



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 11:21 PM EDT

PDB ID : 4L02 / pdb_00004l02
Title : Crystal Structure of SphK1 with inhibitor
Authors : Min, X.; Walker, N.; Wang, Z.
Deposited on : 2013-05-30
Resolution : 2.75 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

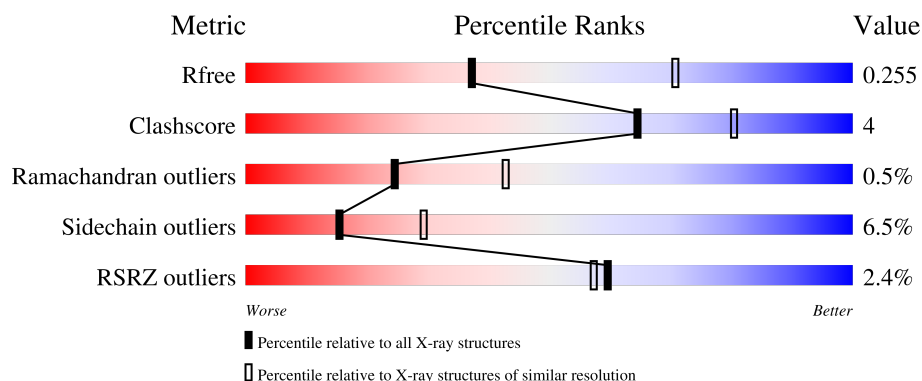
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	384	 2% 80% 12% • 6%
1	B	384	 3% 72% 15% • 11%
1	C	384	 2% 78% 11% • 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8318 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 Sphingosine kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	360	Total	C	N	O	S	0	2	0
			2806	1795	498	491	22			
1	B	340	Total	C	N	O	S	0	1	0
			2668	1711	470	467	20			
1	C	345	Total	C	N	O	S	0	0	0
			2693	1724	477	471	21			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q9NYA1
A	-18	SER	-	expression tag	UNP Q9NYA1
A	-17	TYR	-	expression tag	UNP Q9NYA1
A	-16	TYR	-	expression tag	UNP Q9NYA1
A	-15	HIS	-	expression tag	UNP Q9NYA1
A	-14	HIS	-	expression tag	UNP Q9NYA1
A	-13	HIS	-	expression tag	UNP Q9NYA1
A	-12	HIS	-	expression tag	UNP Q9NYA1
A	-11	HIS	-	expression tag	UNP Q9NYA1
A	-10	HIS	-	expression tag	UNP Q9NYA1
A	-9	ASP	-	expression tag	UNP Q9NYA1
A	-8	TYR	-	expression tag	UNP Q9NYA1
A	-7	ASP	-	expression tag	UNP Q9NYA1
A	-6	ILE	-	expression tag	UNP Q9NYA1
A	-5	PRO	-	expression tag	UNP Q9NYA1
A	-4	THR	-	expression tag	UNP Q9NYA1
A	-3	THR	-	expression tag	UNP Q9NYA1
A	-2	GLU	-	expression tag	UNP Q9NYA1
A	-1	ASN	-	expression tag	UNP Q9NYA1
A	0	LEU	-	expression tag	UNP Q9NYA1
A	1	TYR	-	expression tag	UNP Q9NYA1
A	2	PHE	-	expression tag	UNP Q9NYA1
A	3	GLN	-	expression tag	UNP Q9NYA1

Continued on next page...

Continued from previous page...

