



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 13, 2020 – 10:53 PM BST

PDB ID : 5ZH3  
Title : CRYSTAL STRUCTURE OF PfkRS WITH INHIBITOR CLADO-6  
Authors : Babbar, P.; Malhotra, N.; Sharma, M.; Harlos, K.; Reddy, D.S.; Manickam, Y.; Sharma, A.  
Deposited on : 2018-03-11  
Resolution : 2.86 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](mailto: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 : 2.14.4.dev1  
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) : 2.14.4.dev1

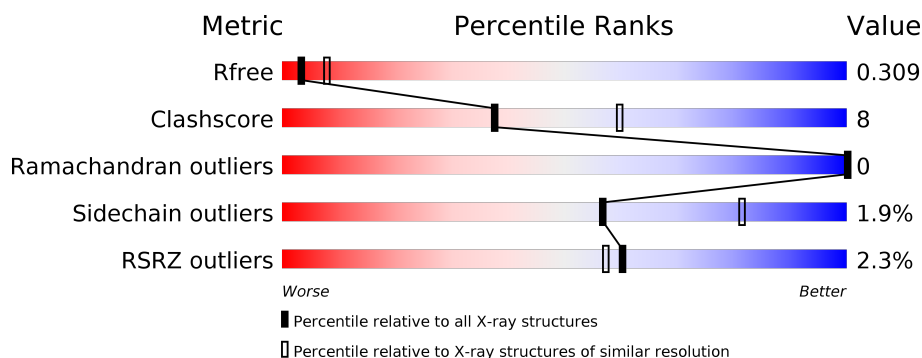
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.86 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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  $\geq 3$ , 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	507	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>17%</div> <div>• 7%</div> </div> </div>
1	B	507	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>17%</div> <div>7%</div> </div> </div>

## 2 Entry composition [i](#)

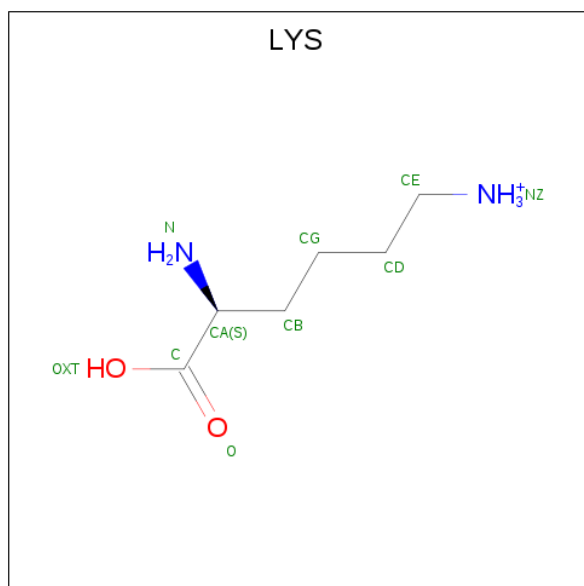
There are 3 unique types of molecules in this entry. The entry contains 7311 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Lysine-tRNA ligase.

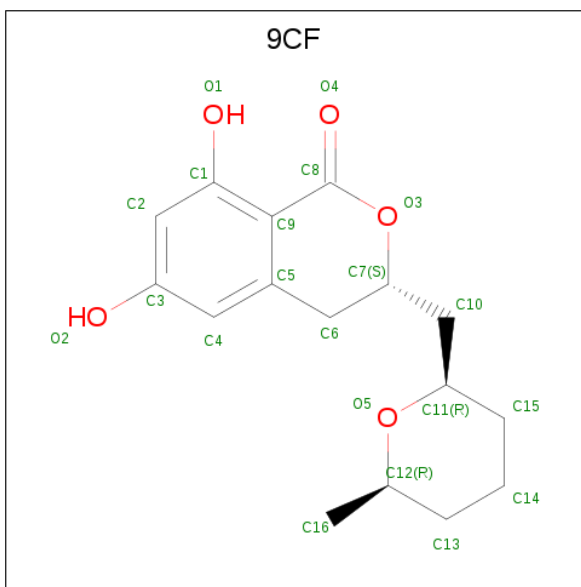
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	469	Total	C	N	O	S	0	0	0
			3592	2318	598	661	15			
1	B	474	Total	C	N	O	S	0	0	0
			3657	2360	613	670	14			

