



# Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 09:38 am BST

PDB ID : 5UKF  
Title : Crystal Structure of the Human Vaccinia-related Kinase 1 Bound to an Oxindole Inhibitor  
Authors : Counago, R.M.; Wells, C.; Zuercher, W.; Willson, T.M.; Bountra, C.; Edwards, A.M.; Arruda, P.; Gileadi, O.; Structural Genomics Consortium (SGC)  
Deposited on : 2017-01-22  
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](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 : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

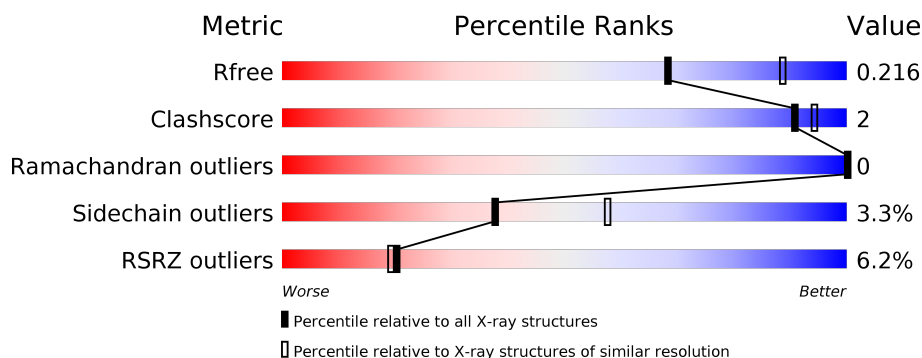
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	363	<div> <div>7%</div> <div> <div></div> <div>85%</div> <div>11%</div> </div> </div>
1	B	363	<div> <div>4%</div> <div> <div></div> <div>68%</div> <div>5%</div> <div>27%</div> </div> </div>
1	C	363	<div> <div>4%</div> <div> <div></div> <div>81%</div> <div>7%</div> <div>13%</div> </div> </div>
1	D	363	<div> <div>5%</div> <div> <div></div> <div>70%</div> <div>26%</div> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9854 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 Serine/threonine-protein kinase VRK1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	0	0
			2473	1586	420	455	12			
1	B	265	Total	C	N	O	S	0	0	0
			2073	1331	352	381	9			
1	C	317	Total	C	N	O	S	0	0	0
			2478	1589	418	459	12			
1	D	268	Total	C	N	O	S	0	0	0
			2130	1369	365	387	9			

There are 48 discrepancies between the modelled and reference sequences:

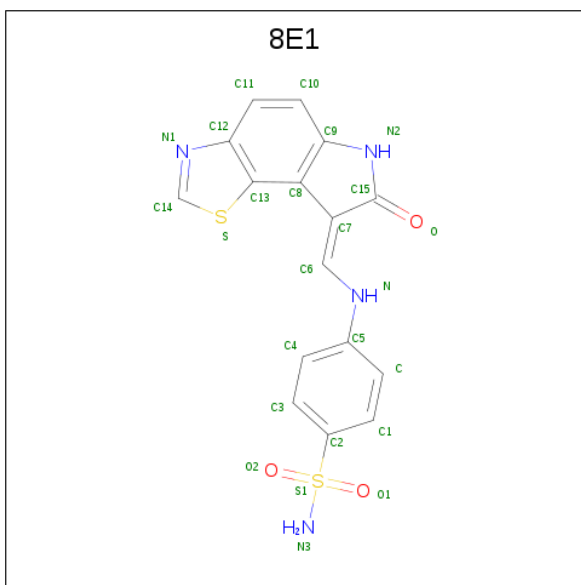
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	MET	-	initiating methionine	UNP Q99986
A	34	ALA	LYS	engineered mutation	UNP Q99986
A	35	ALA	LYS	engineered mutation	UNP Q99986
A	36	ALA	GLU	engineered mutation	UNP Q99986
A	212	ALA	GLU	engineered mutation	UNP Q99986
A	214	ALA	LYS	engineered mutation	UNP Q99986
A	215	ALA	GLU	engineered mutation	UNP Q99986
A	292	ALA	GLU	engineered mutation	UNP Q99986
A	293	ALA	LYS	engineered mutation	UNP Q99986
A	295	ALA	LYS	engineered mutation	UNP Q99986
A	359	ALA	LYS	engineered mutation	UNP Q99986
A	360	ALA	LYS	engineered mutation	UNP Q99986
B	2	MET	-	initiating methionine	UNP Q99986
B	34	ALA	LYS	engineered mutation	UNP Q99986
B	35	ALA	LYS	engineered mutation	UNP Q99986
B	36	ALA	GLU	engineered mutation	UNP Q99986
B	212	ALA	GLU	engineered mutation	UNP Q99986
B	214	ALA	LYS	engineered mutation	UNP Q99986
B	215	ALA	GLU	engineered mutation	UNP Q99986
B	292	ALA	GLU	engineered mutation	UNP Q99986
B	293	ALA	LYS	engineered mutation	UNP Q99986

