



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 09:18 pm BST

PDB ID : 2HYY
Title : Human Abl kinase domain in complex with imatinib (STI571, Glivec)
Authors : Cowan-Jacob, S.W.; Fendrich, G.; Liebetanz, J.; Fabbro, D.; Manley, P.
Deposited on : 2006-08-08
Resolution : 2.40 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

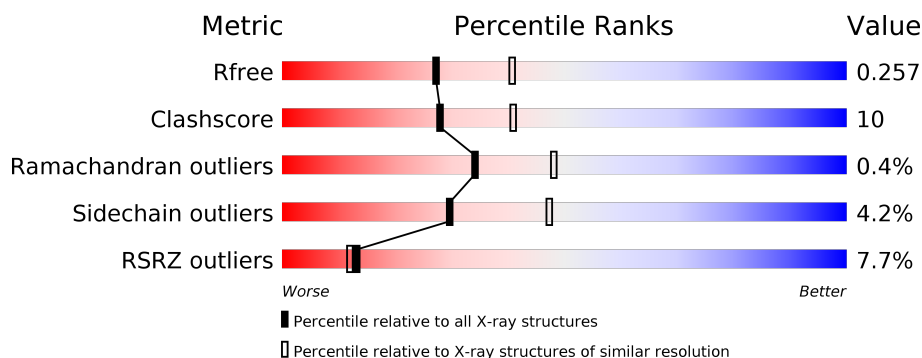
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 on 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	273	<div> <div>10%</div> <div> <div></div> <div>74%</div> <div>21%</div> <div>• •</div> </div> </div>
1	B	273	<div> <div>6%</div> <div> <div></div> <div>79%</div> <div>18%</div> <div>•</div> </div> </div>
1	C	273	<div> <div>6%</div> <div> <div></div> <div>73%</div> <div>20%</div> <div>• •</div> </div> </div>
1	D	273	<div> <div>8%</div> <div> <div></div> <div>69%</div> <div>25%</div> <div>• •</div> </div> </div>

2 Entry composition [i](#)

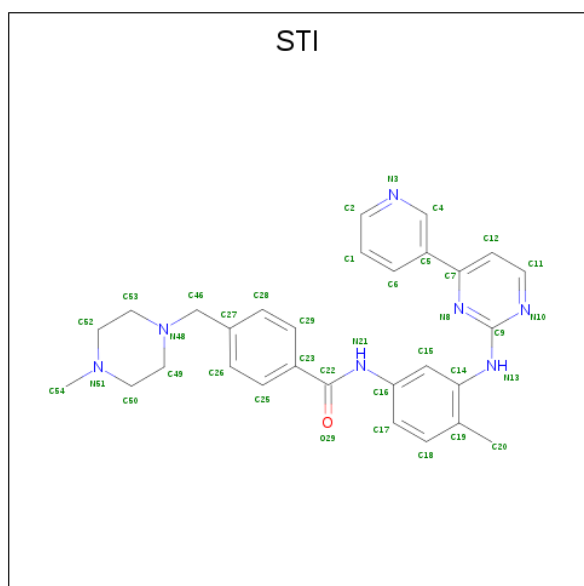
There are 3 unique types of molecules in this entry. The entry contains 8702 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Proto-oncogene tyrosine-protein kinase ABL1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2073	1343	332	382	16			
1	B	265	Total	C	N	O	S	0	0	0
			2100	1359	340	384	17			
1	C	262	Total	C	N	O	S	0	0	0
			2083	1350	338	379	16			
1	D	264	Total	C	N	O	S	0	0	0
			2098	1358	340	384	16			

- Molecule 2 is 4-(4-METHYL-PIPERAZIN-1-YLMETHYL)-N-[4-METHYL-3-(4-PYRIDIN-3-YL-PYRIMIDIN-2-YLAMINO)-PHENYL]-BENZAMIDE (three-letter code: STI) (formula: C₂₉H₃₁N₇O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			37	29	7	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			37	29	7	1		
2	C	1	Total	C	N	O	0	0
			37	29	7	1		
2	D	1	Total	C	N	O	0	0
			37	29	7	1		

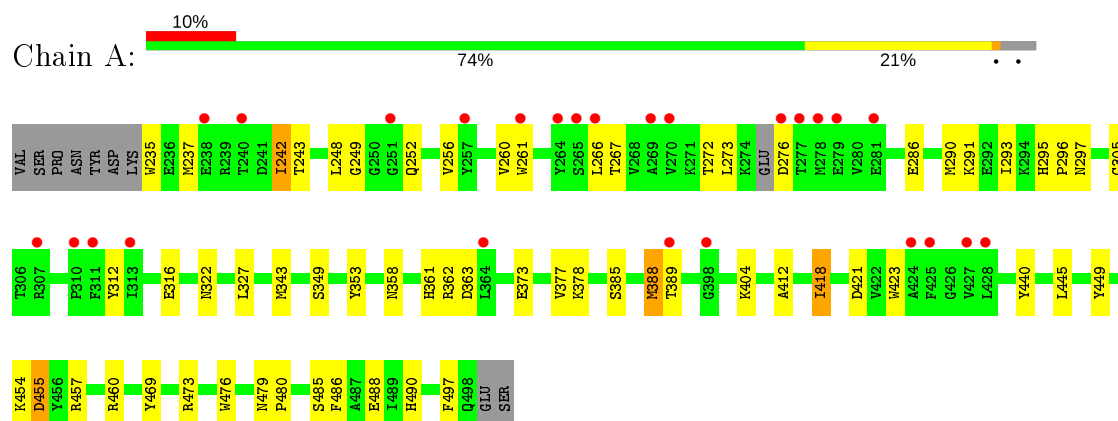
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	51	Total	O	0	0
			51	51		
3	B	45	Total	O	0	0
			45	45		
3	C	57	Total	O	0	0
			57	57		
3	D	47	Total	O	0	0
			47	47		

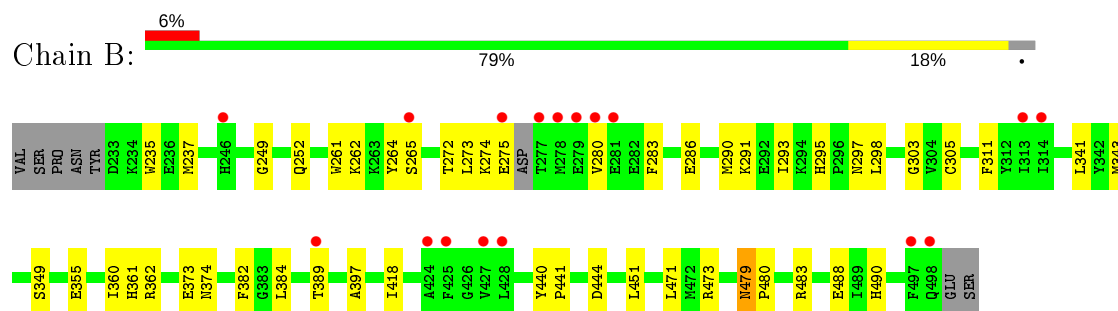
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