- Molecule 2 is LYSINE (three-letter code: LYS) (formula:  $C_6H_{15}N_2O_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			10	6	2	2		
2	B	1	Total	C	N	O	0	0
			10	6	2	2		

- Molecule 3 is (3S)-6,8-dihydroxy-3-[(2R,6R)-6-methyloxan-2-yl]methyl}-3,4-dihydro-1H-2-benzopyran-1-one (three-letter code: 9CF) (formula:  $C_{16}H_{20}O_5$ ).

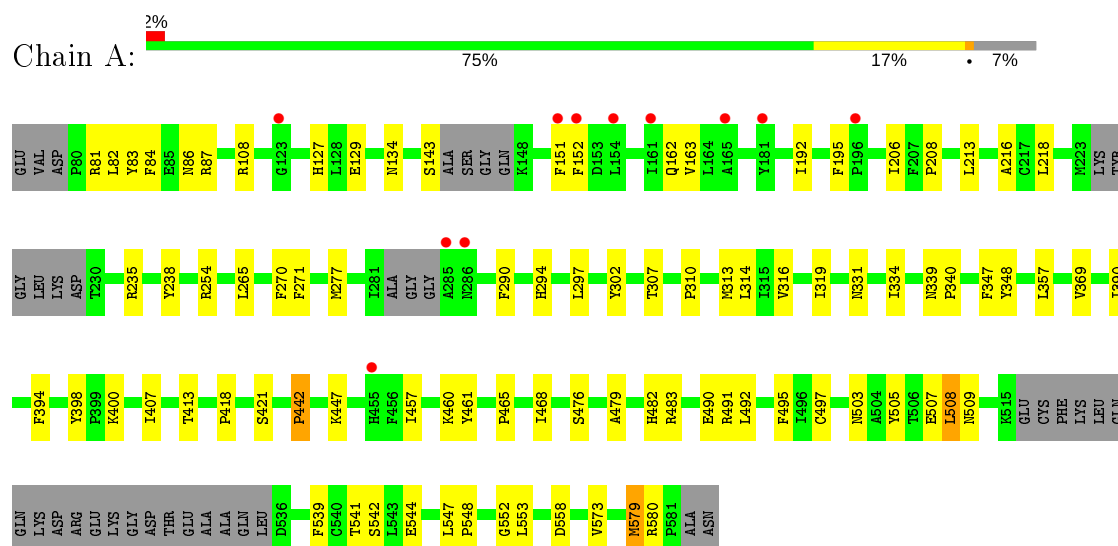


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			21	16	5		
3	B	1	Total	C	O	0	0
			21	16	5		