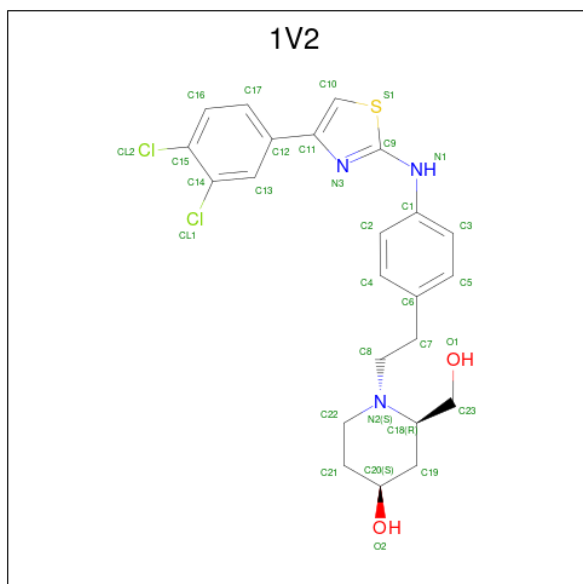
Chain	Residue	Modelled	Actual	Comment	Reference
A	4	GLY	-	expression tag	UNP Q9NYA1
A	5	ALA	-	expression tag	UNP Q9NYA1
A	6	MET	-	expression tag	UNP Q9NYA1
A	7	GLY	-	expression tag	UNP Q9NYA1
A	8	SER	-	expression tag	UNP Q9NYA1
B	-19	MET	-	expression tag	UNP Q9NYA1
B	-18	SER	-	expression tag	UNP Q9NYA1
B	-17	TYR	-	expression tag	UNP Q9NYA1
B	-16	TYR	-	expression tag	UNP Q9NYA1
B	-15	HIS	-	expression tag	UNP Q9NYA1
B	-14	HIS	-	expression tag	UNP Q9NYA1
B	-13	HIS	-	expression tag	UNP Q9NYA1
B	-12	HIS	-	expression tag	UNP Q9NYA1
B	-11	HIS	-	expression tag	UNP Q9NYA1
B	-10	HIS	-	expression tag	UNP Q9NYA1
B	-9	ASP	-	expression tag	UNP Q9NYA1
B	-8	TYR	-	expression tag	UNP Q9NYA1
B	-7	ASP	-	expression tag	UNP Q9NYA1
B	-6	ILE	-	expression tag	UNP Q9NYA1
B	-5	PRO	-	expression tag	UNP Q9NYA1
B	-4	THR	-	expression tag	UNP Q9NYA1
B	-3	THR	-	expression tag	UNP Q9NYA1
B	-2	GLU	-	expression tag	UNP Q9NYA1
B	-1	ASN	-	expression tag	UNP Q9NYA1
B	0	LEU	-	expression tag	UNP Q9NYA1
B	1	TYR	-	expression tag	UNP Q9NYA1
B	2	PHE	-	expression tag	UNP Q9NYA1
B	3	GLN	-	expression tag	UNP Q9NYA1
B	4	GLY	-	expression tag	UNP Q9NYA1
B	5	ALA	-	expression tag	UNP Q9NYA1
B	6	MET	-	expression tag	UNP Q9NYA1
B	7	GLY	-	expression tag	UNP Q9NYA1
B	8	SER	-	expression tag	UNP Q9NYA1
C	-19	MET	-	expression tag	UNP Q9NYA1
C	-18	SER	-	expression tag	UNP Q9NYA1
C	-17	TYR	-	expression tag	UNP Q9NYA1
C	-16	TYR	-	expression tag	UNP Q9NYA1
C	-15	HIS	-	expression tag	UNP Q9NYA1
C	-14	HIS	-	expression tag	UNP Q9NYA1
C	-13	HIS	-	expression tag	UNP Q9NYA1
C	-12	HIS	-	expression tag	UNP Q9NYA1
C	-11	HIS	-	expression tag	UNP Q9NYA1

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	HIS	-	expression tag	UNP Q9NYA1
C	-9	ASP	-	expression tag	UNP Q9NYA1
C	-8	TYR	-	expression tag	UNP Q9NYA1
C	-7	ASP	-	expression tag	UNP Q9NYA1
C	-6	ILE	-	expression tag	UNP Q9NYA1
C	-5	PRO	-	expression tag	UNP Q9NYA1
C	-4	THR	-	expression tag	UNP Q9NYA1
C	-3	THR	-	expression tag	UNP Q9NYA1
C	-2	GLU	-	expression tag	UNP Q9NYA1
C	-1	ASN	-	expression tag	UNP Q9NYA1
C	0	LEU	-	expression tag	UNP Q9NYA1
C	1	TYR	-	expression tag	UNP Q9NYA1
C	2	PHE	-	expression tag	UNP Q9NYA1
C	3	GLN	-	expression tag	UNP Q9NYA1
C	4	GLY	-	expression tag	UNP Q9NYA1
C	5	ALA	-	expression tag	UNP Q9NYA1
C	6	MET	-	expression tag	UNP Q9NYA1
C	7	GLY	-	expression tag	UNP Q9NYA1
C	8	SER	-	expression tag	UNP Q9NYA1

- Molecule 2 is (2R,4S)-1-[2-(4-{[4-(3,4-dichlorophenyl)-1,3-thiazol-2-yl]amino}phenyl)ethyl]-2-(hydroxymethyl)piperidin-4-ol (CCD ID: 1V2) (formula: C₂₃H₂₅Cl₂N₃O₂S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	S	0	0
			31	23	2	3	2	1		

Continued on next page...