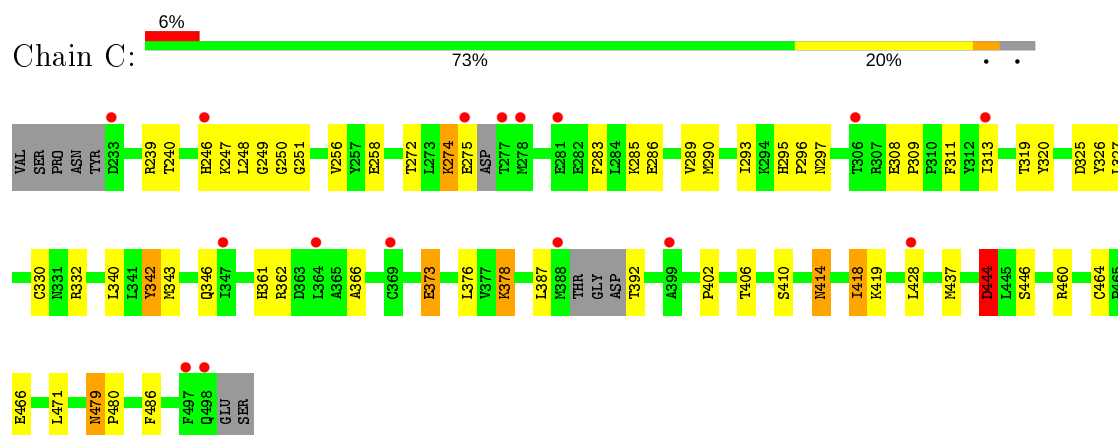
- Molecule 1: Proto-oncogene tyrosine-protein kinase ABL1



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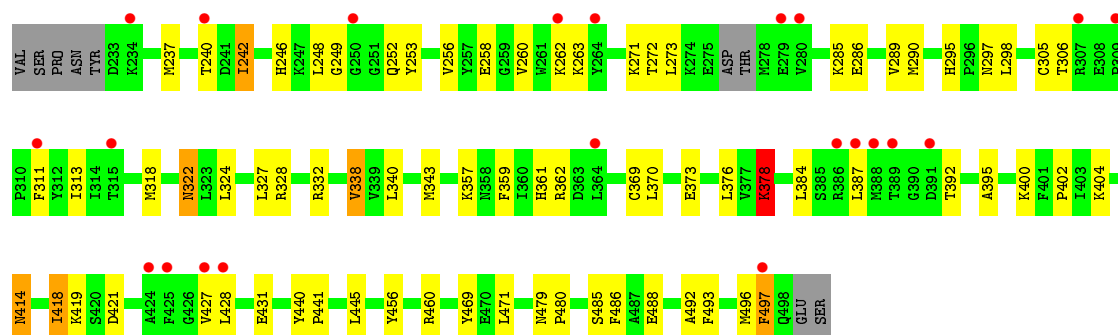


- Molecule 1: Proto-oncogene tyrosine-protein kinase ABL1



● Molecule 1: Proto-oncogene tyrosine-protein kinase ABL1

Chain D:  8% 69% 25%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	141.68Å 148.57Å 115.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.40 24.98 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.5 (25.00-2.40) 99.6 (24.98-2.40)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 2.39Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.204 , 0.267 0.197 , 0.257	Depositor DCC
R_{free} test set	2408 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.024 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8702	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.94% 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: STI

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	0.95	1/2127 (0.0%)	0.90	3/2891 (0.1%)
1	B	1.04	0/2155	0.95	4/2927 (0.1%)
1	C	1.01	2/2137 (0.1%)	0.93	5/2900 (0.2%)
1	D	0.96	0/2153	0.93	4/2923 (0.1%)
All	All	0.99	3/8572 (0.0%)	0.93	16/11641 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	373	GLU	CG-CD	-7.13	1.41	1.51
1	C	342	TYR	CD1-CE1	7.03	1.49	1.39
1	A	353	TYR	CD2-CE2	-5.07	1.31	1.39

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	332	ARG	NE-CZ-NH2	-8.31	116.14	120.30
1	D	378	LYS	CD-CE-NZ	8.20	130.56	111.70
1	A	473	ARG	NE-CZ-NH2	-7.20	116.70	120.30
1	B	444	ASP	CB-CG-OD1	-6.75	112.23	118.30
1	C	444	ASP	CB-CG-OD2	6.73	124.36	118.30
1	D	328	ARG	NE-CZ-NH2	-6.68	116.96	120.30
1	B	483	ARG	NE-CZ-NH2	-6.63	116.98	120.30
1	D	332	ARG	NE-CZ-NH1	-6.59	117.00	120.30
1	A	457	ARG	NE-CZ-NH2	-6.49	117.06	120.30
1	A	363	ASP	CB-CG-OD1	6.17	123.85	118.30
1	D	421	ASP	CB-CG-OD1	6.06	123.75	118.30
1	B	473	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	C	437	MET	CG-SD-CE	-5.54	91.34	100.20
1	C	332	ARG	NE-CZ-NH1	5.50	123.05	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	341	LEU	CB-CG-CD2	-5.32	101.96	111.00
1	C	325	ASP	CB-CG-OD2	-5.23	113.59	118.30

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	2073	0	1971	38	0
1	B	2100	0	2010	31	1
1	C	2083	0	1996	40	0
1	D	2098	0	2011	57	0
2	A	37	0	31	1	0
2	B	37	0	31	2	0
2	C	37	0	31	0	0
2	D	37	0	31	4	0
3	A	51	0	0	0	0
3	B	45	0	0	1	0
3	C	57	0	0	1	0
3	D	47	0	0	3	0
All	All	8702	0	8112	165	1

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 10.