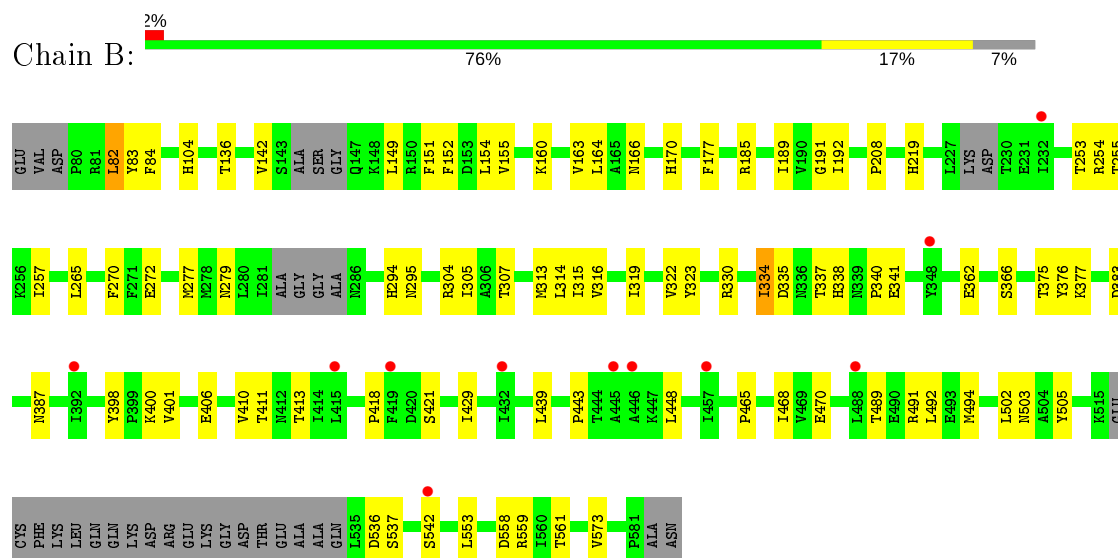
### 3 Residue-property plots

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: Lysine-tRNA ligase



#### • Molecule 1: Lysine-tRNA ligase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.44Å 130.77Å 174.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	53.20 – 2.86 53.20 – 2.86	Depositor EDS
% Data completeness (in resolution range)	99.6 (53.20-2.86) 99.6 (53.20-2.86)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.09 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.251 , 0.311 0.256 , 0.309	Depositor DCC
$R_{free}$ test set	1418 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	77.2	Xtriage
Anisotropy	0.861	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 75.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7311	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	108.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 9CF

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	0.43	1/3683 (0.0%)	0.62	1/5013 (0.0%)
1	B	0.42	0/3747	0.61	1/5091 (0.0%)
All	All	0.42	1/7430 (0.0%)	0.61	2/10104 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	442	PRO	N-CD	5.39	1.55	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	579	MET	O-C-N	-13.01	101.89	122.70
1	B	334	ILE	CB-CA-C	-5.72	100.16	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	579	MET	Mainchain

## 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	3592	0	3309	63	0
1	B	3657	0	3420	54	0
2	A	10	0	12	1	0
2	B	10	0	12	1	0
3	A	21	0	0	0	0
3	B	21	0	0	0	0
All	All	7311	0	6753	108	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 8.

All (108) 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:305:ILE:O	1:B:330:ARG:NH2	1.83	1.10
1:A:143:SER:HB2	1:A:151:PHE:HB2	1.52	0.89
1:A:294:HIS:HD2	1:A:297:LEU:H	1.22	0.88
1:A:82:LEU:HD11	1:A:86:ASN:HD21	1.47	0.80
1:A:82:LEU:O	1:A:82:LEU:HD12	1.85	0.76
1:A:294:HIS:CD2	1:A:297:LEU:H	2.07	0.73
1:B:152:PHE:HB2	1:B:163:VAL:HB	1.70	0.72
1:A:235:ARG:NH2	1:A:580:ARG:O	2.20	0.69
1:A:82:LEU:HD11	1:A:86:ASN:ND2	2.07	0.69
1:A:340:PRO:HG3	1:B:294:HIS:CE1	2.28	0.68
1:A:340:PRO:HG3	1:B:294:HIS:HE1	1.59	0.67
1:A:508:LEU:HD23	1:A:508:LEU:C	2.14	0.67
1:B:502:LEU:HD11	1:B:553:LEU:HD11	1.77	0.67
1:A:508:LEU:HD23	1:A:508:LEU:O	1.95	0.66
1:A:82:LEU:CD1	1:A:86:ASN:ND2	2.59	0.66
1:A:82:LEU:CD1	1:A:86:ASN:HD21	2.12	0.63
1:A:491:ARG:HA	1:A:505:TYR:HB3	1.80	0.63
1:B:254:ARG:HG3	1:B:561:THR:HG21	1.81	0.62
1:A:152:PHE:HB2	1:A:163:VAL:HB	1.82	0.61
1:A:357:LEU:HD11	1:A:552:GLY:H	1.65	0.60
1:B:192:ILE:HG23	1:B:208:PRO:HB3	1.83	0.59