*Continued on next page...*

*Continued from previous page...*

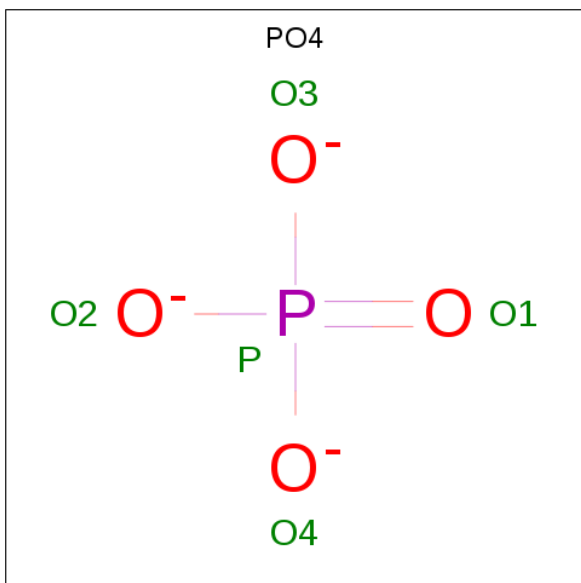
Chain	Residue	Modelled	Actual	Comment	Reference
B	295	ALA	LYS	engineered mutation	UNP Q99986
B	359	ALA	LYS	engineered mutation	UNP Q99986
B	360	ALA	LYS	engineered mutation	UNP Q99986
C	2	MET	-	initiating methionine	UNP Q99986
C	34	ALA	LYS	engineered mutation	UNP Q99986
C	35	ALA	LYS	engineered mutation	UNP Q99986
C	36	ALA	GLU	engineered mutation	UNP Q99986
C	212	ALA	GLU	engineered mutation	UNP Q99986
C	214	ALA	LYS	engineered mutation	UNP Q99986
C	215	ALA	GLU	engineered mutation	UNP Q99986
C	292	ALA	GLU	engineered mutation	UNP Q99986
C	293	ALA	LYS	engineered mutation	UNP Q99986
C	295	ALA	LYS	engineered mutation	UNP Q99986
C	359	ALA	LYS	engineered mutation	UNP Q99986
C	360	ALA	LYS	engineered mutation	UNP Q99986
D	2	MET	-	initiating methionine	UNP Q99986
D	34	ALA	LYS	engineered mutation	UNP Q99986
D	35	ALA	LYS	engineered mutation	UNP Q99986
D	36	ALA	GLU	engineered mutation	UNP Q99986
D	212	ALA	GLU	engineered mutation	UNP Q99986
D	214	ALA	LYS	engineered mutation	UNP Q99986
D	215	ALA	GLU	engineered mutation	UNP Q99986
D	292	ALA	GLU	engineered mutation	UNP Q99986
D	293	ALA	LYS	engineered mutation	UNP Q99986
D	295	ALA	LYS	engineered mutation	UNP Q99986
D	359	ALA	LYS	engineered mutation	UNP Q99986
D	360	ALA	LYS	engineered mutation	UNP Q99986

- Molecule 2 is 4-{{[(Z)-(7-oxo-6,7-dihydro-8H-[1,3]thiazolo[5,4-e]indol-8-ylidene)methyl]amino}benzene-1-sulfonamide (three-letter code: 8E1) (formula: C<sub>16</sub>H<sub>12</sub>N<sub>4</sub>O<sub>3</sub>S<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			25	16	4	3	2		
2	B	1	Total	C	N	O	S	0	0
			25	16	4	3	2		
2	C	1	Total	C	N	O	S	0	0
			25	16	4	3	2		
2	D	1	Total	C	N	O	S	0	0
			25	16	4	3	2		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).

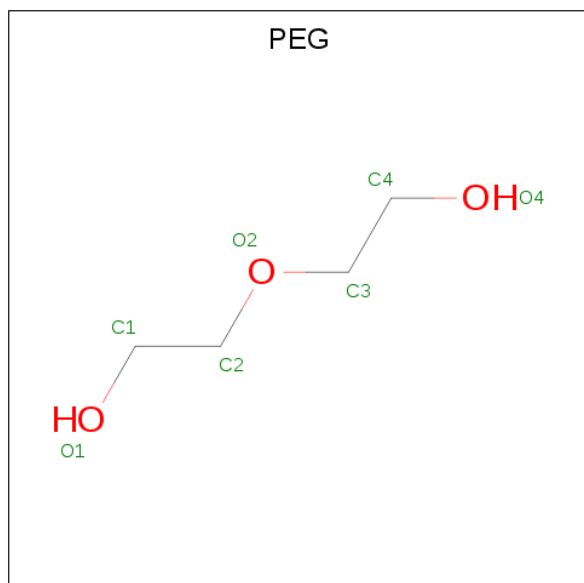


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0
3	B	1	Total O P 5 4 1	0	0
3	B	1	Total O P 5 4 1	0	0
3	B	1	Total O P 5 4 1	0	0
3	B	1	Total O P 5 4 1	0	0
3	C	1	Total O P 5 4 1	0	0
3	C	1	Total O P 5 4 1	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Cl 1 1	0	0
4	A	1	Total Cl 1 1	0	0
4	D	4	Total Cl 4 4	0	0
4	C	1	Total Cl 1 1	0	0

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is water.

