



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 02:41 am BST

PDB ID : 3EMG
Title : Discovery and SAR of novel 4-thiazolyl-2-phenylaminopyrimidines as potent inhibitors of spleen tyrosine kinase (SYK)
Authors : Ter Haar, E.
Deposited on : 2008-09-24
Resolution : 2.60 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

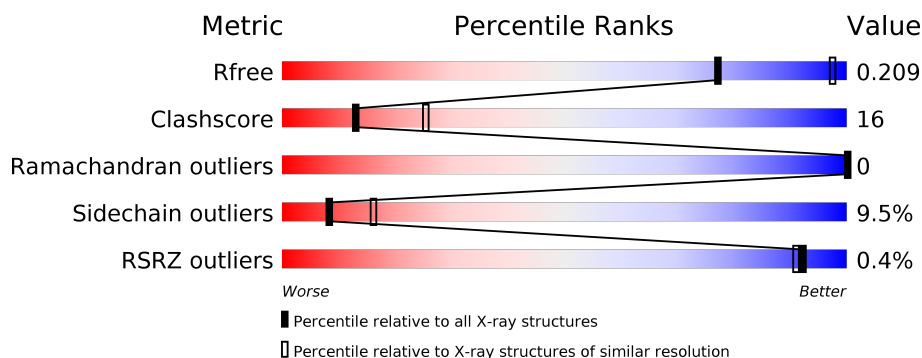
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	291	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2192 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.

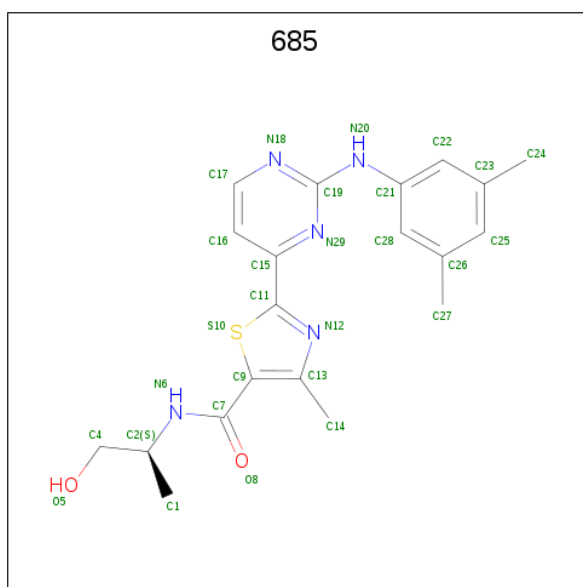
- Molecule 1 is a protein called Tyrosine-protein kinase SYK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	255	2081	1339	349	374	19	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	345	GLY	-	EXPRESSION TAG	UNP P43405
A	346	SER	-	EXPRESSION TAG	UNP P43405
A	347	LEU	-	EXPRESSION TAG	UNP P43405
A	348	HIS	-	EXPRESSION TAG	UNP P43405

- Molecule 2 is 2-{2-[(3,5-dimethylphenyl)amino]pyrimidin-4-yl}-N-[(1S)-2-hydroxy-1-methyl-4-methyl-1,3-thiazole-5-carboxamide (three-letter code: 685) (formula: C₂₀H₂₃N₅O₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	28	20	5	2	1	0	0

- Molecule 3 is water.