All (165) 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:C:286:GLU:HG2	1:C:290:MET:HE2	1.35	1.08
1:D:272:THR:HG22	1:D:273:LEU:H	1.21	1.01
1:C:286:GLU:HG2	1:C:290:MET:CE	1.91	1.00
1:C:479:ASN:HD22	1:C:480:PRO:HD2	1.33	0.93
1:A:479:ASN:HD22	1:A:480:PRO:HD2	1.34	0.91
1:B:286:GLU:HG2	1:B:290:MET:HE2	1.50	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:295:HIS:HD2	1:D:297:ASN:H	1.19	0.88
1:A:295:HIS:HD2	1:A:297:ASN:H	1.18	0.88
1:A:286:GLU:HG2	1:A:290:MET:CE	2.08	0.83
1:B:286:GLU:HG2	1:B:290:MET:CE	2.09	0.81
1:D:479:ASN:HD22	1:D:480:PRO:HD2	1.45	0.80
1:A:295:HIS:CD2	1:A:297:ASN:H	2.02	0.77
1:D:286:GLU:HG2	1:D:290:MET:HE2	1.66	0.77
1:B:479:ASN:HD22	1:B:480:PRO:HD2	1.50	0.76
1:D:272:THR:HG22	1:D:273:LEU:N	2.00	0.73
1:D:295:HIS:CD2	1:D:297:ASN:H	2.06	0.72
1:C:239:ARG:NH2	1:C:308:GLU:O	2.22	0.72
1:C:286:GLU:CG	1:C:290:MET:HE2	2.16	0.72
1:C:444:ASP:OD1	1:C:446:SER:OG	2.07	0.72
1:C:309:PRO:HG3	1:C:311:PHE:CE1	2.25	0.72
1:D:253:TYR:OH	3:D:161:HOH:O	2.08	0.72
1:D:272:THR:CG2	1:D:273:LEU:H	2.01	0.72
1:C:248:LEU:HD12	1:C:256:VAL:HG12	1.76	0.68
1:C:295:HIS:HD2	1:C:297:ASN:H	1.42	0.67
1:A:349:SER:HB2	1:A:490:HIS:CE1	2.29	0.67
1:A:286:GLU:HG2	1:A:290:MET:HE2	1.77	0.66
1:A:327:LEU:HD21	1:A:343:MET:HE1	1.78	0.66
1:C:479:ASN:HD22	1:C:480:PRO:CD	2.08	0.65
1:B:488:GLU:OE1	3:B:191:HOH:O	2.13	0.65
1:D:493:PHE:HA	1:D:496:MET:HE3	1.80	0.64
1:B:360:ILE:HG22	1:B:362:ARG:HG3	1.79	0.64
1:B:249:GLY:O	1:B:252:GLN:HB2	1.98	0.64
1:A:361:HIS:O	1:A:362:ARG:HB2	1.98	0.63
1:B:389:THR:HB	1:C:466:GLU:HB3	1.81	0.63
1:D:242:ILE:HA	1:D:260:VAL:O	1.99	0.61
1:D:285:LYS:O	1:D:289:VAL:HG23	2.00	0.61
1:C:418:ILE:HD11	1:C:486:PHE:HD2	1.65	0.61
1:C:340:LEU:HA	1:C:343:MET:HE2	1.83	0.61
1:D:249:GLY:O	1:D:252:GLN:HB2	2.00	0.61
1:B:355:GLU:HB2	1:B:418:ILE:HG12	1.84	0.60
1:B:295:HIS:HD2	1:B:297:ASN:H	1.50	0.59
1:D:338:VAL:HG23	1:D:497:PHE:HE1	1.67	0.59
1:A:316:GLU:OE1	1:A:378:LYS:NZ	2.33	0.59
1:D:361:HIS:O	1:D:362:ARG:HB2	2.01	0.59
1:A:479:ASN:ND2	1:A:480:PRO:HD2	2.14	0.58
1:C:285:LYS:O	1:C:289:VAL:HG23	2.04	0.58
1:C:295:HIS:CD2	1:C:297:ASN:H	2.20	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:MET:HE1	1:A:242:ILE:HD12	1.86	0.58
1:B:451:LEU:HD21	1:D:456:TYR:HB2	1.85	0.58
1:D:237:MET:HE1	1:D:242:ILE:HD12	1.86	0.57
1:C:296:PRO:O	1:C:378:LYS:HD2	2.04	0.57
1:B:235:TRP:CD1	1:B:291:LYS:HG2	2.39	0.57
1:D:318:MET:HE1	3:D:53:HOH:O	2.05	0.57
1:C:414:ASN:O	1:C:414:ASN:ND2	2.37	0.56
1:D:318:MET:CE	3:D:53:HOH:O	2.54	0.56
1:D:340:LEU:HA	1:D:343:MET:HE2	1.87	0.56
1:A:295:HIS:HD2	1:A:297:ASN:N	1.97	0.56
1:D:376:LEU:HD21	1:D:378:LYS:HE2	1.87	0.56
1:B:361:HIS:O	1:B:362:ARG:HB2	2.06	0.55
1:A:404:LYS:HG2	1:A:440:TYR:CD1	2.42	0.55
1:A:327:LEU:HD21	1:A:343:MET:CE	2.36	0.54
1:A:248:LEU:HD12	1:A:256:VAL:HG12	1.88	0.54
1:B:349:SER:HB2	1:B:490:HIS:CE1	2.43	0.54
1:D:492:ALA:O	1:D:496:MET:HB2	2.08	0.53
1:B:295:HIS:HB3	1:B:298:LEU:HD12	1.91	0.53
1:D:404:LYS:HE3	1:D:445:LEU:HD23	1.90	0.53
1:A:485:SER:OG	1:A:488:GLU:HG3	2.09	0.53
1:D:392:THR:O	1:D:402:PRO:HA	2.08	0.52
1:D:237:MET:CE	1:D:242:ILE:HD12	2.39	0.52
1:A:454:LYS:O	1:A:455:ASP:HB2	2.07	0.52
1:A:296:PRO:O	1:A:378:LYS:HE2	2.09	0.52
1:D:404:LYS:HE3	1:D:445:LEU:CD2	2.40	0.52
1:A:260:VAL:HA	1:A:266:LEU:O	2.10	0.52
1:A:418:ILE:O	1:A:421:ASP:HB2	2.10	0.52
1:A:343:MET:HG2	1:A:377:VAL:HG21	1.90	0.52
1:C:248:LEU:HD21	1:C:258:GLU:HB2	1.93	0.52
1:B:479:ASN:HD22	1:B:480:PRO:CD	2.22	0.51
1:C:460:ARG:NH1	1:C:464:CYS:O	2.41	0.51
1:A:261:TRP:N	1:A:266:LEU:O	2.37	0.51
1:C:418:ILE:HD11	1:C:486:PHE:CD2	2.45	0.51
1:D:246:HIS:NE2	1:D:258:GLU:OE1	2.44	0.51
1:B:272:THR:HG22	1:B:273:LEU:N	2.26	0.51
1:D:262:LYS:O	1:D:263:LYS:C	2.48	0.51
1:C:283:PHE:CE2	1:C:313:ILE:HG13	2.46	0.50
1:B:295:HIS:CD2	1:B:297:ASN:H	2.28	0.50
1:D:289:VAL:HG13	1:D:359:PHE:HE2	1.77	0.49
1:C:319:THR:HB	1:C:320:TYR:CD1	2.48	0.49
1:C:392:THR:O	1:C:402:PRO:HA	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:318:MET:O	2:D:600:STI:H21	2.12	0.49
2:A:600:STI:H151	2:A:600:STI:O29	2.13	0.49
1:D:305:CYS:O	1:D:311:PHE:HA	2.13	0.48
1:B:237:MET:SD	1:B:303:GLY:HA3	2.53	0.48
1:C:286:GLU:HG2	1:C:290:MET:HE1	1.89	0.48
1:D:418:ILE:HD11	1:D:486:PHE:HD2	1.78	0.48
1:B:286:GLU:CG	1:B:290:MET:HE2	2.34	0.48
1:A:449:TYR:CG	1:B:397:ALA:HA	2.49	0.48
1:A:418:ILE:HD11	1:A:486:PHE:HD2	1.78	0.48
1:D:414:ASN:O	1:D:414:ASN:ND2	2.46	0.48
1:A:272:THR:HG22	1:A:273:LEU:N	2.29	0.48
1:A:423:TRP:CE3	1:A:476:TRP:HA	2.49	0.48
1:A:479:ASN:HD22	1:A:480:PRO:CD	2.16	0.47
1:C:327:LEU:HD21	1:C:343:MET:HE1	1.97	0.47
1:D:295:HIS:HB3	1:D:298:LEU:HD12	1.96	0.47
1:D:419:LYS:HD2	1:D:480:PRO:O	2.14	0.47
1:D:384:LEU:HD23	1:D:387:LEU:HD12	1.96	0.47
2:D:600:STI:O29	2:D:600:STI:H151	2.15	0.47
1:B:343:MET:HB2	1:B:343:MET:HE2	1.61	0.46
1:C:272:THR:HA	1:C:311:PHE:O	2.16	0.46
1:A:388:MET:HG3	1:A:389:THR:H	1.81	0.46
1:D:295:HIS:HD2	1:D:297:ASN:N	2.00	0.46
1:D:306:THR:HA	1:D:311:PHE:HD2	1.80	0.46
1:A:235:TRP:CD1	1:A:291:LYS:HG2	2.50	0.45
1:D:289:VAL:HG13	1:D:359:PHE:CE2	2.50	0.45
1:A:237:MET:CE	1:A:242:ILE:HD12	2.46	0.45
1:A:305:CYS:HB2	1:A:312:TYR:HB2	1.97	0.45
1:C:246:HIS:H	1:C:246:HIS:CD2	2.34	0.45
1:D:485:SER:OG	1:D:488:GLU:HG3	2.17	0.45
1:C:376:LEU:HD21	1:C:378:LYS:HE2	1.99	0.45
1:D:327:LEU:HD21	1:D:343:MET:HE1	1.98	0.45
1:B:283:PHE:CE2	1:B:311:PHE:HB3	2.52	0.44
1:B:440:TYR:N	1:B:441:PRO:HD3	2.32	0.44
1:C:249:GLY:O	1:C:250:GLY:C	2.56	0.44
1:D:248:LEU:HD12	1:D:256:VAL:HG12	1.99	0.43
1:C:286:GLU:CG	1:C:290:MET:CE	2.80	0.43
1:C:419:LYS:NZ	3:C:184:HOH:O	2.50	0.43
1:C:361:HIS:O	1:C:362:ARG:HB2	2.18	0.43
1:C:326:TYR:O	1:C:330:CYS:HB3	2.19	0.43
1:B:286:GLU:OE2	2:B:600:STI:H171	2.19	0.43
1:A:237:MET:HE1	1:A:242:ILE:CD1	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:GLY:O	1:A:252:GLN:HB3	2.18	0.42
1:D:460:ARG:HG3	1:D:469:TYR:HB2	2.01	0.42
1:D:414:ASN:HD22	1:D:414:ASN:C	2.21	0.42
1:C:366:ALA:N	1:C:428:LEU:HD13	2.34	0.42
1:D:479:ASN:HD22	1:D:480:PRO:CD	2.25	0.42
1:C:309:PRO:HG3	1:C:311:PHE:HE1	1.83	0.42
1:D:271:LYS:HB2	2:D:600:STI:H181	2.00	0.42
1:B:261:TRP:CZ2	1:B:264:TYR:HE1	2.38	0.42
1:C:406:THR:HG22	1:C:410:SER:HB2	2.01	0.42
1:A:412:ALA:HA	1:B:397:ALA:CB	2.49	0.42
1:C:387:LEU:HG	1:C:387:LEU:H	1.75	0.42
1:D:318:MET:N	2:D:600:STI:N3	2.52	0.42
1:B:382:PHE:HB2	1:B:384:LEU:HG	2.01	0.42
1:D:322:ASN:HA	1:D:369:CYS:O	2.20	0.41
1:A:460:ARG:HG3	1:A:469:TYR:HB2	2.02	0.41
1:B:305:CYS:O	1:B:311:PHE:HA	2.21	0.41
1:C:274:LYS:O	1:C:275:GLU:C	2.58	0.41
1:B:261:TRP:CZ2	1:B:264:TYR:CE1	3.08	0.41
1:A:445:LEU:HD23	1:A:445:LEU:HA	1.87	0.41
1:B:274:LYS:O	1:B:275:GLU:C	2.58	0.41
1:D:290:MET:HE1	1:D:313:ILE:HG21	2.01	0.41
1:D:318:MET:CE	1:D:318:MET:HA	2.50	0.41
1:D:322:ASN:HD22	1:D:324:LEU:H	1.68	0.41
1:D:427:VAL:O	1:D:431:GLU:HG3	2.21	0.41
1:D:318:MET:HG3	1:D:370:LEU:HD13	2.03	0.41
1:D:471:LEU:HD12	1:D:471:LEU:HA	1.90	0.41
1:B:273:LEU:HD21	1:B:280:VAL:HA	2.03	0.41
1:C:249:GLY:O	1:C:251:GLY:N	2.54	0.41
1:D:306:THR:HA	1:D:311:PHE:CD2	2.56	0.41
1:A:327:LEU:CD2	1:A:343:MET:HE1	2.47	0.41
1:A:361:HIS:O	1:A:362:ARG:CB	2.69	0.41
1:D:343:MET:HB2	1:D:343:MET:HE2	1.92	0.41
1:D:395:ALA:HB2	1:D:400:LYS:HD3	2.02	0.41
1:D:440:TYR:N	1:D:441:PRO:CD	2.84	0.40
2:B:600:STI:O29	2:B:600:STI:H151	2.21	0.40
1:C:342:TYR:O	1:C:346:GLN:HG3	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:265:SER:OG	1:B:374:ASN:ND2[4_566]	2.02	0.18