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:503:ASN:O	2:A:601:LYS:HD2	2.04	0.58
1:B:265:LEU:O	1:B:270:PHE:HB2	2.04	0.58
1:A:143:SER:CB	1:A:151:PHE:HB2	2.30	0.57
1:A:398:TYR:HB3	1:A:465:PRO:O	2.06	0.55
1:B:383:ASP:HB3	1:B:387:ASN:HD22	1.71	0.55
1:B:315:ILE:HD11	1:B:322:VAL:HG12	1.89	0.55
1:B:277:MET:HA	1:B:304:ARG:HD3	1.90	0.54
1:B:254:ARG:NH1	1:B:573:VAL:O	2.40	0.54
1:B:155:VAL:HG12	1:B:160:LYS:HB3	1.90	0.53
1:B:491:ARG:HG3	1:B:505:TYR:HD2	1.73	0.53
1:B:82:LEU:O	1:B:82:LEU:HG	2.09	0.53
1:B:429:ILE:HG23	1:B:439:LEU:HD11	1.90	0.53
1:B:253:THR:O	1:B:257:ILE:HG13	2.10	0.52
1:B:491:ARG:HA	1:B:505:TYR:HB3	1.92	0.51
1:A:548:PRO:HB3	1:B:189:ILE:HD13	1.92	0.51
1:A:213:LEU:HD21	1:A:216:ALA:HB2	1.93	0.51
1:A:84:PHE:HE1	1:A:218:LEU:CD1	2.23	0.51
1:A:81:ARG:O	1:A:84:PHE:HB3	2.10	0.51
1:B:470:GLU:HA	1:B:489:THR:O	2.11	0.51
1:B:136:THR:HG22	1:B:191:GLY:HA2	1.92	0.50
1:A:390:ILE:HD11	1:A:497:CYS:SG	2.52	0.50
1:A:347:PHE:CE2	1:A:553:LEU:HD22	2.47	0.49
1:A:482:HIS:HA	1:A:490:GLU:HG2	1.94	0.49
1:B:337:THR:HB	1:B:559:ARG:HH21	1.78	0.49
1:B:418:PRO:HG2	1:B:421:SER:HB3	1.93	0.49
1:A:400:LYS:N	1:A:400:LYS:HD2	2.27	0.48
1:B:375:THR:HG23	1:B:377:LYS:O	2.13	0.48
1:A:238:TYR:CE2	1:B:542:SER:HB3	2.48	0.48
1:A:544:GLU:HG2	1:B:84:PHE:HE1	1.80	0.47
1:A:302:TYR:CD1	1:A:302:TYR:N	2.83	0.47
1:B:305:ILE:O	1:B:330:ARG:CZ	2.59	0.47
1:A:407:ILE:HG13	1:A:457:ILE:HD11	1.97	0.46
1:B:314:LEU:HD22	1:B:319:ILE:HD12	1.97	0.46
1:A:162:GLN:O	1:A:206:ILE:N	2.42	0.46
1:A:348:TYR:HE2	1:A:507:GLU:OE2	1.98	0.46
1:A:509:ASN:HB3	1:A:547:LEU:HD23	1.98	0.46
1:A:314:LEU:HB3	1:A:319:ILE:HD12	1.96	0.46
1:B:272:GLU:HB2	1:B:323:TYR:CZ	2.51	0.46
1:A:460:LYS:HG2	1:A:461:TYR:CE1	2.51	0.46
1:B:149:LEU:HD12	1:B:166:ASN:ND2	2.30	0.46
1:A:108:ARG:HA	1:A:134:ASN:HB3	1.98	0.46