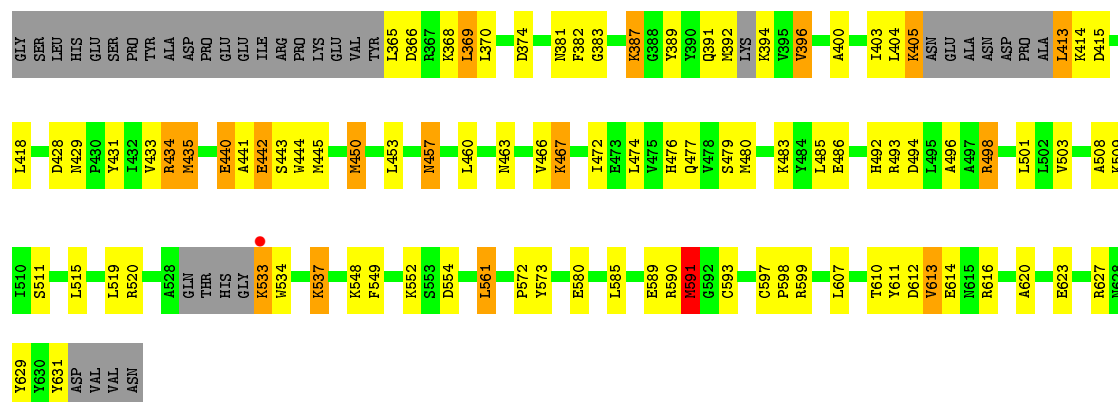
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	83	Total	O	0	0
			83	83		

3 Residue-property plots [i](#)

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: Tyrosine-protein kinase SYK

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.87Å 84.45Å 41.44Å 90.00° 99.63° 90.00°	Depositor
Resolution (Å)	39.30 – 2.60 42.22 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (39.30-2.60) 99.8 (42.22-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.70 (at 2.61Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.200 , 0.222 0.191 , 0.209	Depositor DCC
R_{free} test set	1036 reflections (12.36%)	wwPDB-VP
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	1.737	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.045 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2192	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.34% 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: 685

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	1.44	6/2126 (0.3%)	1.31	23/2859 (0.8%)

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	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	593	CYS	CB-SG	-9.46	1.66	1.82
1	A	548	LYS	CD-CE	6.25	1.66	1.51
1	A	580	GLU	CD-OE2	-5.99	1.19	1.25
1	A	589	GLU	CD-OE1	5.78	1.32	1.25
1	A	433	VAL	CB-CG1	5.45	1.64	1.52
1	A	611	TYR	CD2-CE2	5.01	1.46	1.39

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	441	ALA	N-CA-C	12.77	145.49	111.00
1	A	441	ALA	N-CA-CB	-10.36	95.59	110.10
1	A	498	ARG	NE-CZ-NH2	8.04	124.32	120.30
1	A	498	ARG	NE-CZ-NH1	-7.74	116.43	120.30
1	A	434	ARG	NE-CZ-NH1	-7.33	116.64	120.30
1	A	440	GLU	N-CA-C	6.84	129.46	111.00
1	A	585	LEU	CB-CG-CD1	-6.74	99.54	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	480	MET	CG-SD-CE	6.56	110.70	100.20
1	A	590	ARG	NE-CZ-NH1	-6.45	117.07	120.30
1	A	366	ASP	CB-CG-OD2	6.44	124.10	118.30
1	A	520	ARG	NE-CZ-NH1	-6.30	117.15	120.30
1	A	554	ASP	CB-CG-OD2	-6.07	112.84	118.30
1	A	485	LEU	CB-CG-CD1	-6.05	100.72	111.00
1	A	369	LEU	CA-CB-CG	5.94	128.96	115.30
1	A	435	MET	CG-SD-CE	5.91	109.66	100.20
1	A	442	GLU	N-CA-C	-5.59	95.90	111.00
1	A	612	ASP	CB-CG-OD1	5.57	123.31	118.30
1	A	537	LYS	CD-CE-NZ	-5.50	99.05	111.70
1	A	613	VAL	CB-CA-C	-5.29	101.34	111.40
1	A	486	GLU	OE1-CD-OE2	-5.24	117.02	123.30
1	A	366	ASP	CB-CG-OD1	-5.24	113.59	118.30
1	A	585	LEU	CB-CG-CD2	5.23	119.88	111.00
1	A	591	MET	CG-SD-CE	-5.17	91.92	100.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	394	LYS	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	2081	0	2077	68	1
2	A	28	0	23	1	0
3	A	83	0	0	11	1
All	All	2192	0	2100	69	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 16.

