07	105.22

There are no chirality outliers.

5 of 45 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	351	G98	C16-C17-C18-N7
3	A	351	G98	C3-C11-C12-C13
3	A	351	G98	N1-C1-O3-C16
3	A	351	G98	C2-C1-O3-C16
3	A	351	G98	C17-C16-O3-C1

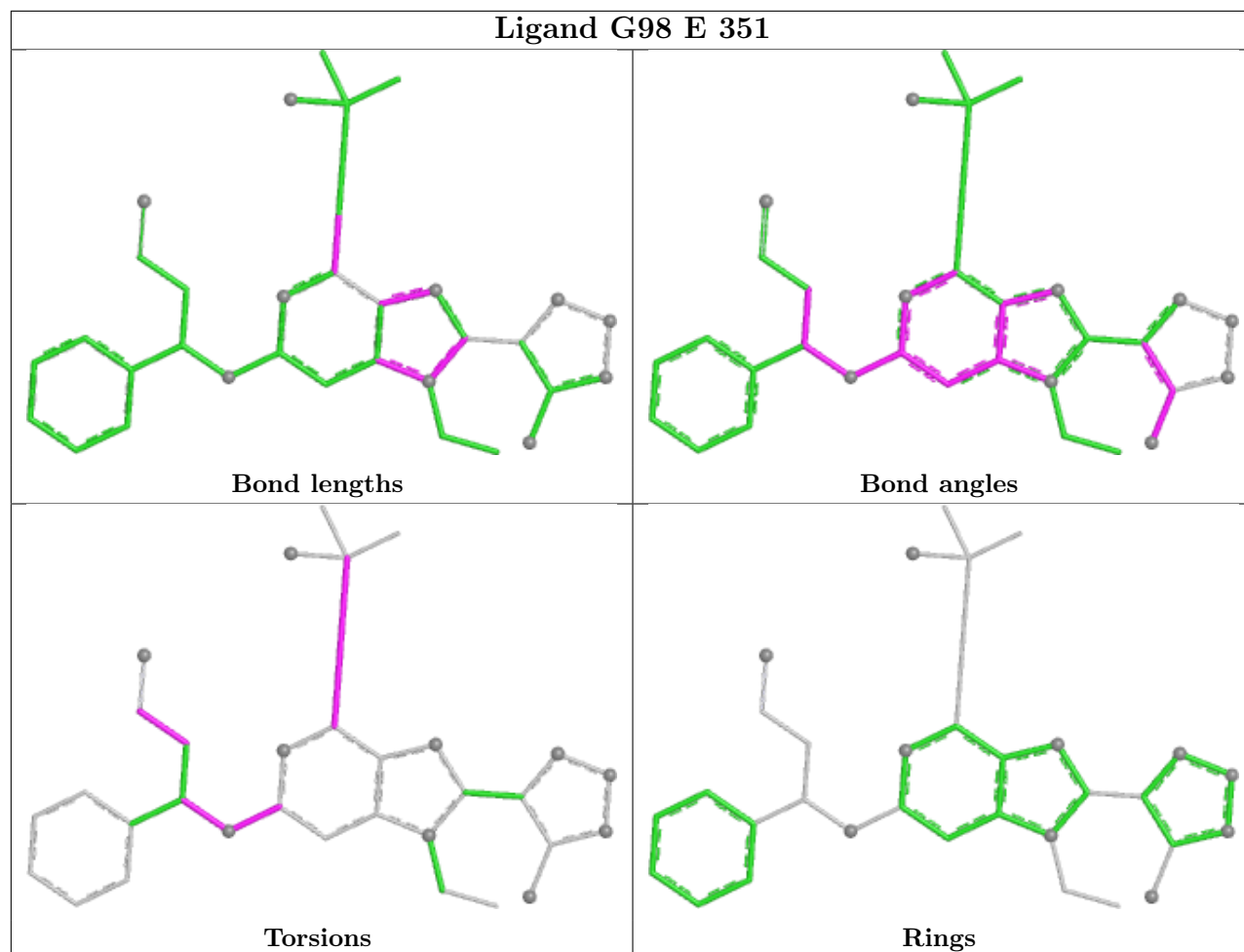
There are no ring outliers.

6 monomers are involved in 24 short contacts:

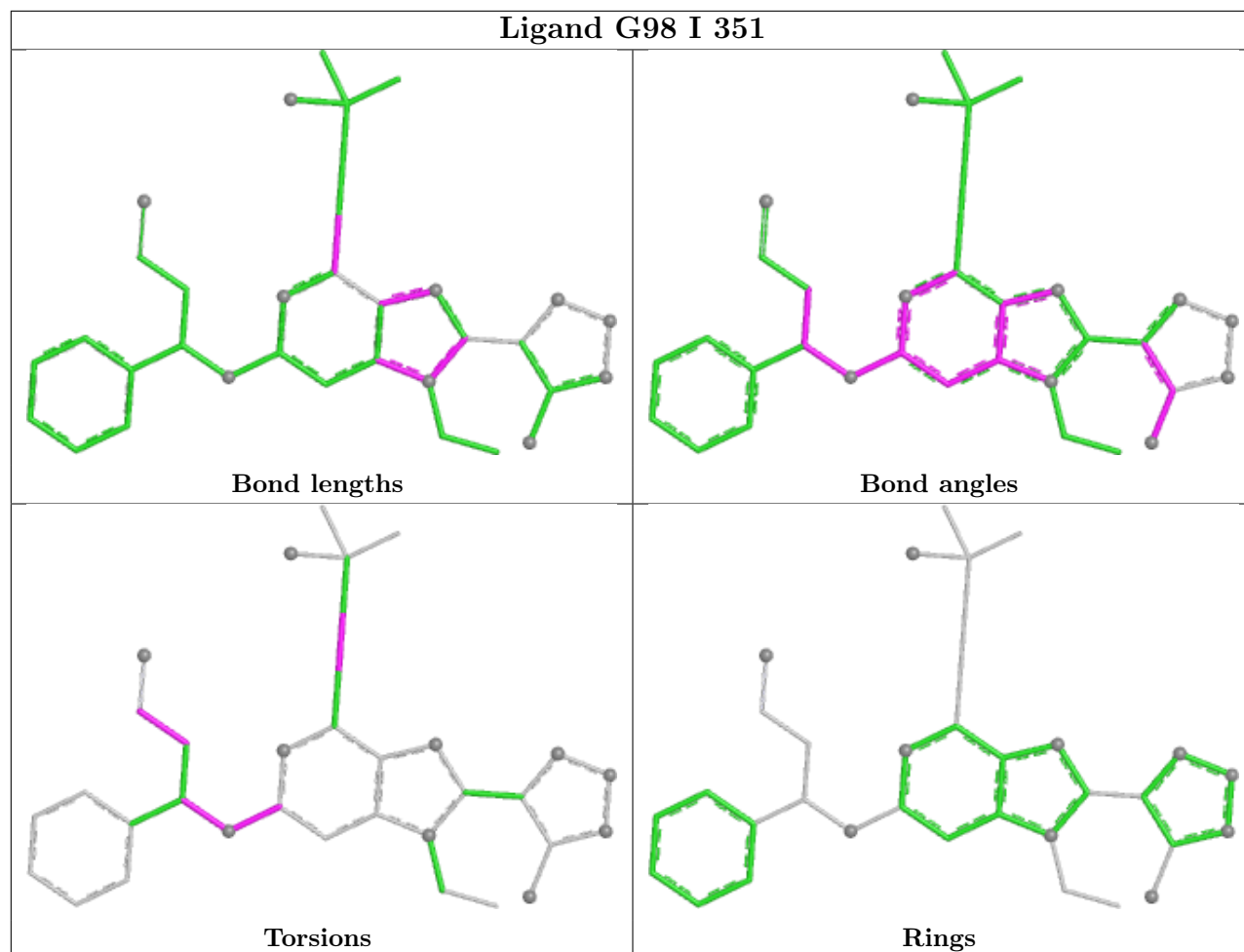
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	351	G98	7	0
3	I	351	G98	2	0
3	L	351	G98	1	0
3	B	351	G98	5	0
3	P	351	G98	3	0
3	A	351	G98	6	0

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.