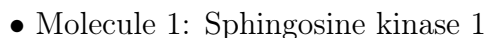
Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	Cl	N	O	S	
			31	23	2	3	2	1	0
2	C	1	Total	C	Cl	N	O	S	
			31	23	2	3	2	1	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	22	Total	O		
			22	22	0	0
3	B	17	Total	O		
			17	17	0	0
3	C	19	Total	O		
			19	19	0	0

- Molecule 1: Sphingosine kinase 1





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	102.58Å 225.68Å 106.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.75 30.00 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-2.75) 99.8 (30.00-2.75)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.76Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.208 , 0.255 0.207 , 0.255	Depositor DCC
R_{free} test set	1645 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.090	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 24.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8318	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: 1V2

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	0.38	0/2879	0.76	0/3906
1	B	0.38	0/2731	0.77	0/3702
1	C	0.38	0/2755	0.78	0/3735
All	All	0.38	0/8365	0.77	0/11343

There are no bond length outliers.

There are no bond angle outliers.

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	2806	0	2852	19	0
1	B	2668	0	2699	28	0
1	C	2693	0	2724	17	0
2	A	31	0	25	1	0
2	B	31	0	25	0	0
2	C	31	0	25	0	0
3	A	22	0	0	0	0
3	B	17	0	0	1	0
3	C	19	0	0	0	0
All	All	8318	0	8350	64	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.

