



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:50 PM UTC

PDB ID : 2QQ7 / pdb_00002qq7
Title : Crystal structure of drug resistant SRC kinase domain with irreversible inhibitor
Authors : Michalczyk, A.; Rode, H.B.; Gruetter, C.; Rauh, D.
Deposited on : 2007-07-26
Resolution : 2.38 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

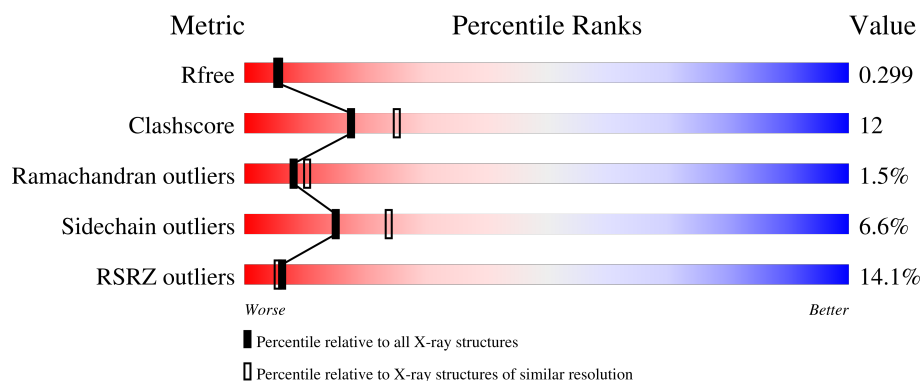
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	286	
1	B	286	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4160 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.

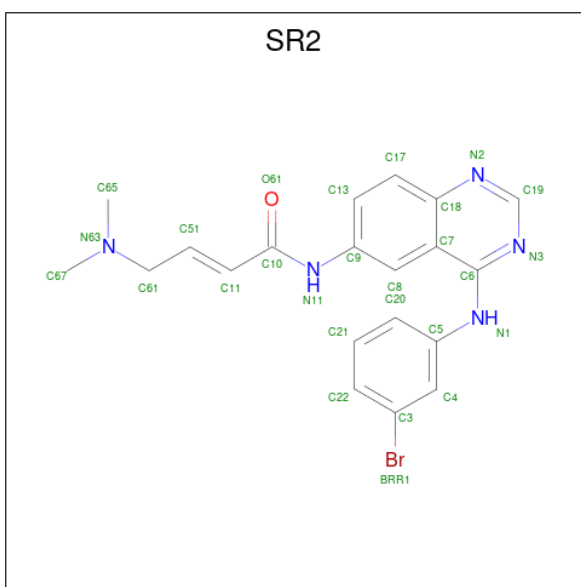
- Molecule 1 is a protein called Proto-oncogene tyrosine-protein kinase Src.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	1	2	0
			1988	1278	334	359	17			
1	B	244	Total	C	N	O	S	0	0	0
			1950	1252	326	355	17			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	248	GLY	-	expression tag	UNP P00523
A	249	HIS	-	expression tag	UNP P00523
A	250	MET	-	expression tag	UNP P00523
A	338	MET	THR	engineered mutation	UNP P00523
A	345	CYS	SER	engineered mutation	UNP P00523
B	248	GLY	-	expression tag	UNP P00523
B	249	HIS	-	expression tag	UNP P00523
B	250	MET	-	expression tag	UNP P00523
B	338	MET	THR	engineered mutation	UNP P00523
B	345	CYS	SER	engineered mutation	UNP P00523

- Molecule 2 is (2E)-N-{4-[(3-bromophenyl)amino]quinazolin-6-yl}-4-(dimethylamino)but-2-enamide (CCD ID: SR2) (formula: C₂₀H₂₀BrN₅O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	Br	C	N	O	0	0
			27	1	20	5	1		
2	B	1	Total	Br	C	N	O	0	0
			27	1	20	5	1		

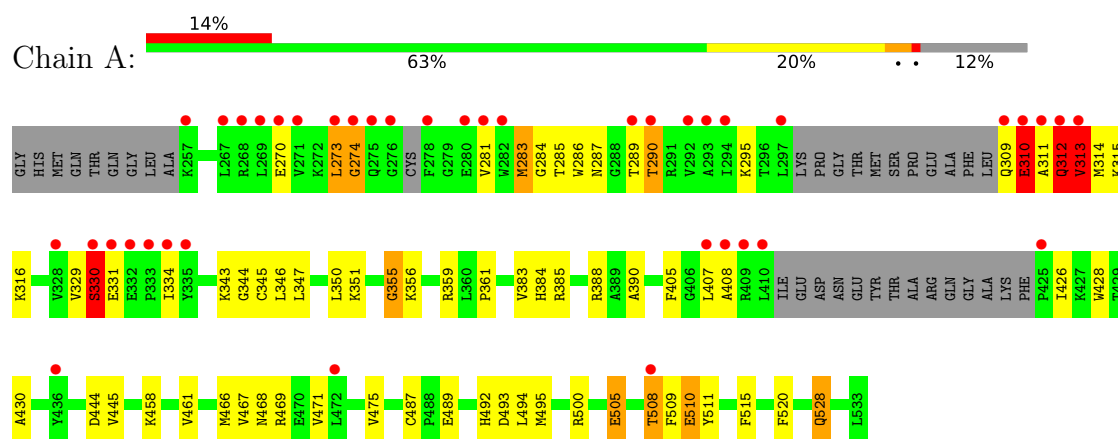
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	93	Total	O	0	0
			93	93		
3	B	75	Total	O	0	0
			75	75		