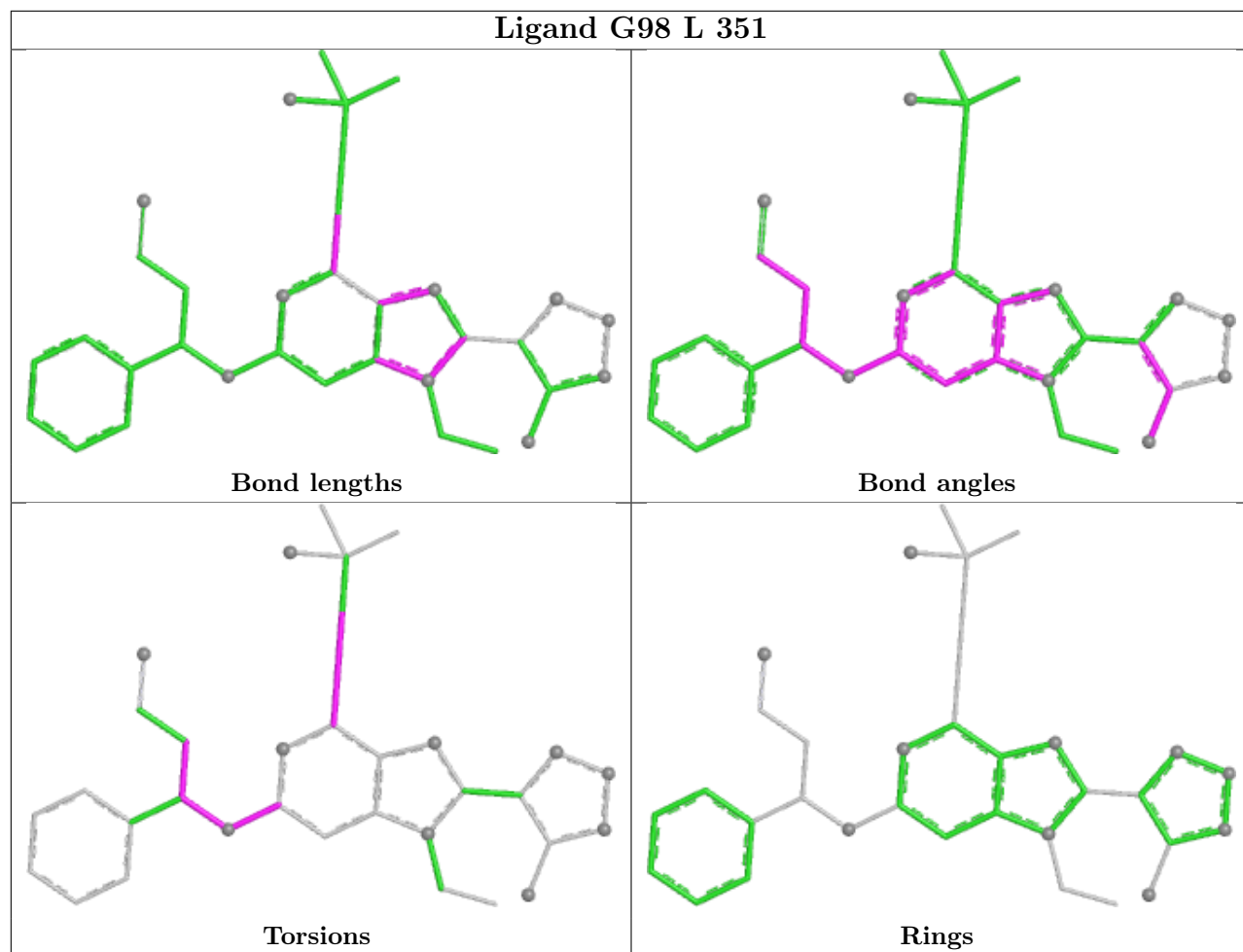
Ligand G98 E 351



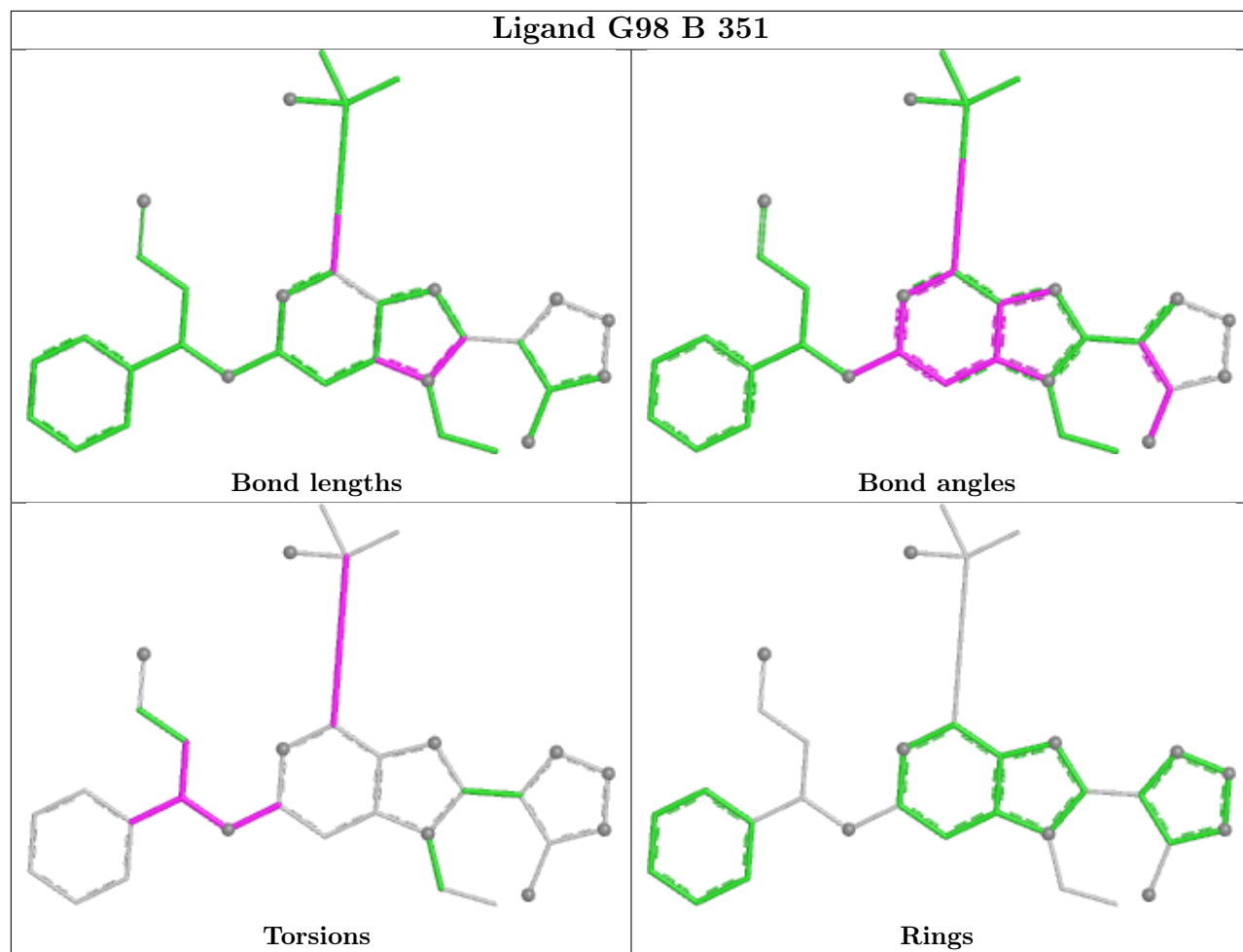
Ligand G98 I 351