All (64) 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:57:ARG:HG3	1:B:57:ARG:HH11	1.34	0.91
1:B:177:VAL:O	1:B:181:SER:HB2	1.92	0.69
1:B:203:LEU:HD12	1:B:296:ARG:HG2	1.77	0.66
1:B:18:LEU:HD11	1:B:52:MET:HE3	1.77	0.65
1:C:335:LYS:HB3	1:C:346:VAL:HG12	1.81	0.61
1:B:57:ARG:HG3	1:B:57:ARG:NH1	2.08	0.59
1:A:286:HIS:HD2	1:A:321:TYR:OH	1.87	0.58
1:B:169:LEU:HD12	1:B:340:VAL:HG22	1.88	0.56
1:B:55:GLU:O	1:B:56:ARG:HB3	2.06	0.55
1:C:169:LEU:HD12	1:C:340:VAL:HG22	1.89	0.55
1:C:18:LEU:HD11	1:C:52:MET:HE3	1.90	0.54
1:C:76:VAL:HB	1:C:106:CYS:HB3	1.90	0.54
1:A:153:LEU:HB3	1:A:165:SER:HB3	1.90	0.54
1:C:88:VAL:HG21	1:C:166:VAL:HG11	1.89	0.53
1:B:88:VAL:HG21	1:B:166:VAL:HG11	1.91	0.52
1:C:80:GLY:HA2	1:C:114:ASN:ND2	2.25	0.52
1:B:363:SER:HB2	3:B:517:HOH:O	2.10	0.51
1:C:77:VAL:CG1	1:C:83:LEU:HB3	2.40	0.51
1:B:286:HIS:HD2	1:B:321:TYR:OH	1.94	0.50
1:A:69:LEU:HB3	1:A:103:LYS:HE3	1.93	0.50
1:A:294:VAL:HG13	1:A:318:TYR:HB2	1.93	0.50
1:A:258:VAL:HG22	1:A:290:VAL:HB	1.94	0.50
1:A:207:ARG:HB3	1:A:332:LYS:HD2	1.92	0.50
1:C:116:LEU:HD11	1:C:361:MET:HE1	1.93	0.49
1:A:170:ALA:HB3	1:A:339:ALA:HB3	1.94	0.49
1:C:22:ASN:C	1:C:22:ASN:HD22	2.21	0.49
1:B:205:THR:HG22	1:B:258:VAL:HG12	1.95	0.48
1:B:149:PRO:HB2	1:B:357:ASN:HB3	1.95	0.48
1:A:199:ARG:HD2	1:A:199:ARG:HA	1.72	0.47
1:B:277:MET:HE1	1:B:312:MET:HE2	1.97	0.47
1:A:328:ARG:HE	1:A:351:GLN:NE2	2.12	0.47
1:A:149:PRO:HB2	1:A:357:ASN:HB3	1.97	0.46
1:A:200:LEU:HD21	1:A:300:LEU:HB2	1.97	0.46
1:C:207:ARG:HB3	1:C:332:LYS:HE3	1.96	0.46
1:B:22:ASN:HD22	1:B:23:PRO:HD2	1.81	0.45
1:A:266:SER:HA	1:A:275:ALA:HB3	1.98	0.45
1:B:157:THR:OG1	1:B:161:LEU:HB3	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:LEU:HD12	1:A:340:VAL:HG22	1.98	0.45
1:B:179:LEU:HG	1:B:344:LEU:HG	1.98	0.44
1:C:115:ALA:HB3	1:C:167:LEU:HD21	1.99	0.44
1:A:112:SER:HB2	1:A:191:ARG:HH21	1.82	0.44
1:A:231:GLN:HG2	1:A:232:GLY:N	2.32	0.44
1:B:210:LEU:HD23	1:B:329:LEU:HD13	2.00	0.43
1:A:356:PRO:O	1:A:357:ASN:C	2.62	0.43
1:C:153:LEU:HB3	1:C:165:SER:HB3	1.99	0.43
1:B:29:LYS:O	1:B:32:GLN:HG2	2.19	0.43
1:B:57:ARG:NH1	1:B:57:ARG:CG	2.79	0.43
1:B:329:LEU:HD23	1:B:338:PHE:HE2	1.84	0.43
1:C:158:ALA:HB2	1:C:349:ALA:HB3	2.00	0.43
2:A:401:1V2:N3	2:A:401:1V2:H8	2.34	0.43
1:C:83:LEU:HD23	1:C:83:LEU:HA	1.91	0.42
1:A:19:VAL:HG22	1:A:76:VAL:HG22	2.01	0.42
1:B:274:ALA:HB2	1:B:306:MET:HE3	2.02	0.42
1:B:193:THR:HG23	1:B:303:PHE:HE2	1.84	0.42
1:C:55:GLU:HB2	1:C:56:ARG:HD3	2.02	0.41
1:B:55:GLU:O	1:B:56:ARG:CB	2.69	0.41
1:B:201:ALA:HB1	1:C:300:LEU:HD21	2.00	0.41
1:B:109:PRO:HG3	1:B:134:LEU:HG	2.02	0.41
1:B:193:THR:HG22	1:B:197:PHE:CZ	2.55	0.41
1:A:99:THR:HA	1:A:102:GLN:HE21	1.86	0.41
1:B:103:LYS:HD3	1:B:104:PRO:HD2	2.04	0.40
1:A:196:THR:O	1:A:200:LEU:HB2	2.21	0.40
1:B:80:GLY:HA2	1:B:114:ASN:OD1	2.21	0.40
1:C:71:ARG:HD3	1:C:72:TRP:CE2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	360/384 (94%)	349 (97%)	10 (3%)	1 (0%)	36	56
1	B	333/384 (87%)	322 (97%)	8 (2%)	3 (1%)	14	28
1	C	341/384 (89%)	331 (97%)	9 (3%)	1 (0%)	36	56
All	All	1034/1152 (90%)	1002 (97%)	27 (3%)	5 (0%)	24	43

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	79	SER
1	A	79	SER
1	B	79	SER
1	B	56	ARG
1	B	189	GLU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	301/322 (94%)	280 (93%)	21 (7%)	14	26
1	B	286/322 (89%)	267 (93%)	19 (7%)	15	29
1	C	287/322 (89%)	269 (94%)	18 (6%)	16	30
All	All	874/966 (90%)	816 (93%)	58 (7%)	15	29