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	259/273 (95%)	237 (92%)	20 (8%)	2 (1%)	19	29
1	B	261/273 (96%)	251 (96%)	10 (4%)	0	100	100
1	C	256/273 (94%)	242 (94%)	13 (5%)	1 (0%)	34	48
1	D	260/273 (95%)	243 (94%)	16 (6%)	1 (0%)	34	48
All	All	1036/1092 (95%)	973 (94%)	59 (6%)	4 (0%)	34	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	388	MET
1	C	274	LYS
1	A	497	PHE
1	D	497	PHE

5.3.2 Protein sidechains [i](#)

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	211/239 (88%)	200 (95%)	11 (5%)	23	38
1	B	216/239 (90%)	211 (98%)	5 (2%)	50	70

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	214/239 (90%)	204 (95%)	10 (5%)	26	42
1	D	216/239 (90%)	206 (95%)	10 (5%)	27	43
All	All	857/956 (90%)	821 (96%)	36 (4%)	30	47

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	242	ILE
1	A	243	THR
1	A	267	THR
1	A	276	ASP
1	A	293	ILE
1	A	322	ASN
1	A	358	ASN
1	A	373	GLU
1	A	385	SER
1	A	418	ILE
1	A	455	ASP
1	B	262	LYS
1	B	293	ILE
1	B	373	GLU
1	B	471	LEU
1	B	479	ASN
1	C	240	THR
1	C	247	LYS
1	C	293	ILE
1	C	373	GLU
1	C	378	LYS
1	C	414	ASN
1	C	418	ILE
1	C	444	ASP
1	C	471	LEU
1	C	479	ASN
1	D	240	THR
1	D	242	ILE
1	D	322	ASN
1	D	338	VAL
1	D	357	LYS
1	D	373	GLU
1	D	378	LYS
1	D	414	ASN