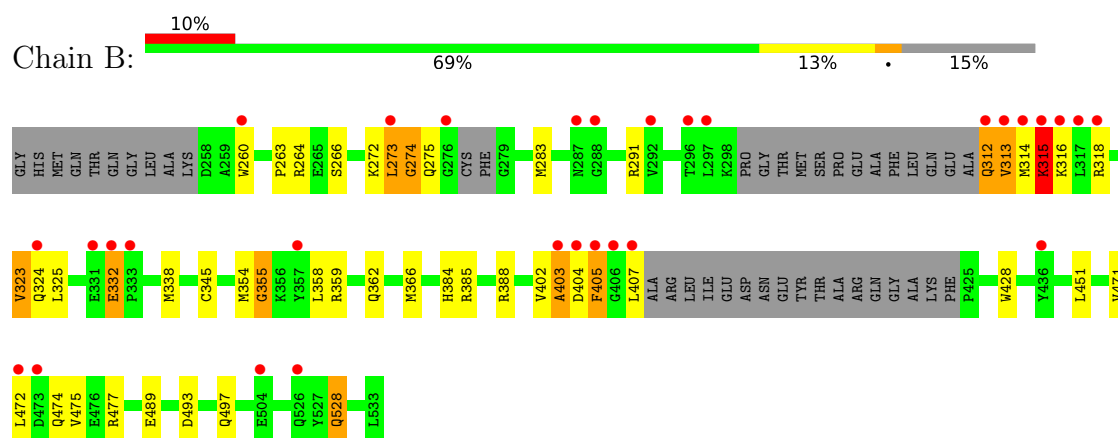
3 Residue-property plots [i](#)

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: Proto-oncogene tyrosine-protein kinase Src



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	42.22Å 63.36Å 75.06Å 78.58° 89.08° 89.87°	Depositor
Resolution (Å)	73.52 – 2.38 73.52 – 2.38	Depositor EDS
% Data completeness (in resolution range)	98.0 (73.52-2.38) 98.0 (73.52-2.38)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	13.44 (at 2.37Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.214 , 0.266 (Not available) , 0.299	Depositor DCC
R_{free} test set	1518 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.178	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.034 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4160	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.64% 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: SR2

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	1.07	2/2040 (0.1%)	1.19	16/2762 (0.6%)
1	B	0.99	0/1995	1.11	6/2699 (0.2%)
All	All	1.03	2/4035 (0.0%)	1.15	22/5461 (0.4%)

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	5
All	All	0	8

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	445	VAL	CA-CB	7.15	1.62	1.54
1	A	383	VAL	C-O	-6.05	1.18	1.24

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	430	ALA	CA-C-N	9.64	129.26	119.24
1	A	430	ALA	C-N-CA	9.64	129.26	119.24
1	B	528	GLN	CA-C-N	6.38	126.40	119.89
1	B	528	GLN	C-N-CA	6.38	126.40	119.89
1	A	461	VAL	CA-C-N	6.06	126.40	119.92
1	A	461	VAL	C-N-CA	6.06	126.40	119.92
1	A	313	VAL	CB-CA-C	6.01	121.14	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	273	LEU	N-CA-C	5.88	120.24	112.72
1	A	528	GLN	CA-C-N	5.73	125.68	119.78
1	A	528	GLN	C-N-CA	5.73	125.68	119.78
1	B	313	VAL	CA-C-N	5.49	131.58	121.70
1	B	313	VAL	C-N-CA	5.49	131.58	121.70
1	A	444	ASP	CA-C-N	5.47	127.66	120.60
1	A	444	ASP	C-N-CA	5.47	127.66	120.60
1	A	356	LYS	N-CA-C	5.41	117.94	111.71
1	A	346	LEU	N-CA-C	5.40	117.17	111.28
1	B	313	VAL	N-CA-C	5.40	120.57	109.34
1	A	487	CYS	CA-C-N	-5.33	114.47	119.85
1	A	487	CYS	C-N-CA	-5.33	114.47	119.85
1	A	355	GLY	CA-C-N	5.13	127.67	120.28
1	A	355	GLY	C-N-CA	5.13	127.67	120.28
1	B	332	GLU	N-CA-C	5.03	118.03	108.12