All (69) 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:A:533:LYS:CB	1:A:534:TRP:HA	1.54	1.24
1:A:533:LYS:HB2	1:A:534:TRP:HA	1.22	1.14
1:A:392:MET:HG3	3:A:148:HOH:O	1.49	1.12
1:A:413:LEU:HD12	1:A:415:ASP:HB2	1.18	1.11
1:A:533:LYS:HB3	1:A:534:TRP:HA	1.38	0.99
1:A:413:LEU:CD1	1:A:415:ASP:HB2	1.95	0.95
1:A:533:LYS:CB	1:A:534:TRP:CA	2.47	0.92
1:A:368:LYS:O	3:A:129:HOH:O	1.90	0.89
1:A:404:LEU:HD11	3:A:110:HOH:O	1.75	0.86
1:A:392:MET:CG	3:A:148:HOH:O	2.11	0.85
1:A:533:LYS:HB2	1:A:534:TRP:CD1	2.14	0.83
1:A:620:ALA:HB3	3:A:21:HOH:O	1.79	0.82
1:A:533:LYS:HB2	1:A:534:TRP:CA	2.06	0.81
1:A:533:LYS:HB2	1:A:534:TRP:HD1	1.44	0.79
1:A:387:LYS:HE2	3:A:3:HOH:O	1.82	0.79
1:A:492:HIS:HD2	1:A:494:ASP:H	1.30	0.76
1:A:418:LEU:HD12	3:A:161:HOH:O	1.85	0.76
1:A:369:LEU:HA	3:A:129:HOH:O	1.86	0.75
1:A:515:LEU:HD11	1:A:533:LYS:HG3	1.70	0.72
1:A:457:ASN:HD22	1:A:457:ASN:H	1.36	0.72
1:A:370:LEU:HD12	1:A:389:TYR:O	1.90	0.71
1:A:392:MET:SD	3:A:148:HOH:O	2.48	0.70
1:A:573:TYR:H	1:A:591:MET:CE	2.06	0.69
2:A:685:685:N29	2:A:685:685:H22	2.07	0.68
1:A:440:GLU:O	1:A:440:GLU:HG2	1.96	0.66
1:A:492:HIS:CD2	1:A:494:ASP:H	2.13	0.64
1:A:428:ASP:OD1	1:A:434:ARG:NH1	2.29	0.64
1:A:383:GLY:HA2	1:A:405:LYS:HB2	1.79	0.63
1:A:573:TYR:CG	1:A:591:MET:HE1	2.34	0.61
1:A:403:ILE:HG12	1:A:445:MET:HG2	1.83	0.61
1:A:573:TYR:N	1:A:591:MET:HE1	2.19	0.57
1:A:493:ARG:HD2	1:A:549:PHE:CD2	2.39	0.57
1:A:533:LYS:HE3	1:A:533:LYS:HA	1.86	0.57
1:A:573:TYR:H	1:A:591:MET:HE1	1.70	0.55
1:A:382:PHE:CD1	1:A:383:GLY:N	2.73	0.55
1:A:383:GLY:HA3	1:A:403:ILE:O	2.06	0.55
1:A:573:TYR:H	1:A:591:MET:HE2	1.70	0.55
1:A:476:HIS:HE1	1:A:623:GLU:OE2	1.88	0.55
1:A:610:THR:HB	1:A:616:ARG:HG3	1.90	0.54
1:A:496:ALA:HB2	3:A:156:HOH:O	2.07	0.53
1:A:533:LYS:HB3	1:A:534:TRP:CA	2.25	0.52
1:A:501:LEU:HD12	1:A:511:SER:HB3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:503:VAL:CG2	1:A:509:LYS:HD2	2.40	0.52
1:A:440:GLU:O	1:A:440:GLU:CG	2.55	0.51
1:A:404:LEU:HD12	1:A:404:LEU:C	2.31	0.51
1:A:404:LEU:O	1:A:404:LEU:HD12	2.13	0.49
1:A:561:LEU:O	1:A:561:LEU:HD12	2.13	0.49
1:A:477:GLN:NE2	1:A:508:ALA:H	2.11	0.49
1:A:404:LEU:CD1	3:A:110:HOH:O	2.50	0.48
1:A:515:LEU:CD1	1:A:533:LYS:HG3	2.43	0.48
1:A:405:LYS:HD3	1:A:405:LYS:HA	1.60	0.48
1:A:457:ASN:ND2	1:A:498:ARG:HA	2.29	0.47
1:A:472:ILE:HD13	1:A:629:TYR:HE2	1.80	0.46
1:A:597:CYS:HA	1:A:598:PRO:HD3	1.75	0.45
1:A:463:ASN:O	1:A:466:VAL:HG23	2.17	0.45
1:A:479:SER:O	1:A:483:LYS:N	2.49	0.44
1:A:537:LYS:HG3	1:A:573:TYR:HD2	1.83	0.43
1:A:404:LEU:HG	1:A:444:TRP:HB2	2.00	0.43
1:A:442:GLU:HG2	1:A:443:SER:N	2.34	0.43
1:A:391:GLN:HA	1:A:396:VAL:HG23	2.01	0.43
1:A:400:ALA:HB2	1:A:450:MET:HG2	2.01	0.43
1:A:627:ARG:O	1:A:631:TYR:HD1	2.02	0.43
1:A:429:ASN:ND2	1:A:431:TYR:H	2.18	0.42
1:A:537:LYS:HG2	1:A:572:PRO:HG2	2.02	0.42
1:A:607:LEU:HA	1:A:607:LEU:HD12	1.77	0.41
1:A:460:LEU:HD23	1:A:460:LEU:HA	1.79	0.41
1:A:457:ASN:N	1:A:457:ASN:HD22	2.10	0.41
1:A:467:LYS:HE3	1:A:467:LYS:HB3	1.79	0.41
1:A:453:LEU:HD23	1:A:453:LEU:HA	1.83	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:A:391:GLN:OE1	3:A:127:HOH:O[2_656]	1.86	0.34