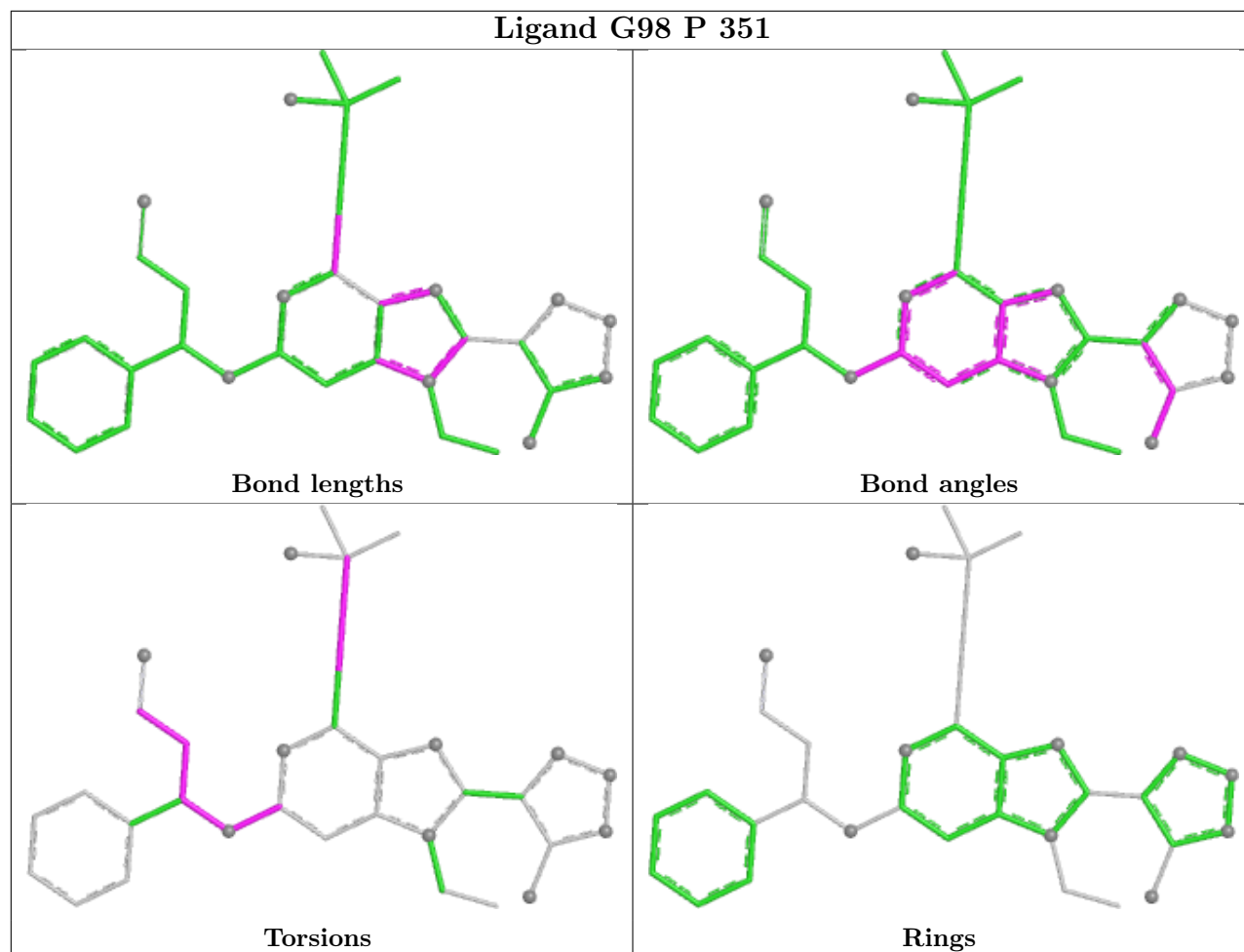
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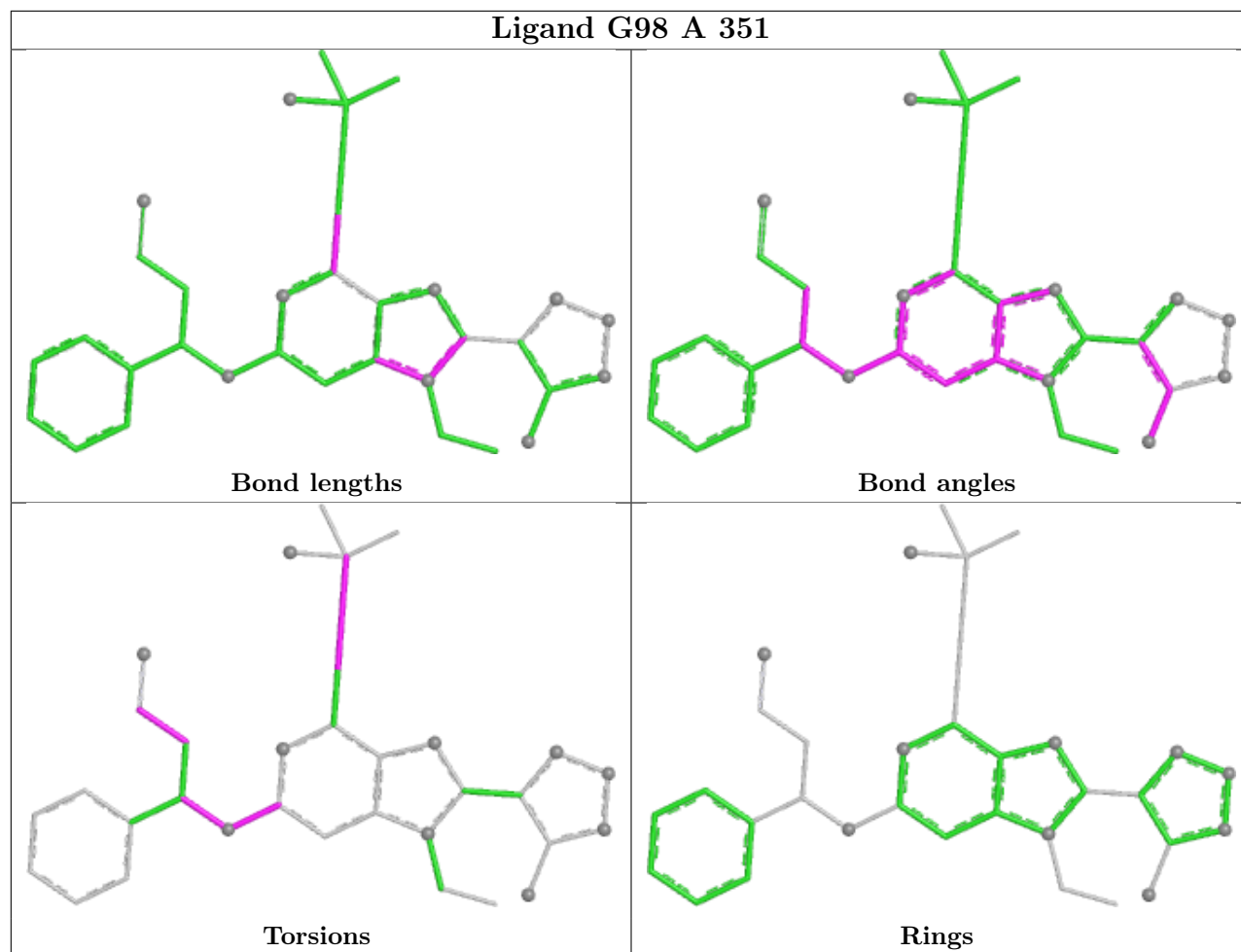