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Mol	Chain	Res	Type
1	D	418	ILE
1	D	428	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	295	HIS
1	A	322	ASN
1	A	479	ASN
1	A	490	HIS
1	B	295	HIS
1	B	322	ASN
1	B	479	ASN
1	C	246	HIS
1	C	375	HIS
1	C	479	ASN
1	C	490	HIS
1	D	252	GLN
1	D	295	HIS
1	D	322	ASN
1	D	414	ASN
1	D	479	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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
2	STI	D	600	-	41,41,41	1.29	7 (17%)	56,56,56	1.85	11 (19%)
2	STI	A	600	-	41,41,41	1.42	7 (17%)	56,56,56	2.02	17 (30%)
2	STI	C	600	-	41,41,41	1.27	7 (17%)	56,56,56	2.21	20 (35%)
2	STI	B	600	-	41,41,41	1.49	8 (19%)	56,56,56	1.89	15 (26%)

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
2	STI	D	600	-	-	1/20/30/30	0/5/5/5
2	STI	A	600	-	-	1/20/30/30	0/5/5/5
2	STI	C	600	-	-	2/20/30/30	0/5/5/5
2	STI	B	600	-	-	3/20/30/30	0/5/5/5

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	600	STI	C29-C28	3.27	1.44	1.38
2	A	600	STI	C26-C25	2.96	1.44	1.38
2	D	600	STI	C9-N10	-2.92	1.30	1.34
2	A	600	STI	C5-C7	2.89	1.53	1.48
2	B	600	STI	C26-C25	2.76	1.43	1.38
2	C	600	STI	C4-N3	2.73	1.40	1.34
2	A	600	STI	C29-C28	2.69	1.43	1.38
2	A	600	STI	C29-C23	2.66	1.43	1.39
2	B	600	STI	C15-C14	2.66	1.43	1.39
2	B	600	STI	C5-C7	2.52	1.52	1.48
2	A	600	STI	C17-C18	2.51	1.43	1.38
2	D	600	STI	C52-C53	2.46	1.60	1.51
2	B	600	STI	C52-C53	2.41	1.60	1.51
2	D	600	STI	C5-C7	2.40	1.52	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	600	STI	C25-C23	2.36	1.43	1.39
2	C	600	STI	C16-N21	2.19	1.46	1.41
2	A	600	STI	C52-N51	2.18	1.51	1.46
2	D	600	STI	C29-C28	2.12	1.42	1.38
2	D	600	STI	C12-C7	-2.11	1.34	1.38
2	C	600	STI	C12-C7	-2.09	1.34	1.38
2	A	600	STI	C12-C7	-2.07	1.34	1.38
2	D	600	STI	C29-C23	2.07	1.42	1.39
2	D	600	STI	C23-C22	-2.06	1.45	1.50
2	C	600	STI	C29-C28	2.05	1.42	1.38
2	B	600	STI	C1-C6	2.05	1.43	1.38
2	C	600	STI	C6-C5	2.03	1.43	1.39
2	C	600	STI	C26-C25	2.02	1.42	1.38
2	B	600	STI	C23-C22	-2.02	1.46	1.50
2	C	600	STI	C15-C14	2.01	1.42	1.39

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	600	STI	C11-N10-C9	6.14	120.90	115.45
2	D	600	STI	N10-C9-N8	-6.12	120.75	126.55
2	C	600	STI	C11-N10-C9	6.03	120.80	115.45
2	A	600	STI	N10-C9-N8	-5.95	120.91	126.55
2	D	600	STI	C11-N10-C9	5.75	120.55	115.45
2	C	600	STI	C54-N51-C52	-5.74	102.07	110.66
2	A	600	STI	C50-N51-C52	5.18	116.77	109.52
2	C	600	STI	N10-C9-N8	-5.01	121.80	126.55
2	B	600	STI	C11-N10-C9	4.27	119.24	115.45
2	B	600	STI	C50-N51-C52	4.25	115.47	109.52
2	B	600	STI	N10-C9-N8	-4.22	122.55	126.55
2	B	600	STI	C53-C52-N51	-4.13	106.14	110.80
2	C	600	STI	C20-C19-C14	4.07	125.86	121.25
2	C	600	STI	C50-C49-N48	-4.01	102.42	110.64
2	B	600	STI	C14-N13-C9	-3.80	118.13	129.60
2	A	600	STI	C14-N13-C9	-3.69	118.47	129.60
2	C	600	STI	C14-N13-C9	-3.62	118.69	129.60
2	D	600	STI	C7-N8-C9	3.60	119.66	116.69
2	C	600	STI	C12-C11-N10	-3.37	119.77	123.96
2	C	600	STI	C17-C16-C15	-3.33	115.70	119.65
2	A	600	STI	C7-N8-C9	3.14	119.28	116.69
2	C	600	STI	C19-C14-N13	3.14	124.65	118.70
2	B	600	STI	N13-C9-N10	3.10	125.89	116.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	600	STI	C49-N48-C53	2.96	115.49	108.83
2	C	600	STI	C23-C22-N21	2.93	122.38	115.92
2	D	600	STI	N13-C9-N10	2.91	125.33	116.28
2	B	600	STI	C49-N48-C53	2.91	115.38	108.83
2	B	600	STI	C12-C11-N10	-2.89	120.36	123.96
2	A	600	STI	C12-C11-N10	-2.88	120.38	123.96
2	D	600	STI	O29-C22-C23	-2.85	115.85	120.94
2	C	600	STI	C49-C50-N51	-2.77	107.67	110.80
2	C	600	STI	C4-C5-C7	-2.75	116.79	121.22
2	A	600	STI	N13-C9-N10	2.74	124.80	116.28
2	A	600	STI	C49-C50-N51	-2.73	107.72	110.80
2	B	600	STI	C18-C17-C16	-2.71	117.17	120.30
2	A	600	STI	C20-C19-C14	2.67	124.28	121.25
2	C	600	STI	O29-C22-C23	-2.67	116.18	120.94
2	C	600	STI	N13-C9-N10	2.60	124.36	116.28
2	A	600	STI	C54-N51-C52	-2.59	106.78	110.66
2	C	600	STI	C53-C52-N51	-2.47	108.02	110.80
2	B	600	STI	O29-C22-C23	-2.46	116.55	120.94
2	D	600	STI	C19-C14-N13	2.46	123.35	118.70
2	C	600	STI	C15-C14-N13	-2.44	116.00	121.05
2	D	600	STI	C11-C12-C7	-2.43	114.93	117.22
2	B	600	STI	C54-N51-C50	-2.39	107.08	110.66
2	B	600	STI	C49-C50-N51	-2.37	108.12	110.80
2	D	600	STI	C6-C1-C2	-2.36	115.42	118.91
2	D	600	STI	C50-C49-N48	-2.32	105.89	110.64
2	A	600	STI	C28-C29-C23	-2.31	118.09	120.78
2	A	600	STI	C15-C16-N21	2.30	127.69	120.18
2	C	600	STI	C15-C16-N21	2.30	127.69	120.18
2	A	600	STI	C6-C1-C2	-2.28	115.53	118.91
2	B	600	STI	C17-C16-N21	-2.24	112.88	120.40
2	A	600	STI	C16-N21-C22	-2.15	121.01	126.58
2	B	600	STI	C15-C16-N21	2.14	127.17	120.18
2	C	600	STI	C16-C15-C14	2.14	124.40	120.17
2	D	600	STI	C26-C25-C23	-2.13	118.30	120.78
2	C	600	STI	C26-C25-C23	2.10	123.22	120.78
2	C	600	STI	C12-C7-N8	2.08	124.67	121.97
2	A	600	STI	C53-C52-N51	-2.07	108.46	110.80
2	A	600	STI	C25-C26-C27	-2.02	118.25	121.03
2	B	600	STI	C54-N51-C52	-2.01	107.65	110.66
2	D	600	STI	C12-C11-N10	-2.01	121.46	123.96

There are no chirality outliers.