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	247/291 (85%)	230 (93%)	17 (7%)	0	100	100

There are no Ramachandran outliers to report.

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	222/253 (88%)	201 (90%)	21 (10%)	8	16

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	365	LEU
1	A	374	ASP
1	A	381	ASN
1	A	387	LYS
1	A	396	VAL
1	A	405	LYS
1	A	413	LEU
1	A	414	LYS
1	A	435	MET
1	A	450	MET
1	A	457	ASN
1	A	467	LYS
1	A	474	LEU
1	A	519	LEU
1	A	533	LYS
1	A	552	LYS
1	A	561	LEU
1	A	591	MET
1	A	599	ARG

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Mol	Chain	Res	Type
1	A	613	VAL
1	A	614	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	429	ASN
1	A	457	ASN
1	A	461	GLN
1	A	476	HIS
1	A	477	GLN
1	A	492	HIS
1	A	505	GLN
1	A	545	ASN
1	A	570	GLN
1	A	628	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	685	A	685	-	26,30,30	3.55	15 (57%)	31,42,42	3.08	12 (38%)

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	685	A	685	-	-	1/11/18/18	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	685	685	C11-S10	-10.89	1.58	1.73
2	A	685	685	C19-N18	-7.26	1.24	1.34
2	A	685	685	C11-N12	-5.35	1.23	1.31
2	A	685	685	O8-C7	-5.11	1.12	1.23
2	A	685	685	C15-N29	-4.04	1.28	1.34
2	A	685	685	C17-N18	-3.58	1.27	1.34
2	A	685	685	C7-N6	-3.15	1.27	1.34
2	A	685	685	C22-C21	-2.72	1.35	1.39
2	A	685	685	C4-C2	-2.58	1.47	1.51
2	A	685	685	C25-C26	-2.36	1.35	1.39
2	A	685	685	C25-C23	-2.35	1.35	1.39
2	A	685	685	C15-C11	-2.32	1.41	1.48
2	A	685	685	C16-C17	-2.29	1.33	1.38
2	A	685	685	C19-N29	-2.19	1.28	1.34
2	A	685	685	C16-C15	-2.03	1.34	1.38