Ligand G98 B 351



Ligand G98 P 351





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	A	341/350 (97%)	1.01	52 (15%) 5 5	19, 36, 48, 76	1 (0%)
1	B	337/350 (96%)	1.05	47 (13%) 6 5	19, 36, 47, 58	1 (0%)
1	E	341/350 (97%)	1.15	48 (14%) 6 5	19, 36, 46, 54	1 (0%)
1	I	345/350 (98%)	0.88	29 (8%) 17 15	19, 36, 47, 55	1 (0%)
1	L	336/350 (96%)	1.12	48 (14%) 6 5	19, 37, 47, 55	1 (0%)
1	P	341/350 (97%)	1.11	43 (12%) 8 7	19, 37, 48, 68	2 (0%)
2	C	20/20 (100%)	2.47	8 (40%) 1 1	37, 42, 53, 53	0
2	F	20/20 (100%)	1.91	5 (25%) 2 1	36, 41, 54, 54	0
2	G	20/20 (100%)	1.78	7 (35%) 1 1	39, 43, 56, 56	0
2	J	20/20 (100%)	1.93	8 (40%) 1 1	39, 42, 52, 53	0
2	N	20/20 (100%)	1.69	7 (35%) 1 1	38, 42, 55, 56	0
2	Q	20/20 (100%)	2.09	9 (45%) 0 1	39, 43, 54, 55	0
All	All	2161/2220 (97%)	1.10	311 (14%) 6 5	19, 37, 48, 76	7 (0%)