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	310	GLU	Peptide
1	A	312	GLN	Peptide
1	A	330	SER	Peptide
1	B	273	LEU	Peptide
1	B	315	LYS	Peptide
1	B	316	LYS	Peptide
1	B	402	VAL	Peptide
1	B	403	ALA	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	1988	0	1937	57	0
1	B	1950	0	1921	40	0
2	A	27	0	19	5	0
2	B	27	0	19	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	93	0	0	0	0
3	B	75	0	0	0	0
All	All	4160	0	3896	98	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 12.

All (98) 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:404:ASP:N	1:B:405:PHE:HA	1.45	1.25
1:B:403:ALA:HB1	1:B:404:ASP:HA	1.21	1.20
1:B:404:ASP:H	1:B:405:PHE:CA	1.56	1.19
1:A:312:GLN:HB3	1:A:313:VAL:HG23	1.32	1.12
1:B:403:ALA:CB	1:B:404:ASP:HA	1.97	0.94
1:A:508:THR:HG21	1:A:510:GLU:HG2	1.49	0.93
1:B:315:LYS:N	1:B:315:LYS:HD2	1.80	0.93
1:A:355:GLY:O	1:A:458:LYS:HE2	1.72	0.90
1:A:330:SER:CB	1:A:331:GLU:HB2	2.05	0.86
1:A:508:THR:CG2	1:A:510:GLU:H	1.89	0.85
1:B:404:ASP:N	1:B:405:PHE:CA	2.26	0.85
1:B:404:ASP:H	1:B:405:PHE:HA	0.71	0.84
1:A:311:ALA:O	1:A:312:GLN:HB2	1.75	0.83
1:A:309:GLN:O	1:A:313:VAL:HG21	1.79	0.81
1:B:358:LEU:O	1:B:359:ARG:NH1	2.14	0.80
1:A:508:THR:HG22	1:A:510:GLU:N	1.96	0.80
1:A:295:LYS:HB2	2:A:1345:SR2:BRR1	2.39	0.77
1:A:309:GLN:O	1:A:313:VAL:CG2	2.34	0.76
1:B:403:ALA:HB1	1:B:404:ASP:CA	2.11	0.76
1:B:313:VAL:H	1:B:314:MET:HB2	1.56	0.71
1:A:508:THR:HG21	1:A:510:GLU:CG	2.20	0.69
1:A:312:GLN:CB	1:A:313:VAL:HG23	2.18	0.69
1:A:345:CYS:SG	2:A:1345:SR2:H673	2.34	0.67
1:A:508:THR:CG2	1:A:510:GLU:N	2.57	0.67
1:B:272:LYS:HG2	1:B:274:GLY:HA2	1.77	0.66
1:A:385:ARG:HD2	1:A:407:LEU:O	1.95	0.66
1:A:508:THR:HG22	1:A:510:GLU:H	1.57	0.66
1:A:329:VAL:O	1:A:331:GLU:N	2.29	0.65
1:B:260:TRP:CD2	1:B:315:LYS:HE2	2.32	0.64
1:B:315:LYS:HD2	1:B:315:LYS:H	1.62	0.63
1:B:315:LYS:HG3	1:B:325:LEU:HD23	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:345:CYS:HB3	2:A:1345:SR2:O61	2.00	0.62
1:B:312:GLN:N	1:B:314:MET:O	2.35	0.60
1:A:283:MET:HG3	1:A:284:GLY:N	2.16	0.60
1:A:508:THR:HG23	1:A:510:GLU:H	1.65	0.60
1:B:273:LEU:H	1:B:274:GLY:HA2	1.66	0.60
1:A:273:LEU:H	1:A:274:GLY:HA3	1.68	0.59
1:A:329:VAL:O	1:A:330:SER:C	2.46	0.59
1:B:263:PRO:O	1:B:266:SER:OG	2.20	0.58
1:B:345:CYS:HB3	2:B:1345:SR2:O61	2.04	0.58
1:A:286:TRP:HB3	1:A:290:THR:HB	1.86	0.57
1:B:403:ALA:CB	1:B:404:ASP:CA	2.76	0.56
1:B:474:GLN:OE1	1:B:477:ARG:NH1	2.38	0.56
1:A:508:THR:HG22	1:A:511:TYR:H	1.70	0.56
1:B:264:ARG:HH22	1:B:332:GLU:H	1.56	0.54