All (7) torsion outliers are listed below:

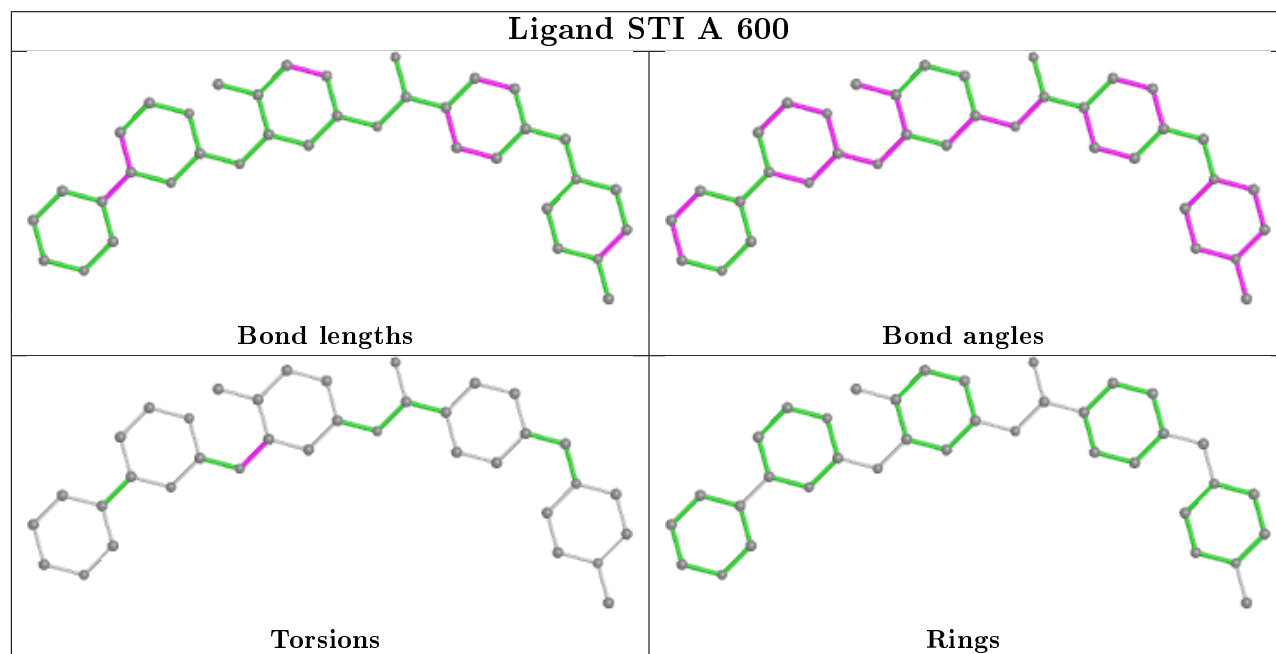
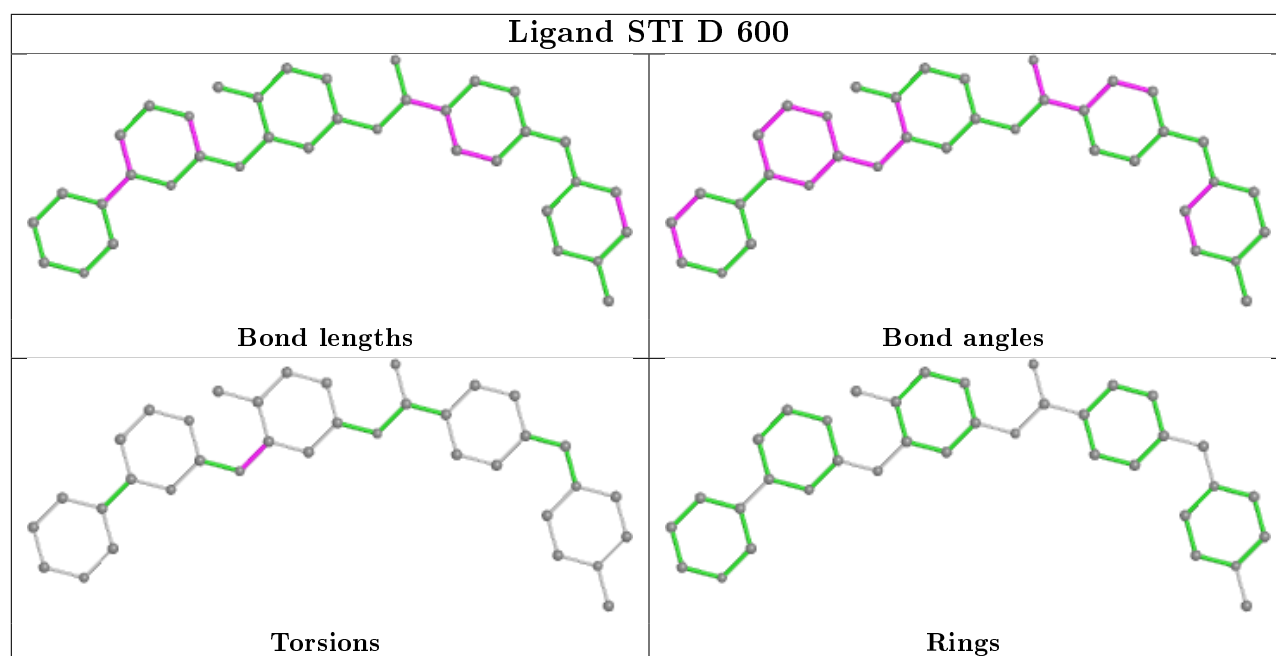
Mol	Chain	Res	Type	Atoms
2	D	600	STI	C15-C14-N13-C9
2	C	600	STI	C15-C14-N13-C9
2	B	600	STI	C15-C14-N13-C9
2	B	600	STI	C19-C14-N13-C9
2	B	600	STI	O29-C22-N21-C16
2	A	600	STI	C15-C14-N13-C9
2	C	600	STI	O29-C22-N21-C16