The worst 5 of 311 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	P	34	ALA	8.3
2	F	24	ASP	7.2
2	C	5	THR	6.9
1	L	15	VAL	6.6
1	E	9	GLY	6.5

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	SEP	A	10	10/11	0.53	0.20	70,72,72,73	0
1	SEP	P	10	10/11	0.59	0.16	67,67,71,71	0
1	SEP	I	10	10/11	0.73	0.18	51,51,51,52	0
1	SEP	E	10	10/11	0.74	0.18	48,50,51,51	0
1	SEP	L	338	10/11	0.77	0.14	48,49,53,53	0
1	SEP	E	338	10/11	0.84	0.13	46,47,47,48	0
1	SEP	I	338	10/11	0.90	0.10	47,48,50,51	0
1	SEP	B	338	10/11	0.92	0.09	45,46,47,48	0
1	TPO	L	197	11/12	0.93	0.09	35,35,36,36	0
1	TPO	P	197	11/12	0.93	0.09	35,35,36,38	0
1	SEP	P	338	10/11	0.94	0.09	44,46,47,47	0
1	TPO	B	197	11/12	0.95	0.10	33,34,34,34	0
1	SEP	A	338	10/11	0.96	0.10	41,44,46,47	0
1	TPO	A	197	11/12	0.96	0.11	32,34,34,35	0
1	TPO	E	197	11/12	0.97	0.08	32,34,34,34	0
1	TPO	I	197	11/12	0.97	0.07	28,31,33,33	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

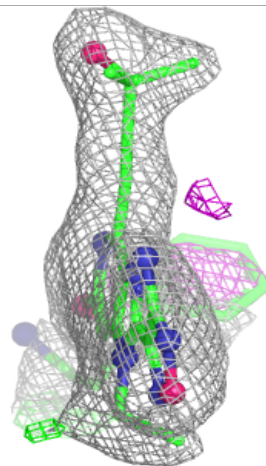
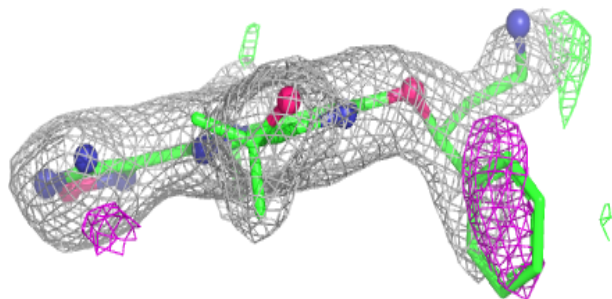
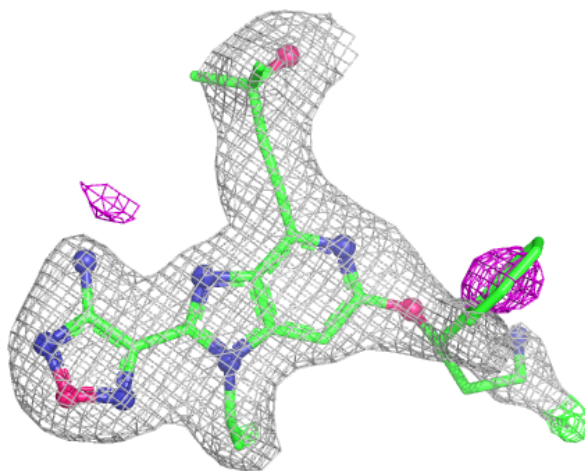
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	G98	L	351	34/34	0.87	0.15	39,42,51,51	0
3	G98	P	351	34/34	0.88	0.12	30,33,44,46	0
3	G98	B	351	34/34	0.89	0.11	25,28,37,40	0
3	G98	I	351	34/34	0.89	0.11	24,31,40,40	0
3	G98	A	351	34/34	0.90	0.10	21,24,35,37	0
3	G98	E	351	34/34	0.91	0.11	17,23,35,35	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.

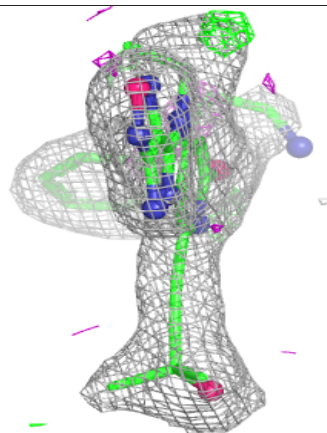
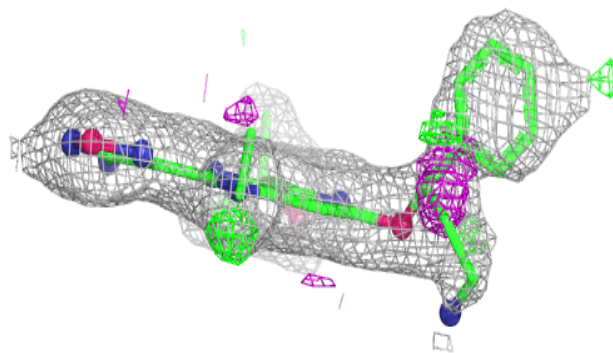
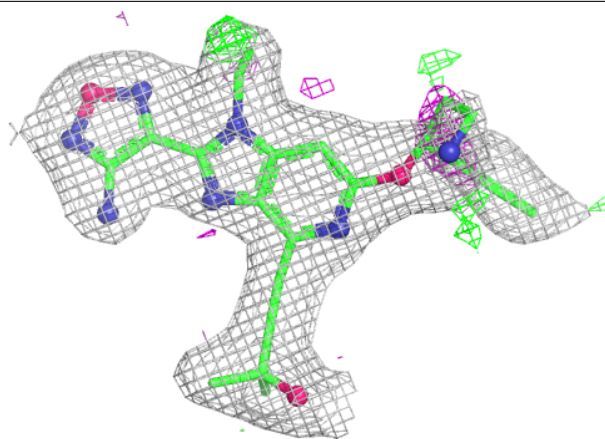
Electron density around G98 L 351:

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



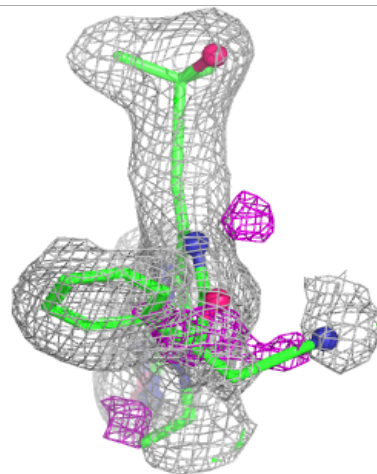
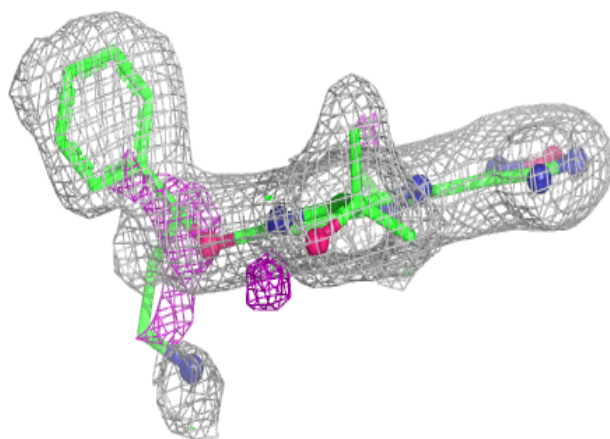
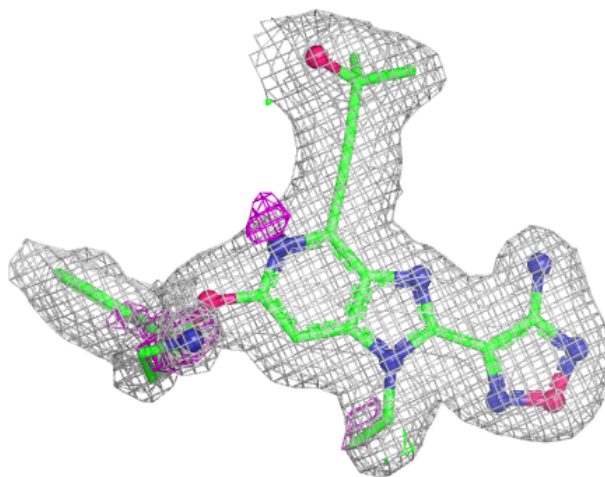
Electron density around G98 P 351:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



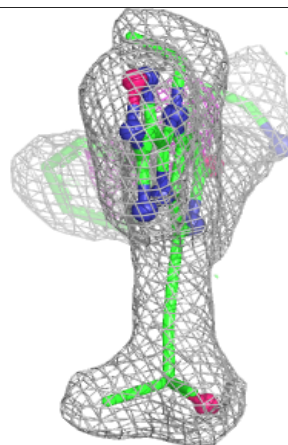
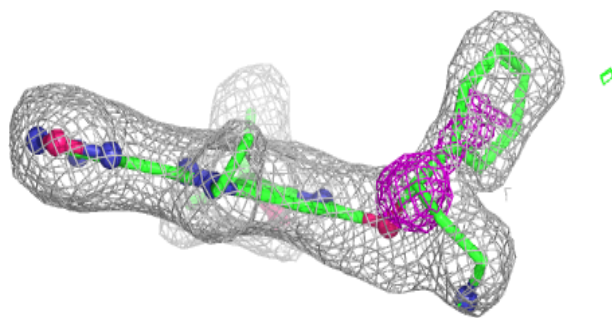
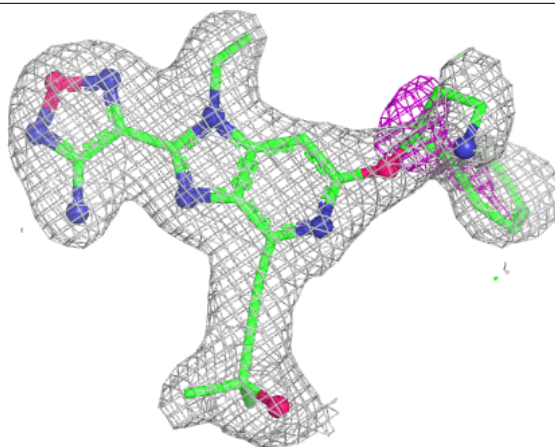
Electron density around G98 B 351:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