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	LEU
1	A	35[A]	ARG
1	A	35[B]	ARG
1	A	57	ARG
1	A	76	VAL
1	A	83	LEU
1	A	98	GLU
1	A	105	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	132	GLU
1	A	152	LEU
1	A	153	LEU
1	A	169	LEU
1	A	179	LEU
1	A	210	LEU
1	A	258	VAL
1	A	294	VAL
1	A	296	ARG
1	A	322	VAL
1	A	337	VAL
1	A	358	TYR
1	A	362	VAL
1	B	32	GLN
1	B	35	ARG
1	B	50	THR
1	B	57	ARG
1	B	105	LEU
1	B	132	GLU
1	B	147	LEU
1	B	152	LEU
1	B	153	LEU
1	B	169	LEU
1	B	179	LEU
1	B	185	ARG
1	B	199	ARG
1	B	242	GLU
1	B	245	VAL
1	B	294	VAL
1	B	337	VAL
1	B	346	VAL
1	B	354	VAL
1	C	17	VAL
1	C	29	LYS
1	C	35	ARG
1	C	50	THR
1	C	71	ARG
1	C	76	VAL
1	C	83	LEU
1	C	105	LEU
1	C	147	LEU
1	C	152	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	153	LEU
1	C	162	ARG
1	C	210	LEU
1	C	215	VAL
1	C	245	VAL
1	C	301	ARG
1	C	345	MET
1	C	348	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	85	HIS
1	A	102	GLN
1	A	131	ASN
1	A	286	HIS
1	A	351	GLN
1	B	22	ASN
1	B	32	GLN
1	B	102	GLN
1	B	286	HIS
1	B	351	GLN
1	C	22	ASN
1	C	58	ASN
1	C	102	GLN
1	C	114	ASN
1	C	286	HIS
1	C	351	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	1V2	C	401	-	34,34,34	1.59	3 (8%)	46,47,47	2.18	7 (15%)
2	1V2	B	401	-	34,34,34	1.58	3 (8%)	46,47,47	2.20	5 (10%)
2	1V2	A	401	-	34,34,34	1.66	3 (8%)	46,47,47	2.28	5 (10%)

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	1V2	C	401	-	-	2/15/28/28	0/4/4/4
2	1V2	B	401	-	-	2/15/28/28	0/4/4/4
2	1V2	A	401	-	-	1/15/28/28	0/4/4/4

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	1V2	C9-S1	-5.49	1.66	1.74
2	B	401	1V2	C12-C11	-5.43	1.39	1.47
2	A	401	1V2	C12-C11	-5.42	1.39	1.47
2	C	401	1V2	C12-C11	-5.30	1.39	1.47
2	B	401	1V2	C9-S1	-4.85	1.67	1.74
2	C	401	1V2	C9-S1	-4.84	1.67	1.74
2	A	401	1V2	C1-N1	-3.56	1.34	1.41
2	B	401	1V2	C1-N1	-3.46	1.34	1.41

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	1V2	C1-N1	-3.41	1.34	1.41

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	1V2	C10-S1-C9	9.21	95.56	88.40
2	B	401	1V2	C10-S1-C9	8.91	95.33	88.40
2	C	401	1V2	C10-S1-C9	8.42	94.94	88.40
2	A	401	1V2	S1-C9-N3	-7.65	107.66	115.69
2	B	401	1V2	S1-C9-N3	-7.53	107.78	115.69
2	C	401	1V2	S1-C9-N3	-7.18	108.15	115.69
2	A	401	1V2	C9-N3-C11	6.56	115.26	109.96
2	B	401	1V2	C9-N3-C11	6.51	115.21	109.96
2	C	401	1V2	C9-N3-C11	6.15	114.92	109.96
2	A	401	1V2	C11-C10-S1	-3.24	107.23	111.16
2	B	401	1V2	C11-C10-S1	-3.20	107.29	111.16
2	C	401	1V2	C12-C11-C10	-3.20	122.75	126.54
2	B	401	1V2	C12-C11-C10	-2.96	123.03	126.54
2	A	401	1V2	C12-C11-C10	-2.90	123.10	126.54
2	C	401	1V2	C11-C10-S1	-2.84	107.72	111.16
2	C	401	1V2	C19-C18-N2	2.46	112.95	110.51
2	C	401	1V2	N1-C9-N3	2.01	128.50	122.51

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	1V2	C7-C8-N2-C22
2	B	401	1V2	N3-C9-N1-C1
2	C	401	1V2	C4-C6-C7-C8
2	C	401	1V2	C5-C6-C7-C8
2	A	401	1V2	C4-C6-C7-C8