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:83:TYR:OH	1:B:185:ARG:HD2	2.16	0.46
1:B:494:MET:HB2	1:B:502:LEU:HD22	1.98	0.46
1:A:83:TYR:CE1	1:A:87:ARG:HG3	2.51	0.45
1:B:503:ASN:HD22	2:B:601:LYS:HD2	1.81	0.45
1:A:192:ILE:HG23	1:A:208:PRO:HB3	1.97	0.45
1:A:539:PHE:O	1:A:542:SER:OG	2.32	0.45
1:A:442:PRO:HD2	1:A:447:LYS:NZ	2.31	0.45
1:B:294:HIS:HD2	1:B:295:ASN:N	2.14	0.45
1:A:277:MET:HB2	1:B:341:GLU:OE1	2.16	0.45
1:A:334:ILE:HD11	1:B:294:HIS:CE1	2.52	0.44
1:B:84:PHE:CD1	1:B:219:HIS:CE1	3.05	0.44
1:A:541:THR:O	1:A:544:GLU:HB2	2.16	0.44
1:B:536:ASP:OD1	1:B:537:SER:N	2.51	0.44
1:B:401:VAL:O	1:B:468:ILE:HA	2.18	0.44
1:B:366:SER:HB2	1:B:376:TYR:CE2	2.53	0.43
1:B:362:GLU:OE2	1:B:400:LYS:NZ	2.52	0.43
1:B:142:VAL:O	1:B:142:VAL:HG23	2.19	0.43
1:B:398:TYR:HB3	1:B:465:PRO:O	2.19	0.43
1:B:406:GLU:O	1:B:410:VAL:HG23	2.19	0.42
1:A:265:LEU:O	1:A:270:PHE:HB2	2.19	0.42
1:A:271:PHE:HE1	1:B:255:THR:HG1	1.66	0.42
1:B:154:LEU:O	1:B:160:LYS:HA	2.19	0.42
1:A:84:PHE:HE1	1:A:218:LEU:HD13	1.85	0.42
1:A:206:ILE:O	1:A:208:PRO:HD3	2.20	0.42
1:A:307:THR:O	1:A:310:PRO:HD2	2.19	0.42
1:B:315:ILE:CD1	1:B:322:VAL:HG12	2.49	0.42
1:B:334:ILE:HG12	1:B:340:PRO:HD3	2.02	0.41
1:A:290:PHE:CE2	1:A:331:ASN:ND2	2.88	0.41
1:A:83:TYR:CD1	1:A:83:TYR:C	2.94	0.41
1:B:335:ASP:OD1	1:B:338:HIS:N	2.54	0.41
1:B:362:GLU:HG2	1:B:398:TYR:CD2	2.55	0.41
1:A:129:GLU:HA	1:A:195:PHE:CD2	2.55	0.41
1:A:254:ARG:NH1	1:A:573:VAL:O	2.53	0.41
1:A:357:LEU:CD1	1:A:552:GLY:H	2.32	0.41
1:B:170:HIS:CD2	1:B:177:PHE:HB2	2.56	0.41
1:A:339:ASN:HA	1:A:340:PRO:HD2	1.90	0.41
1:A:369:VAL:HG21	1:A:394:PHE:CD2	2.57	0.40
1:A:468:ILE:HD12	1:A:495:PHE:CE1	2.56	0.40
1:A:476:SER:HB3	1:A:479:ALA:HB3	2.04	0.40
1:A:418:PRO:HG2	1:A:421:SER:HB3	2.03	0.40
1:B:313:MET:O	1:B:316:VAL:HB	2.21	0.40

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:LEU:HD11	1:A:552:GLY:N	2.33	0.40
1:A:483:ARG:HB3	1:B:104:HIS:CD2	2.56	0.40
1:A:313:MET:O	1:A:316:VAL:HB	2.21	0.40
1:A:347:PHE:CD1	1:A:347:PHE:C	2.95	0.40
1:B:443:PRO:HB3	1:B:448:LEU:HD21	2.03	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	459/507 (90%)	440 (96%)	19 (4%)	0	100	100
1	B	464/507 (92%)	443 (96%)	21 (4%)	0	100	100
All	All	923/1014 (91%)	883 (96%)	40 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	359/457 (79%)	354 (99%)	5 (1%)	67	86
1	B	370/457 (81%)	361 (98%)	9 (2%)	49	77
All	All	729/914 (80%)	715 (98%)	14 (2%)	57	81

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	127	HIS
1	A	413	THR
1	A	492	LEU
1	A	508	LEU
1	A	558	ASP
1	B	82	LEU
1	B	151	PHE
1	B	164	LEU
1	B	279	ASN
1	B	307	THR
1	B	411	THR
1	B	413	THR
1	B	492	LEU
1	B	558	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	ASN
1	A	94	GLN
1	A	294	HIS
1	B	94	GLN
1	B	176	ASN
1	B	294	HIS
1	B	339	ASN
1	B	387	ASN
1	B	430	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 monosaccharides in this entry.