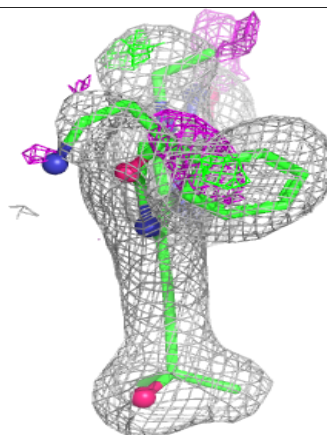
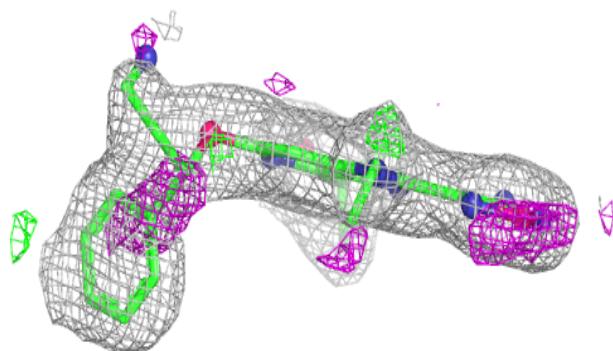
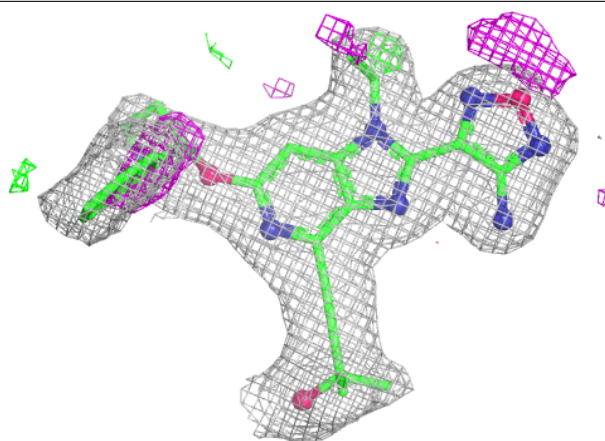
Electron density around G98 I 351:

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



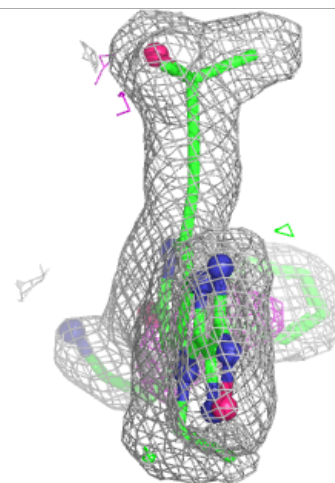
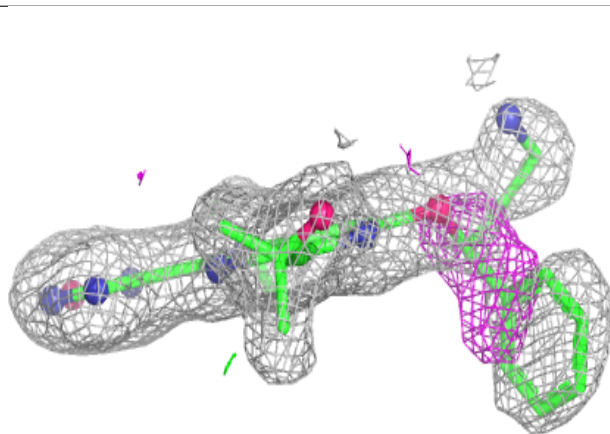
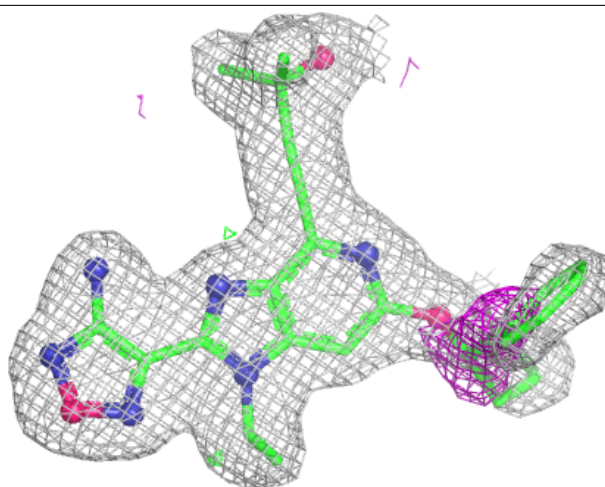
Electron density around G98 A 351:

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



Electron density around G98 E 351:

$2mF_o-DF_c$ (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.