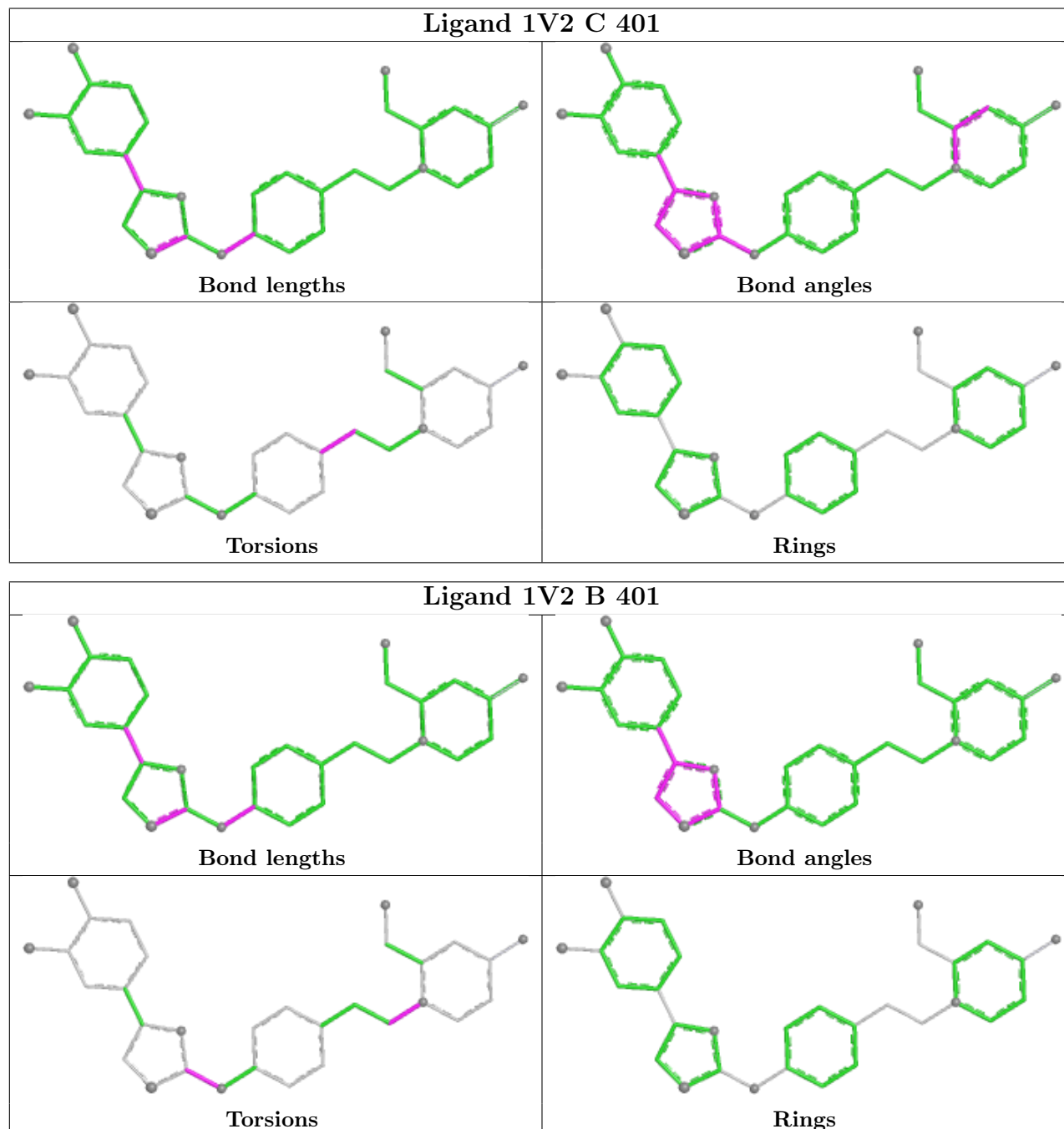
There are no ring outliers.