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	685	685	O5-C4-C2	-9.25	92.61	112.37
2	A	685	685	C1-C2-C4	-7.23	101.09	111.54
2	A	685	685	C15-N29-C19	6.48	122.03	116.69
2	A	685	685	N18-C19-N29	-5.11	121.70	126.55
2	A	685	685	O8-C7-N6	-3.96	115.17	122.45
2	A	685	685	C24-C23-C25	-3.80	115.31	120.94
2	A	685	685	C17-N18-C19	3.13	118.23	115.45
2	A	685	685	C17-C16-C15	-2.93	114.45	117.22
2	A	685	685	C16-C15-N29	-2.42	118.83	121.97
2	A	685	685	C16-C15-C11	2.27	123.01	120.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	685	685	C25-C23-C22	2.15	120.65	118.09
2	A	685	685	C24-C23-C22	2.09	124.03	120.94

There are no chirality outliers.

All (1) torsion outliers are listed below:

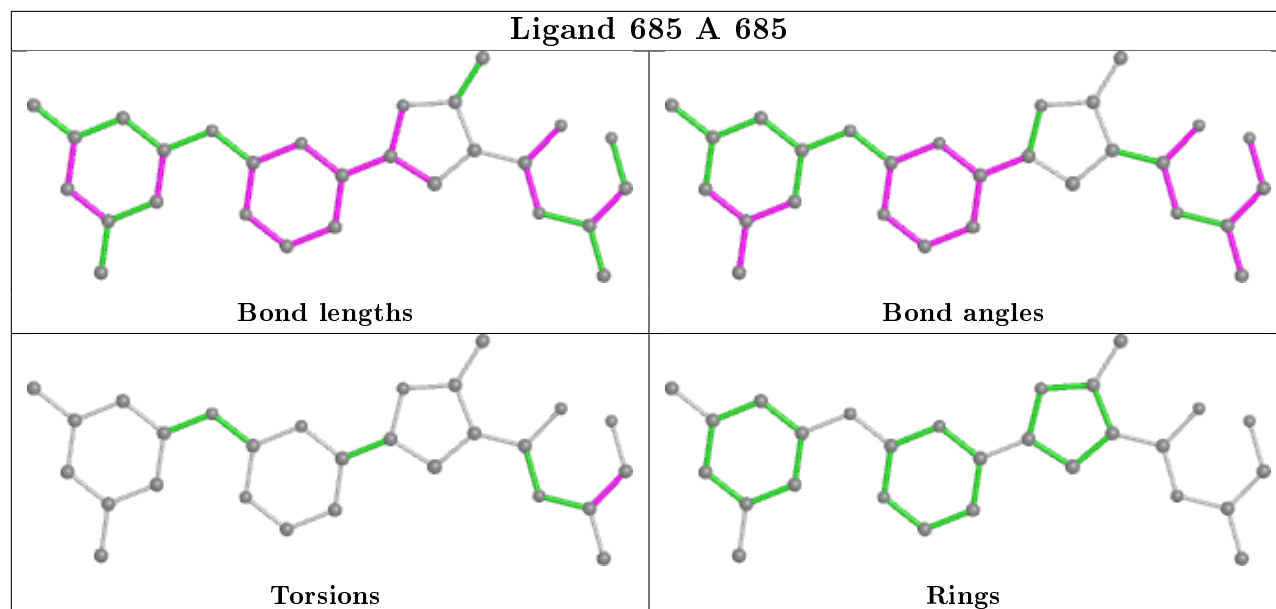
Mol	Chain	Res	Type	Atoms
2	A	685	685	N6-C2-C4-O5

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	685	685	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 [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	255/291 (87%)	-0.29	1 (0%) 92 91	5, 21, 46, 70	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	533	LYS	2.8