1:A:500:ARG:HD3	1:A:505:GLU:HB3	1.90	0.54
1:B:260:TRP:CZ2	1:B:315:LYS:HG2	2.44	0.53
1:B:384:HIS:NE2	1:B:403:ALA:O	2.33	0.52
1:B:405:PHE:CD1	1:B:405:PHE:N	2.78	0.52
1:A:310:GLU:O	1:A:313:VAL:HB	2.08	0.51
1:B:314:MET:HE1	1:B:407:LEU:HD12	1.93	0.51
1:A:344:GLY:HA2	2:A:1345:SR2:C13	2.41	0.51
1:A:384:HIS:O	1:A:385:ARG:HB2	2.12	0.50
1:A:388[A]:ARG:NH1	1:A:390:ALA:HB3	2.26	0.50
1:B:362:GLN:O	1:B:366:MET:HG3	2.12	0.49
1:A:311:ALA:O	1:A:312:GLN:CB	2.54	0.49
1:A:508:THR:CG2	1:A:509:PHE:N	2.74	0.48
1:A:283:MET:HG3	1:A:284:GLY:H	1.79	0.48
1:B:264:ARG:NH2	1:B:332:GLU:H	2.11	0.48
2:B:1345:SR2:N3	2:B:1345:SR2:H4	2.29	0.47
1:B:323:VAL:HG22	1:B:338:MET:HE2	1.96	0.47
1:A:345:CYS:SG	2:A:1345:SR2:C67	3.01	0.47
1:A:285:THR:HG22	1:A:286:TRP:H	1.80	0.46
1:B:388:ARG:HB3	1:B:428:TRP:CD1	2.51	0.46
1:B:471:VAL:O	1:B:475:VAL:HG23	2.16	0.46
1:B:338:MET:HE1	1:B:405:PHE:CE2	2.51	0.46
1:B:404:ASP:C	1:B:405:PHE:CG	2.94	0.46
1:A:330:SER:O	1:A:334:ILE:HG12	2.16	0.45
1:A:361:PRO:HA	1:A:520:PHE:CE2	2.52	0.45
1:B:384:HIS:O	1:B:385:ARG:HB2	2.16	0.45
1:A:508:THR:CG2	1:A:510:GLU:CG	2.93	0.45
1:A:273:LEU:N	1:A:274:GLY:HA3	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:404:ASP:C	1:B:405:PHE:CD1	2.96	0.44
1:B:315:LYS:HB2	1:B:318:ARG:NH2	2.32	0.44
1:A:508:THR:HG23	1:A:509:PHE:N	2.33	0.44
1:B:493:ASP:O	1:B:497:GLN:HG3	2.18	0.44
1:A:274:GLY:HA2	1:A:281:VAL:HB	1.99	0.43
1:A:313:VAL:C	1:A:315:LYS:N	2.75	0.43
1:A:388[A]:ARG:HB3	1:A:428:TRP:CD1	2.53	0.43
1:A:313:VAL:O	1:A:315:LYS:N	2.52	0.43
1:B:260:TRP:CG	1:B:315:LYS:HE2	2.54	0.42
1:B:283:MET:SD	1:B:291:ARG:NH1	2.93	0.42
1:A:330:SER:CB	1:A:331:GLU:CB	2.89	0.42
1:A:385:ARG:CD	1:A:407:LEU:O	2.65	0.42
1:A:494:LEU:HD22	1:A:515:PHE:CE1	2.55	0.42
1:A:510:GLU:HG3	1:A:511:TYR:N	2.34	0.41
1:A:313:VAL:C	1:A:315:LYS:H	2.27	0.41
1:A:467:VAL:HG12	1:A:468:ASN:H	1.85	0.41
1:A:492:HIS:O	1:A:495:MET:HB2	2.20	0.41
1:A:350:LEU:O	1:A:355:GLY:HA3	2.20	0.41
1:B:315:LYS:N	1:B:315:LYS:CD	2.64	0.41
1:A:471:VAL:O	1:A:475:VAL:HG23	2.20	0.41
1:A:405:PHE:HD2	1:A:408:ALA:CB	2.33	0.41
1:B:354:MET:O	1:B:355:GLY:C	2.62	0.41
1:A:343:LYS:HA	1:A:343:LYS:HD3	1.85	0.41
1:A:466:MET:HE2	1:A:466:MET:HB3	1.95	0.41
1:A:329:VAL:HG12	1:A:331:GLU:HB3	2.02	0.40
1:A:287:ASN:C	1:A:289:THR:H	2.29	0.40

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	245/286 (86%)	229 (94%)	11 (4%)	5 (2%)	6	6
1	B	236/286 (82%)	219 (93%)	15 (6%)	2 (1%)	16	23
All	All	481/572 (84%)	448 (93%)	26 (5%)	7 (2%)	8	10