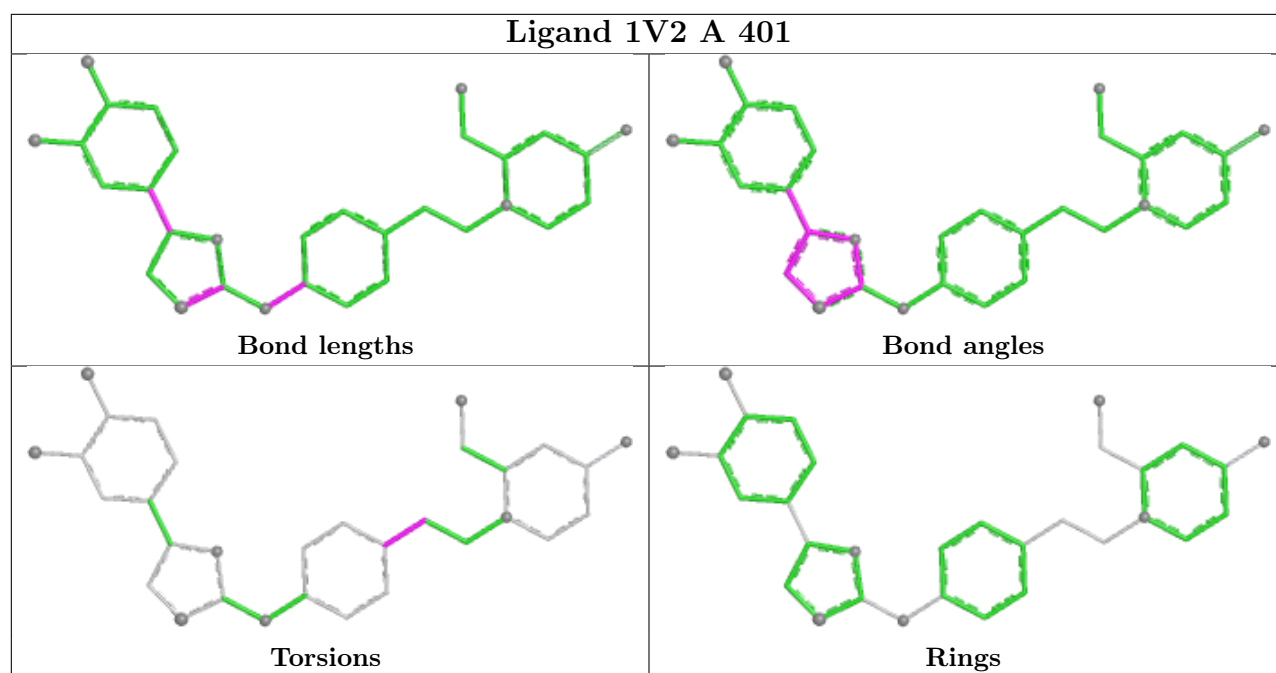
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	1V2	1	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	360/384 (93%)	-0.09	8 (2%) 62 60	18, 34, 65, 78	2 (0%)
1	B	340/384 (88%)	0.18	11 (3%) 50 48	21, 41, 71, 84	1 (0%)
1	C	345/384 (89%)	0.01	6 (1%) 69 67	27, 40, 63, 86	0
All	All	1045/1152 (90%)	0.03	25 (2%) 59 57	18, 38, 67, 86	3 (0%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	228	VAL	4.7
1	B	332	LYS	3.3
1	C	334	GLY	3.1
1	B	188	GLY	3.0
1	B	197	PHE	3.0
1	B	184	TYR	3.0
1	B	185	ARG	2.8
1	B	335	LYS	2.7
1	A	6[A]	MET	2.7
1	B	363	SER	2.5
1	C	217	ARG	2.5
1	A	227	VAL	2.4
1	B	344	LEU	2.4
1	C	216	GLY	2.4
1	B	198	LEU	2.4
1	C	348	GLU	2.3
1	A	52	MET	2.2
1	A	55	GLU	2.2
1	A	333	ASP	2.2
1	B	52	MET	2.2
1	A	222	THR	2.2
1	A	229	VAL	2.2
1	C	336	GLY	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	57	ARG	2.0
1	B	203	LEU	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](#)