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](#)

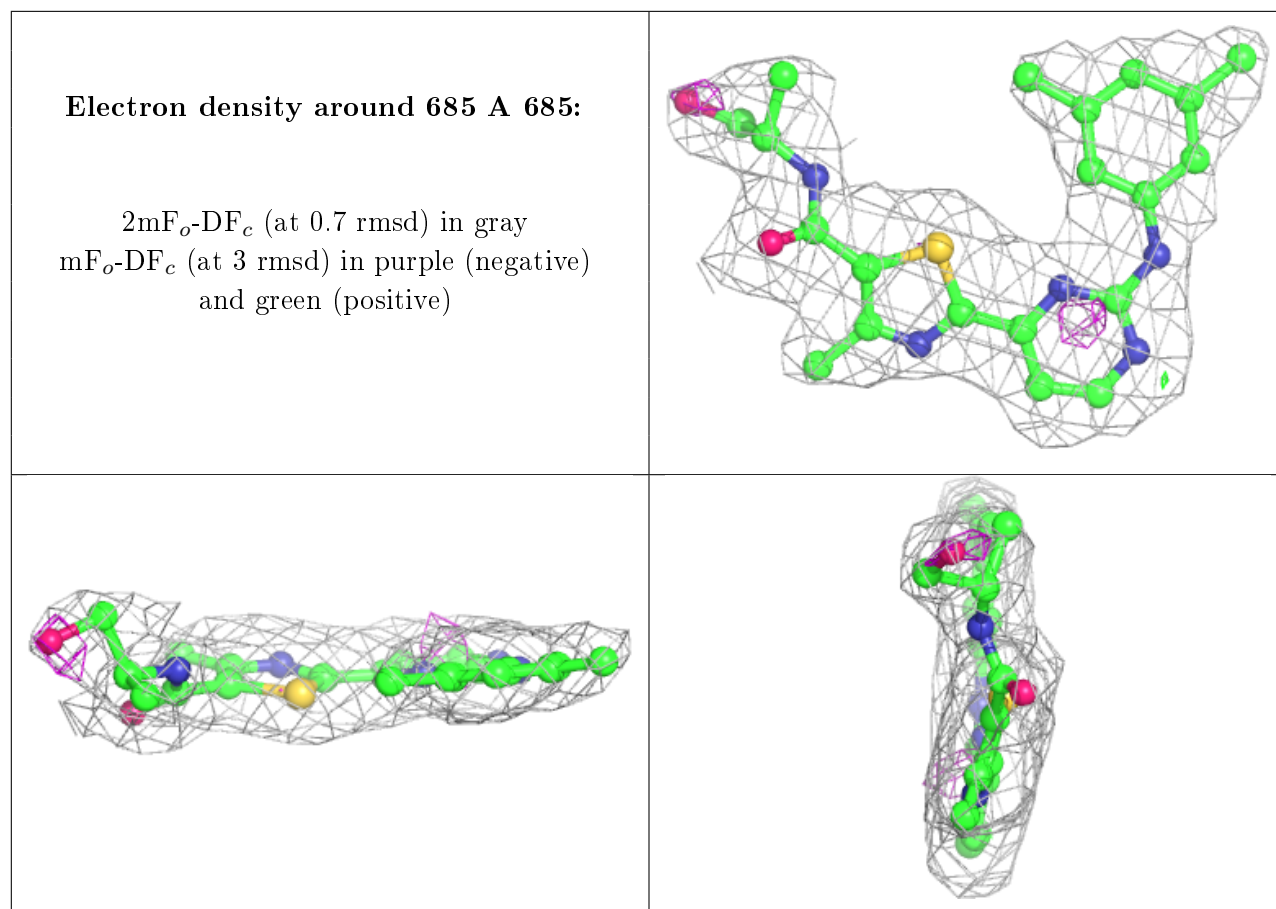
There are no carbohydrates in this entry.

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	685	A	685	28/28	0.94	0.18	21,21,21,21	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.