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	312	GLN
1	B	355	GLY
1	A	313	VAL
1	A	330	SER
1	A	274	GLY
1	A	314	MET
1	B	274	GLY

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	205/245 (84%)	187 (91%)	18 (9%)	9	14
1	B	207/245 (84%)	197 (95%)	10 (5%)	23	37
All	All	412/490 (84%)	384 (93%)	28 (7%)	15	23

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

Mol	Chain	Res	Type
1	A	270	GLU
1	A	283	MET
1	A	290	THR
1	A	310	GLU
1	A	313	VAL
1	A	316	LYS
1	A	347	LEU
1	A	351	LYS
1	A	359	ARG

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Mol	Chain	Res	Type
1	A	426	ILE
1	A	469	ARG
1	A	489	GLU
1	A	493[A]	ASP
1	A	493[B]	ASP
1	A	505	GLU
1	A	508	THR
1	A	510	GLU
1	A	528	GLN
1	B	275	GLN
1	B	312	GLN
1	B	315	LYS
1	B	323	VAL
1	B	324	GLN
1	B	405	PHE
1	B	451	LEU
1	B	472	LEU
1	B	489	GLU
1	B	528	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	275	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SR2	B	1345	1	28,29,29	2.51	6 (21%)	38,39,39	2.44	11 (28%)
2	SR2	A	1345	1	28,29,29	2.30	6 (21%)	38,39,39	2.57	12 (31%)

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	SR2	B	1345	1	-	9/14/14/14	0/3/3/3
2	SR2	A	1345	1	-	1/14/14/14	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1345	SR2	C11-C51	10.37	1.60	1.32
2	A	1345	SR2	C11-C51	8.75	1.56	1.32
2	A	1345	SR2	C6-C7	-3.78	1.40	1.44
2	B	1345	SR2	C11-C10	3.62	1.55	1.48
2	A	1345	SR2	C9-N11	-3.44	1.34	1.41
2	A	1345	SR2	C5-N1	-3.40	1.33	1.40
2	B	1345	SR2	C9-N11	-3.23	1.35	1.41
2	A	1345	SR2	C11-C10	2.84	1.54	1.48
2	B	1345	SR2	C6-C7	-2.84	1.41	1.44
2	B	1345	SR2	C5-N1	-2.59	1.35	1.40
2	B	1345	SR2	C61-C51	2.52	1.56	1.49
2	A	1345	SR2	C61-C51	2.20	1.55	1.49

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1345	SR2	C61-C51-C11	-7.55	113.11	124.82
2	A	1345	SR2	C19-N3-C6	6.72	121.81	116.60
2	B	1345	SR2	C61-C51-C11	-6.64	114.53	124.82
2	B	1345	SR2	C19-N3-C6	6.44	121.60	116.60
2	A	1345	SR2	N2-C19-N3	-5.81	120.78	128.67
2	B	1345	SR2	N2-C19-N3	-5.48	121.24	128.67
2	B	1345	SR2	C19-N2-C18	4.76	120.84	115.43
2	A	1345	SR2	C19-N2-C18	4.56	120.61	115.43
2	B	1345	SR2	C7-C6-N3	-4.13	118.29	121.35
2	A	1345	SR2	C8-C7-C6	-4.10	121.51	124.84
2	B	1345	SR2	C6-C7-C18	3.72	118.06	115.86
2	A	1345	SR2	N1-C6-N3	3.38	122.52	118.71
2	A	1345	SR2	C7-C6-N3	-3.32	118.89	121.35
2	B	1345	SR2	C7-C18-N2	-2.70	119.96	122.82
2	B	1345	SR2	C5-C4-C3	2.69	121.29	118.66
2	B	1345	SR2	C9-N11-C10	-2.63	123.12	128.07
2	B	1345	SR2	C8-C7-C6	-2.47	122.83	124.84
2	A	1345	SR2	C6-C7-C18	2.40	117.28	115.86
2	B	1345	SR2	N1-C6-N3	2.38	121.39	118.71
2	A	1345	SR2	C51-C61-N63	-2.25	108.80	112.25
2	A	1345	SR2	C7-C18-N2	-2.08	120.62	122.82
2	A	1345	SR2	BRR1-C3-C22	2.05	122.09	119.30
2	A	1345	SR2	C51-C11-C10	-2.00	117.85	122.94

There are no chirality outliers.