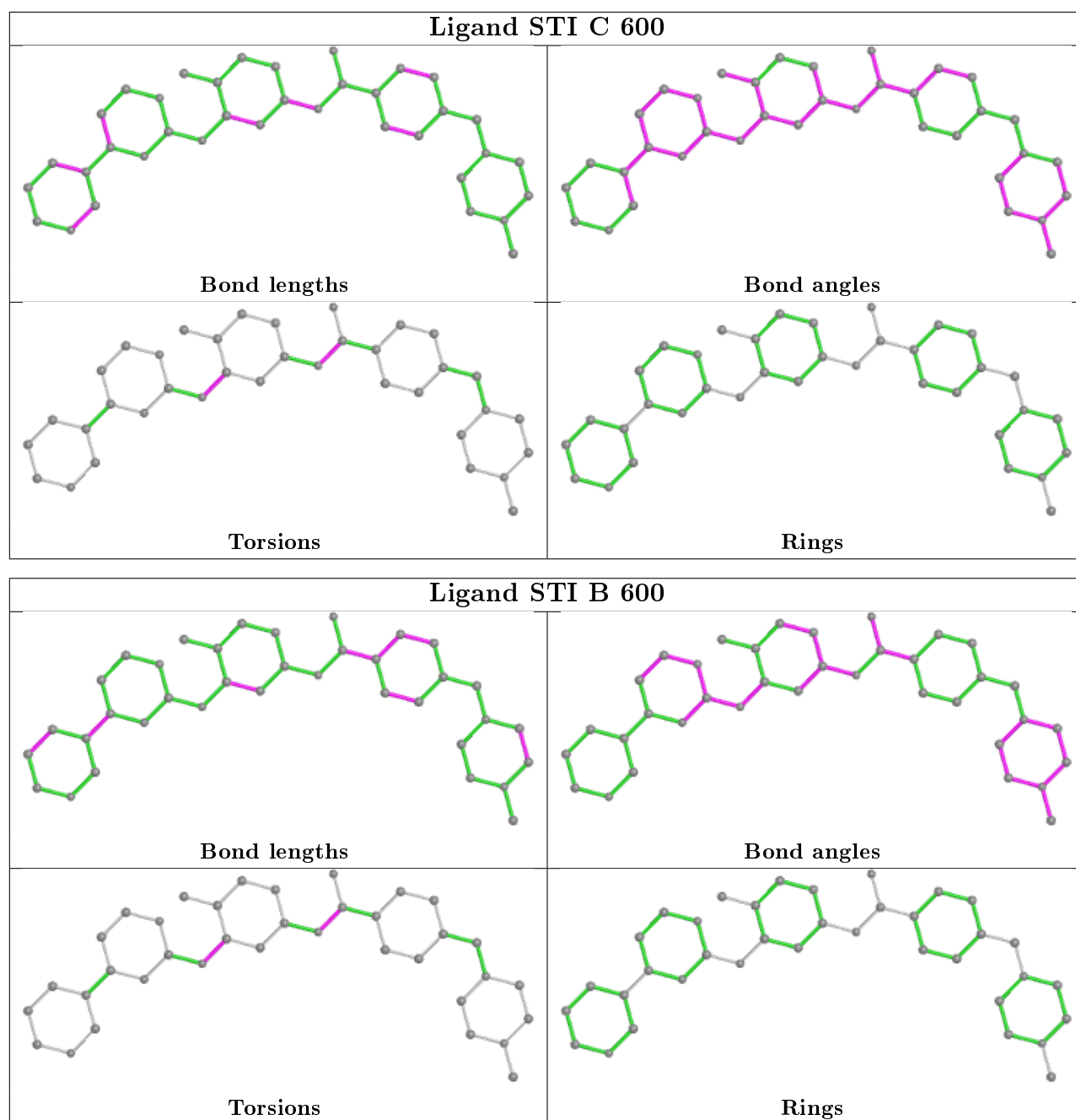
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	600	STI	4	0
2	A	600	STI	1	0
2	B	600	STI	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.





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	263/273 (96%)	0.28	26 (9%) 7 6	22, 40, 86, 100	0
1	B	265/273 (97%)	0.11	17 (6%) 19 18	20, 37, 63, 93	0
1	C	262/273 (95%)	0.09	16 (6%) 21 20	22, 36, 63, 85	0
1	D	264/273 (96%)	0.27	22 (8%) 11 10	23, 40, 76, 83	0
All	All	1054/1092 (96%)	0.19	81 (7%) 13 12	20, 38, 76, 100	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	278	MET	6.0
1	A	277	THR	5.5
1	A	389	THR	5.3
1	A	265	SER	5.1
1	B	278	MET	5.1
1	D	387	LEU	4.3
1	B	389	THR	4.2
1	D	389	THR	3.8
1	A	238	GLU	3.8
1	A	278	MET	3.7
1	D	391	ASP	3.5
1	D	309	PRO	3.5
1	D	264	TYR	3.5
1	C	246	HIS	3.5
1	D	424	ALA	3.4
1	A	427	VAL	3.3
1	A	307	ARG	3.3
1	C	388	MET	3.3
1	B	427	VAL	3.2
1	B	277	THR	3.2
1	D	307	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	C	497	PHE	3.1
1	A	270	VAL	3.1
1	D	279	GLU	3.1
1	C	233	ASP	3.0
1	A	279	GLU	3.0
1	D	497	PHE	3.0
1	A	264	TYR	3.0
1	D	386	ARG	3.0
1	D	388	MET	3.0
1	C	428	LEU	2.9
1	A	364	LEU	2.9
1	B	246	HIS	2.9
1	A	428	LEU	2.9
1	C	364	LEU	2.8
1	A	266	LEU	2.7
1	C	281	GLU	2.7
1	C	277	THR	2.7
1	A	425	PHE	2.7
1	B	425	PHE	2.7
1	D	427	VAL	2.6
1	A	240	THR	2.6
1	A	310	PRO	2.6
1	D	428	LEU	2.6
1	D	425	PHE	2.6
1	D	315	THR	2.5
1	D	311	PHE	2.5
1	D	240	THR	2.5
1	A	398	GLY	2.5
1	A	276	ASP	2.5
1	B	265	SER	2.5
1	B	275	GLU	2.5
1	A	313	ILE	2.5
1	A	261	TRP	2.4
1	A	251	GLY	2.4
1	B	424	ALA	2.4
1	B	313	ILE	2.4
1	D	364	LEU	2.4
1	B	497	PHE	2.4
1	D	234	LYS	2.4
1	C	347	ILE	2.4
1	D	250	GLY	2.3
1	A	269	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	428	LEU	2.3
1	B	281	GLU	2.3
1	A	281	GLU	2.3
1	B	314	ILE	2.3
1	C	313	ILE	2.2
1	B	498	GLN	2.2
1	C	399	ALA	2.1
1	B	279	GLU	2.1
1	A	257	TYR	2.1
1	B	280	VAL	2.1
1	C	498	GLN	2.1
1	A	424	ALA	2.1
1	D	280	VAL	2.1
1	A	311	PHE	2.1
1	C	369	CYS	2.1
1	C	275	GLU	2.0
1	C	306	THR	2.0
1	D	262	LYS	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 carbohydrates in this entry.