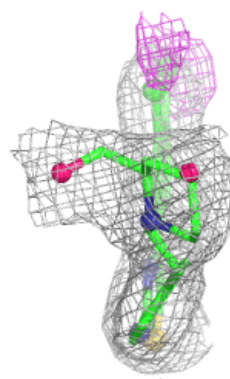
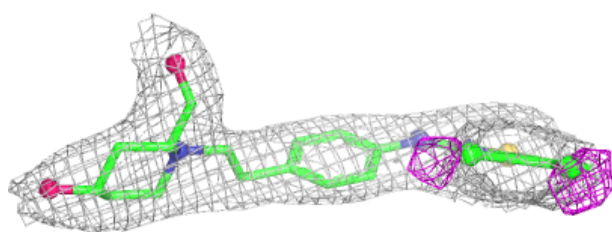
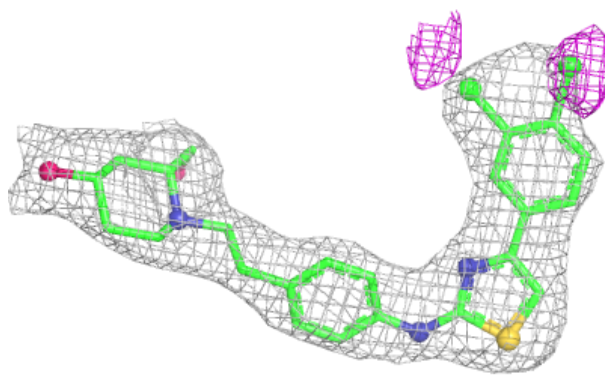
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	1V2	B	401	31/31	0.89	0.11	36,48,67,75	0
2	1V2	C	401	31/31	0.89	0.12	30,39,65,72	0
2	1V2	A	401	31/31	0.90	0.10	26,29,48,54	0

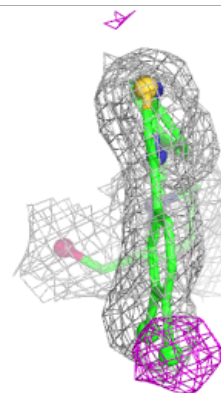
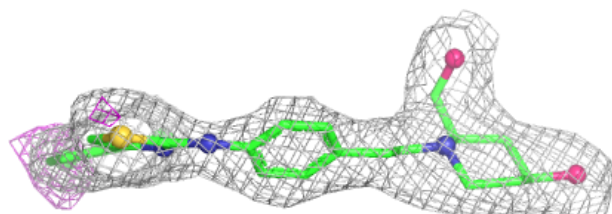
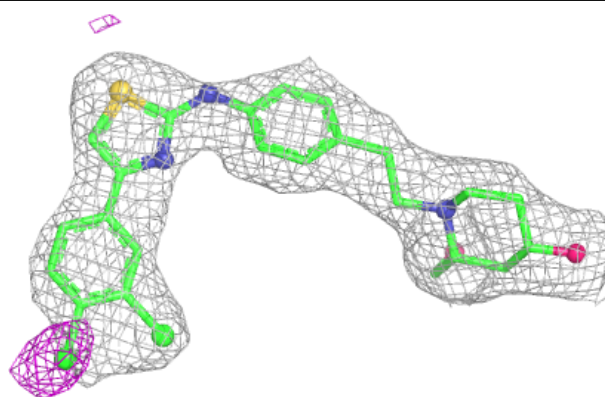
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 1V2 B 401:

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

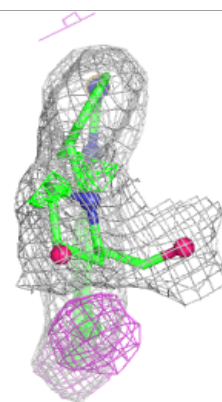
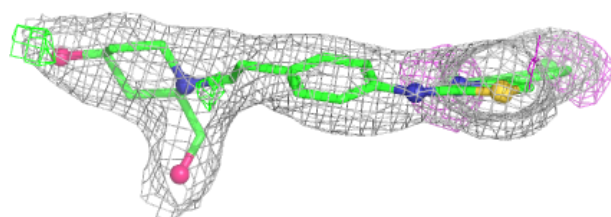
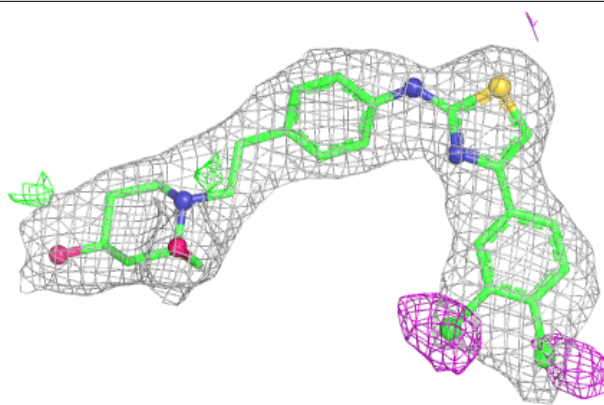
**Electron density around 1V2 C 401:**

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



Electron density around 1V2 A 401:

$2mF_o - DF_c$ (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.