All (10) torsion outliers are listed below:

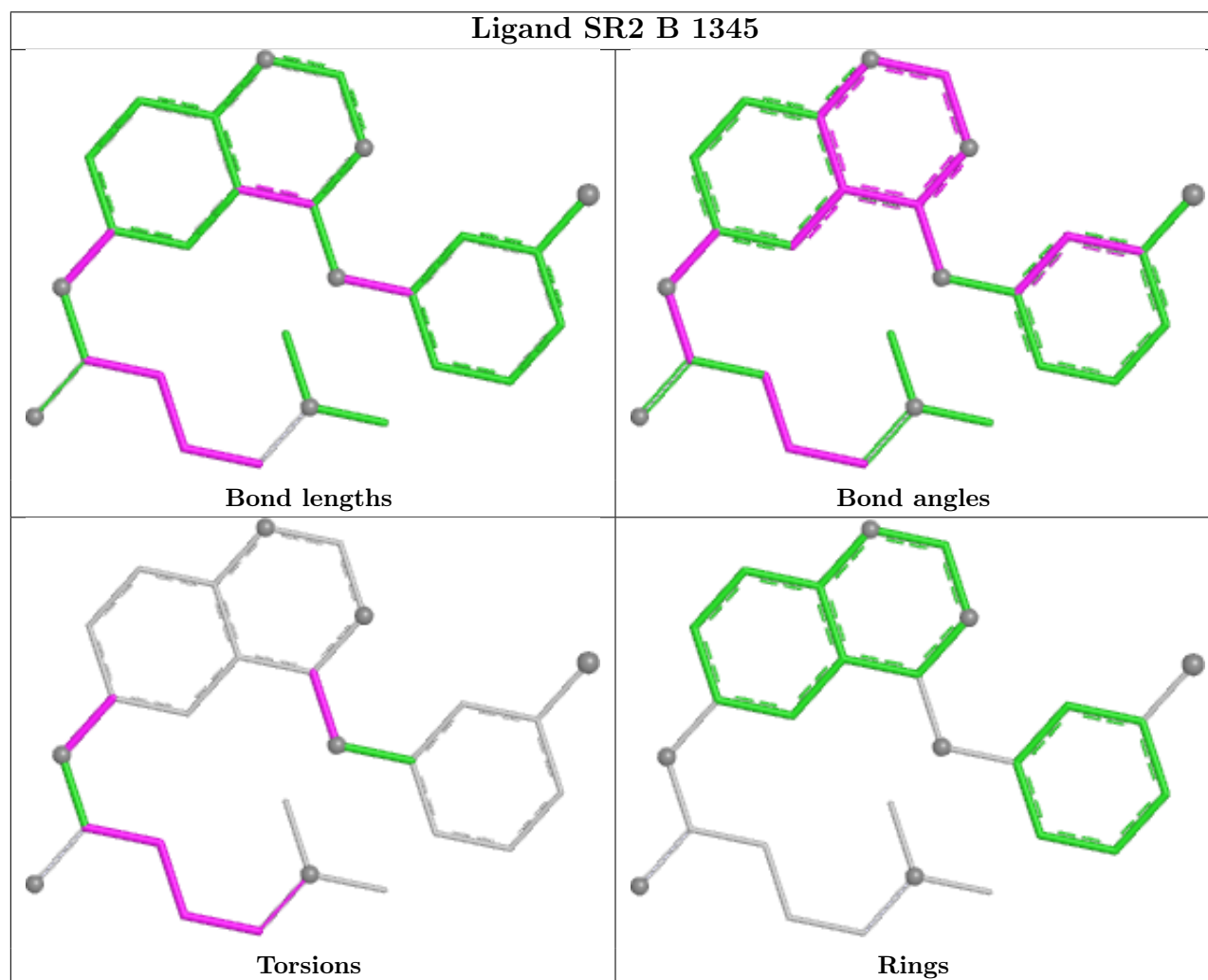
Mol	Chain	Res	Type	Atoms
2	A	1345	SR2	C10-C11-C51-C61
2	B	1345	SR2	C7-C6-N1-C5
2	B	1345	SR2	N3-C6-N1-C5
2	B	1345	SR2	C10-C11-C51-C61
2	B	1345	SR2	O61-C10-C11-C51
2	B	1345	SR2	N11-C10-C11-C51
2	B	1345	SR2	C51-C61-N63-C65
2	B	1345	SR2	C51-C61-N63-C67
2	B	1345	SR2	C11-C51-C61-N63
2	B	1345	SR2	C13-C9-N11-C10