6.4 Ligands [i](#)

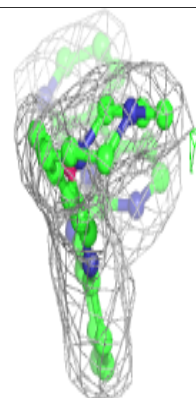
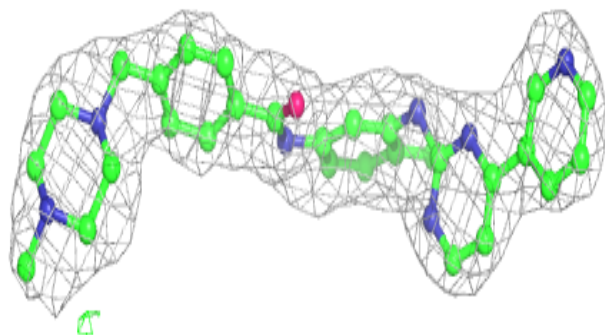
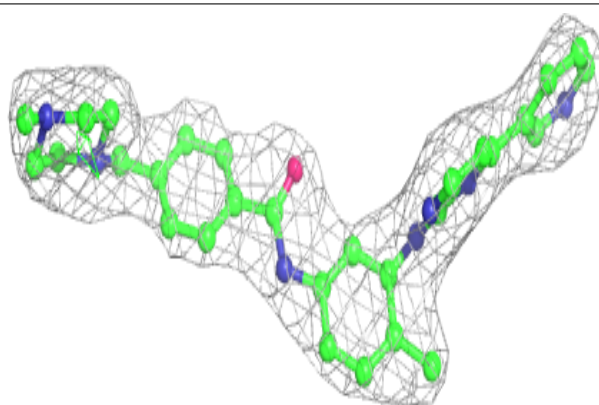
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	STI	A	600	37/37	0.93	0.16	26,38,42,45	0
2	STI	D	600	37/37	0.94	0.15	24,30,38,39	0
2	STI	C	600	37/37	0.94	0.19	18,26,32,33	0
2	STI	B	600	37/37	0.96	0.17	22,27,30,31	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.

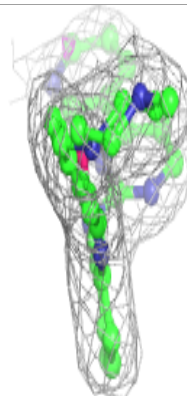
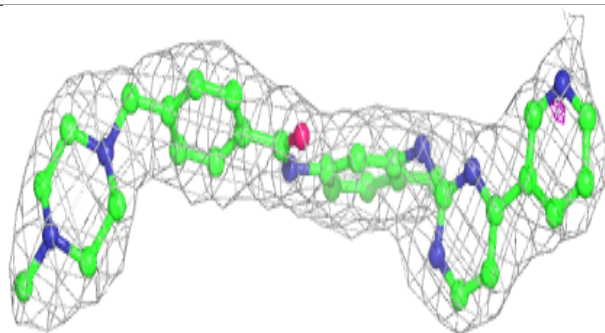
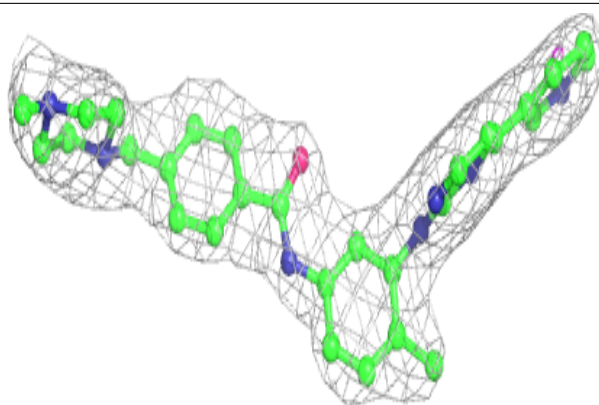
Electron density around STI A 600:

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

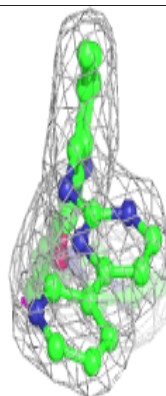
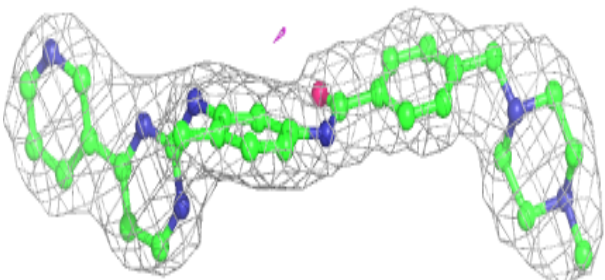
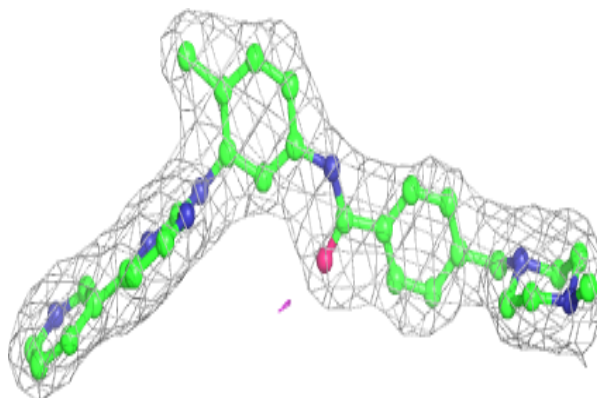


Electron density around STI D 600:

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

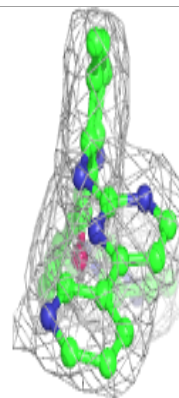
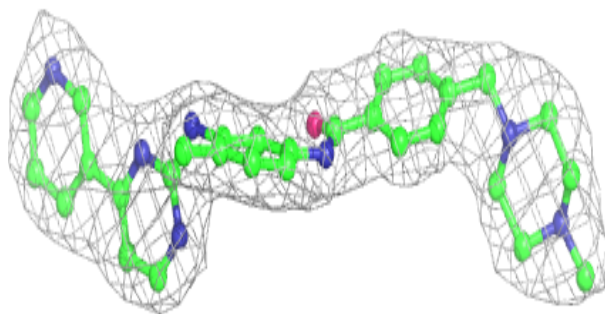
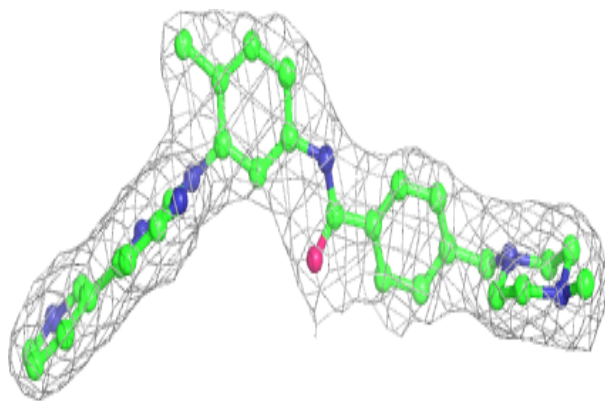
**Electron density around STI C 600:**

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



Electron density around STI B 600:

$2mF_o-DF_c$ (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.