## 5.6 Ligand geometry

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	LYS	B	601	-	5,9,9	0.26	0	4,10,10	0.49	0
3	9CF	B	602	-	23,23,23	1.73	5 (21%)	33,33,33	2.41	11 (33%)
2	LYS	A	601	-	5,9,9	0.42	0	4,10,10	0.25	0
3	9CF	A	602	-	23,23,23	1.54	2 (8%)	33,33,33	2.02	12 (36%)

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	LYS	B	601	-	-	0/5/9/9	-
3	9CF	B	602	-	-	2/4/26/26	0/3/3/3
2	LYS	A	601	-	-	0/5/9/9	-
3	9CF	A	602	-	-	4/4/26/26	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	9CF	C6-C5	-6.37	1.39	1.51
3	B	602	9CF	C6-C5	-5.37	1.41	1.51
3	B	602	9CF	O5-C12	3.16	1.50	1.44
3	B	602	9CF	O5-C11	2.51	1.48	1.44
3	B	602	9CF	C15-C11	2.47	1.57	1.51
3	A	602	9CF	C9-C8	-2.42	1.40	1.46
3	B	602	9CF	C9-C8	-2.00	1.41	1.46

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	9CF	C7-O3-C8	-7.16	107.66	118.69
3	B	602	9CF	O3-C7-C10	5.56	115.02	106.17
3	A	602	9CF	O3-C8-O4	4.63	122.76	117.60
3	B	602	9CF	O3-C8-O4	4.34	122.44	117.60
3	A	602	9CF	C14-C13-C12	-3.71	108.47	111.04
3	A	602	9CF	O5-C11-C15	3.69	113.82	109.86
3	B	602	9CF	O5-C11-C15	3.61	113.74	109.86
3	A	602	9CF	C10-C7-C6	-3.51	106.83	113.44
3	A	602	9CF	C10-C11-C15	-3.38	106.03	113.83
3	B	602	9CF	O3-C8-C9	-3.38	114.39	119.09
3	B	602	9CF	C1-C9-C8	3.18	123.55	119.81
3	A	602	9CF	C6-C5-C9	2.79	122.78	118.03
3	A	602	9CF	O3-C8-C9	-2.75	115.26	119.09
3	B	602	9CF	O5-C12-C16	2.64	112.48	106.88
3	A	602	9CF	O5-C11-C10	-2.62	101.49	107.70
3	B	602	9CF	C10-C11-C15	-2.54	107.97	113.83
3	B	602	9CF	C14-C13-C12	-2.53	109.29	111.04
3	A	602	9CF	O1-C1-C9	-2.52	116.42	121.14
3	B	602	9CF	C6-C5-C9	2.44	122.19	118.03
3	A	602	9CF	O3-C7-C6	2.25	113.08	110.20
3	A	602	9CF	C12-O5-C11	2.08	118.28	114.00
3	B	602	9CF	C16-C12-C13	-2.08	107.44	114.42
3	A	602	9CF	O3-C7-C10	2.07	109.46	106.17

There are no chirality outliers.

All (6) torsion outliers are listed below:

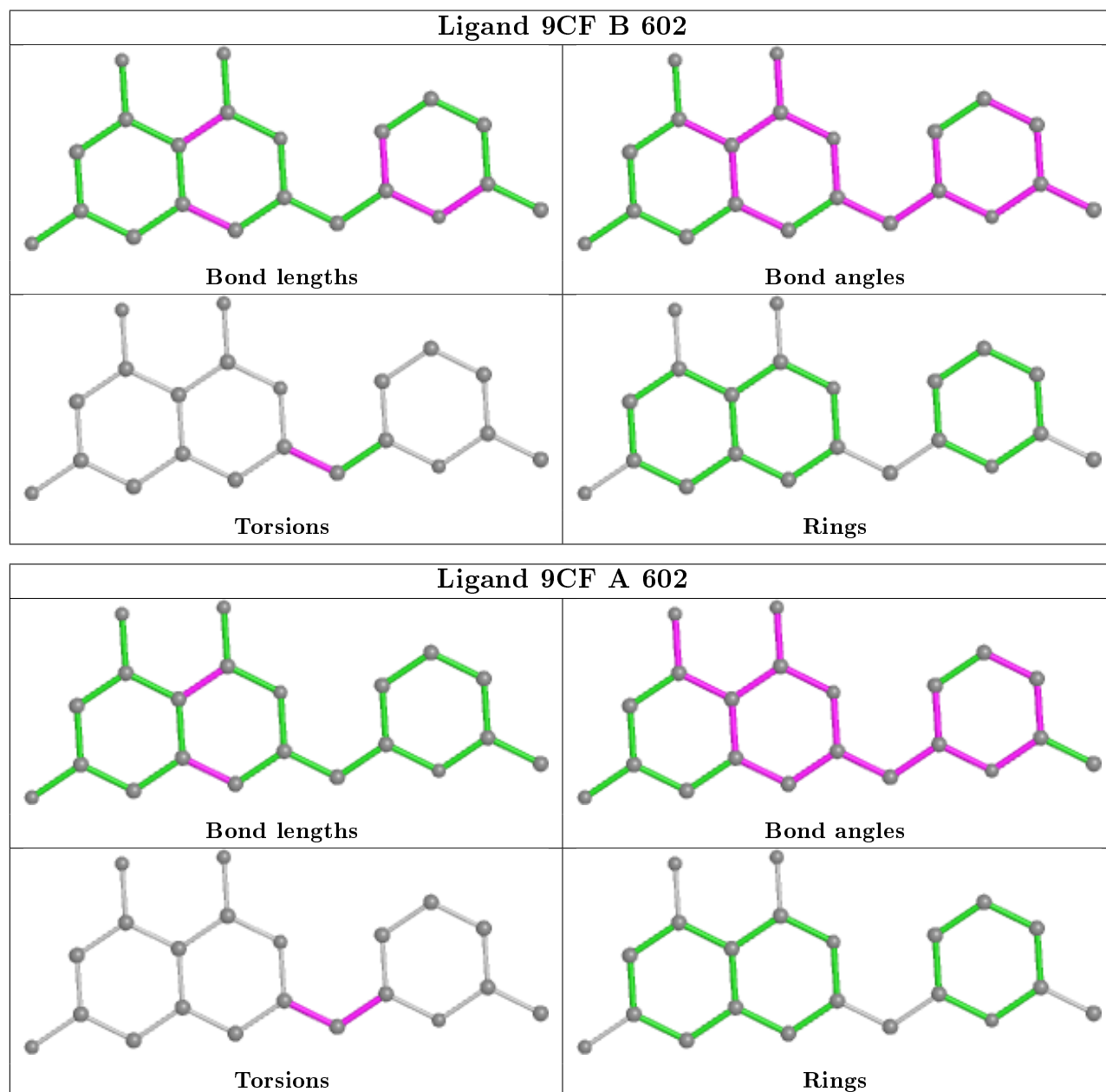
Mol	Chain	Res	Type	Atoms
3	B	602	9CF	C11-C10-C7-O3
3	A	602	9CF	C7-C10-C11-C15
3	A	602	9CF	C7-C10-C11-O5
3	A	602	9CF	C11-C10-C7-O3
3	B	602	9CF	C11-C10-C7-C6
3	A	602	9CF	C11-C10-C7-C6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	LYS	1	0
2	A	601	LYS	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

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	469/507 (92%)	0.04	11 (2%)	60	57	30, 109, 151, 166	0
1	B	474/507 (93%)	0.07	11 (2%)	60	57	70, 106, 145, 176	0
All	All	943/1014 (92%)	0.06	22 (2%)	60	57	30, 107, 149, 176	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	285	ALA	3.5
1	B	419	PHE	3.3
1	B	446	ALA	2.9
1	B	445	ALA	2.9
1	B	432	ILE	2.9
1	B	415	LEU	2.8
1	B	348	TYR	2.6
1	B	232	ILE	2.5
1	A	154	LEU	2.5
1	A	161	ILE	2.5
1	B	488	LEU	2.5
1	A	181	TYR	2.5
1	A	286	ASN	2.4
1	B	457	ILE	2.3
1	A	123	GLY	2.3
1	A	151	PHE	2.3
1	A	455	HIS	2.3
1	A	165	ALA	2.3
1	A	196	PRO	2.3
1	B	542	SER	2.2
1	A	152	PHE	2.0
1	B	392	ILE	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