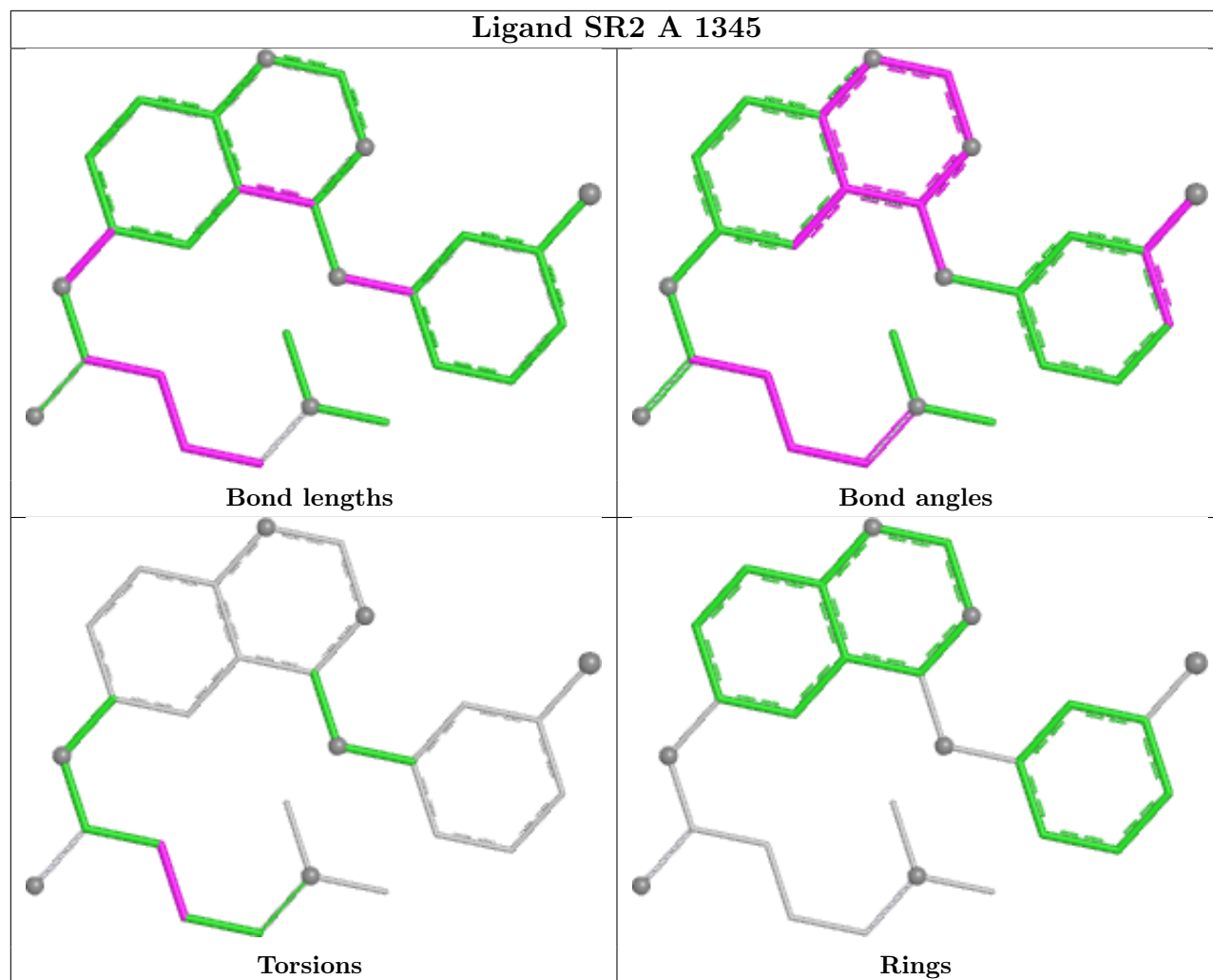
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1345	SR2	2	0
2	A	1345	SR2	5	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	251/286 (87%)	0.77	40 (15%) 5 4	12, 29, 65, 76	4 (1%)
1	B	244/286 (85%)	0.75	30 (12%) 8 7	13, 30, 59, 73	3 (1%)
All	All	495/572 (86%)	0.76	70 (14%) 6 5	12, 30, 64, 76	7 (1%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	404	ASP	7.7
1	B	403	ALA	6.1
1	A	409	ARG	5.8
1	A	309	GLN	5.5
1	B	333	PRO	5.5
1	A	330	SER	5.2
1	B	315	LYS	5.1
1	A	410	LEU	4.8
1	A	278	PHE	4.8
1	A	310	GLU	4.7
1	B	316	LYS	4.5
1	A	407	LEU	4.2
1	A	311	ALA	4.1
1	A	331	GLU	4.1
1	B	324	GLN	4.0
1	A	276	GLY	4.0
1	A	297	LEU	3.9
1	B	313	VAL	3.8
1	A	313	VAL	3.8
1	A	335	TYR	3.6
1	A	273	LEU	3.5
1	B	260	TRP	3.5
1	B	406	GLY	3.4
1	B	317	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	314	MET	3.2
1	B	312	GLN	3.1
1	A	408	ALA	3.0
1	A	280	GLU	3.0
1	B	504	GLU	2.9
1	A	268	ARG	2.8
1	B	407	LEU	2.8
1	B	405	PHE	2.8
1	A	332	GLU	2.8
1	A	333	PRO	2.8
1	B	288	GLY	2.8
1	B	296	THR	2.7
1	A	508	THR	2.7
1	A	334	ILE	2.7
1	B	276	GLY	2.7
1	A	257	LYS	2.6
1	A	425	PRO	2.6
1	A	270	GLU	2.6
1	B	331	GLU	2.5
1	A	271	VAL	2.5
1	A	436	TYR	2.5
1	A	267	LEU	2.5
1	A	293	ALA	2.5
1	A	274	GLY	2.5
1	A	281	VAL	2.5
1	A	290	THR	2.4
1	A	275	GLN	2.4
1	A	312	GLN	2.4
1	A	294	ILE	2.4
1	B	297	LEU	2.4
1	A	269	LEU	2.3
1	B	287	ASN	2.3
1	A	282	TRP	2.3
1	A	328	VAL	2.3
1	B	273	LEU	2.2
1	B	472	LEU	2.2
1	B	318	ARG	2.2
1	B	436	TYR	2.2
1	B	292	VAL	2.2
1	B	473	ASP	2.1
1	A	289	THR	2.1
1	A	292	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	332	GLU	2.1
1	B	526	GLN	2.1
1	A	472	LEU	2.0
1	B	357	TYR	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](#)

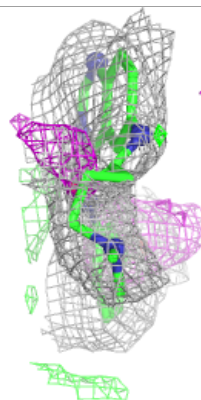
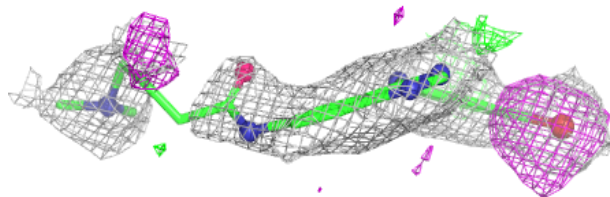
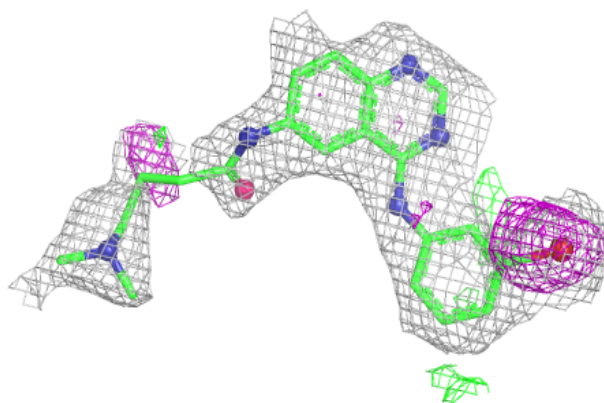
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	SR2	B	1345	27/27	0.80	0.20	45,60,73,76	0
2	SR2	A	1345	27/27	0.81	0.20	45,66,76,79	0

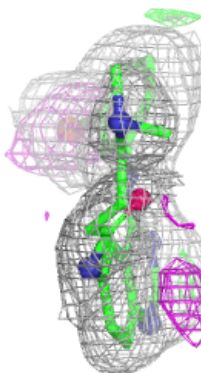
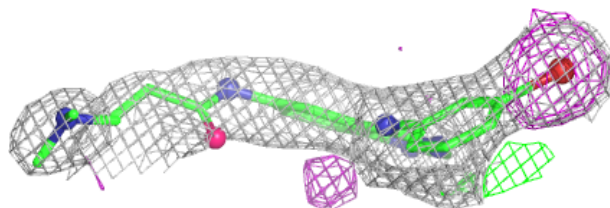
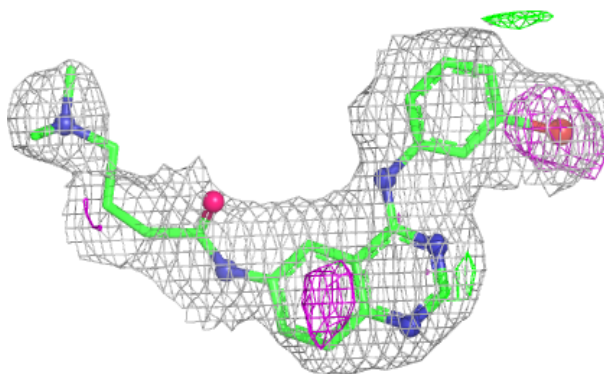
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around SR2 B 1345:

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

**Electron density around SR2 A 1345:**

$2mF_o-DF_c$ (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.