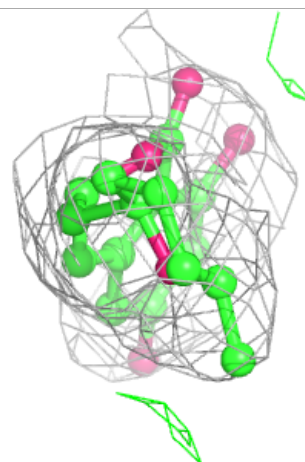
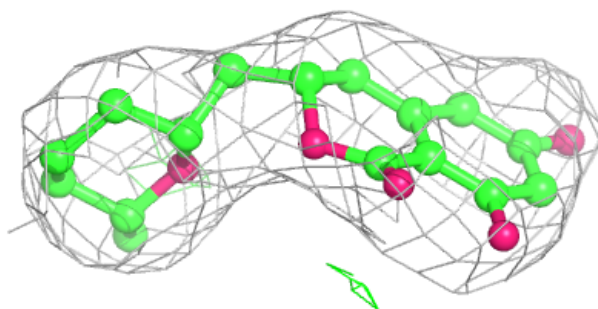
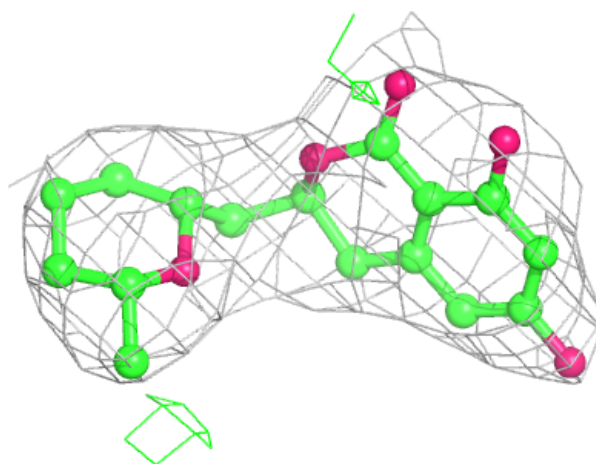
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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	LYS	B	601	10/10	0.90	0.37	59,62,66,67	10
2	LYS	A	601	10/10	0.91	0.36	48,52,53,58	10
3	9CF	A	602	21/21	0.94	0.20	83,103,119,120	0
3	9CF	B	602	21/21	0.95	0.22	68,90,102,105	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.

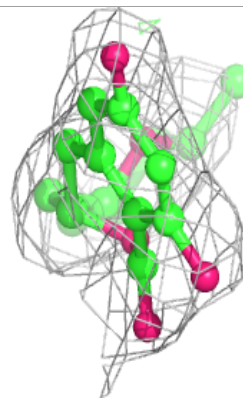
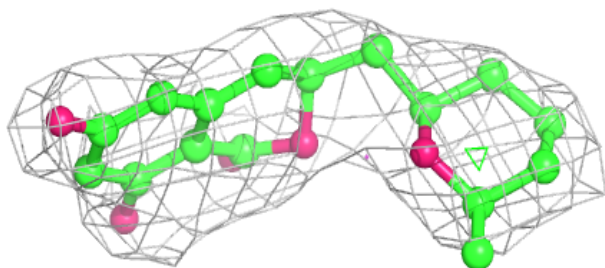
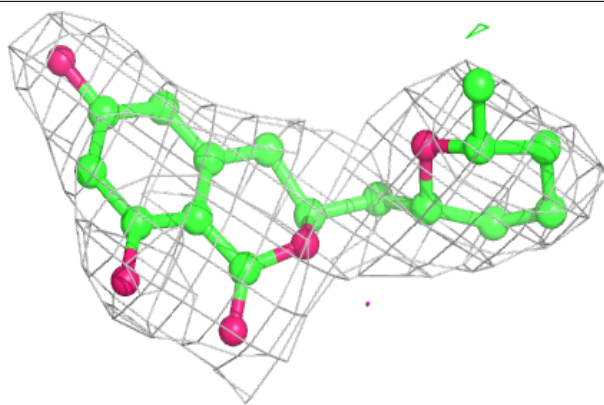
**Electron density around 9CF A 602:**

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



**Electron density around 9CF B 602:**

$2mF_o-DF_c$  (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.