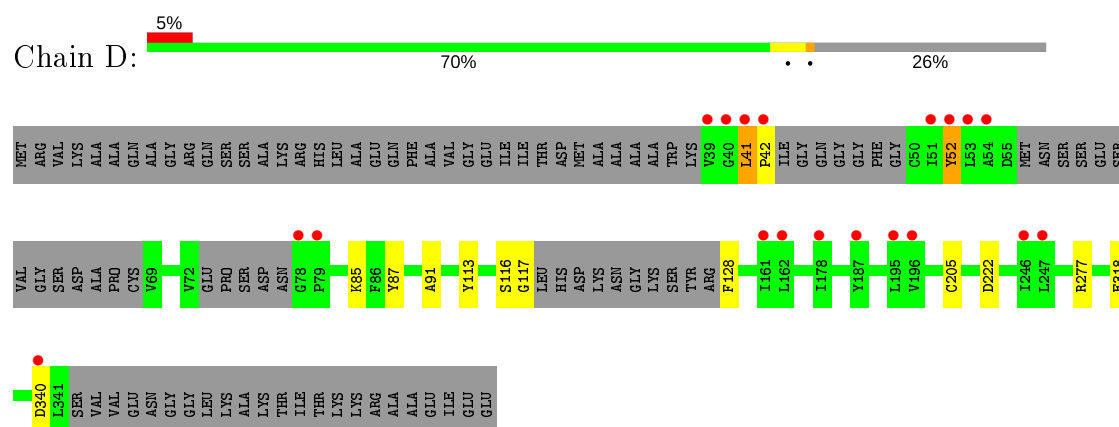
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	132	Total	O	0	0
			132	132		
6	B	122	Total	O	0	0
			122	122		
6	C	141	Total	O	0	0
			141	141		
6	D	156	Total	O	0	0
			156	156		



- Molecule 1: Serine/threonine-protein kinase VRK1







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	92.14Å 96.05Å 192.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.19 – 2.40 29.56 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (24.19-2.40) 100.0 (29.56-2.40)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.15 (at 2.39Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, $R_{free}$	0.177 , 0.214 0.187 , 0.216	Depositor DCC
$R_{free}$ test set	3417 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.7	Xtriage
Anisotropy	0.576	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 56.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.022 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9854	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, PEG, 8E1, CL

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	0.50	0/2534	0.65	0/3444
1	B	0.48	0/2120	0.61	0/2875
1	C	0.49	0/2538	0.61	0/3446
1	D	0.52	0/2179	0.65	0/2946
All	All	0.50	0/9371	0.63	0/12711

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 [i](#)

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	2473	0	2352	7	0
1	B	2073	0	1972	10	0
1	C	2478	0	2383	9	0
1	D	2130	0	2059	4	0
2	A	25	0	0	0	0
2	B	25	0	0	0	0
2	C	25	0	0	0	0
2	D	25	0	0	0	0
3	A	5	0	0	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	20	0	0	0	0
3	C	10	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	4	0	0	0	0
5	D	7	0	10	0	0
6	A	132	0	0	0	0
6	B	122	0	0	0	0
6	C	141	0	0	0	0
6	D	156	0	0	0	0
All	All	9854	0	8776	30	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 2.

All (30) 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:109:GLY:HA3	1:B:164:ILE:HD11	1.84	0.60
1:C:185:LEU:HD11	1:C:341:LEU:HD12	1.89	0.55
1:C:169:HIS:CG	1:C:240:ARG:HG2	2.47	0.50
1:C:204:TYR:CE2	1:C:205:CYS:HB2	2.47	0.49
1:B:138:LEU:CD2	1:B:157:LEU:HD21	2.43	0.49
1:B:161:ILE:HG21	1:B:247:LEU:HD13	1.95	0.48
1:B:204:TYR:CE2	1:B:205:CYS:HB2	2.49	0.48
1:D:41:LEU:H	1:D:41:LEU:HD12	1.79	0.47
1:B:155:LEU:HD21	1:B:298:GLU:HG3	1.97	0.47
1:A:138:LEU:HA	1:A:138:LEU:HD23	1.67	0.46
1:C:174:VAL:HG13	1:C:204:TYR:CD1	2.50	0.46
1:B:138:LEU:HD21	1:B:157:LEU:HD21	1.96	0.46
1:A:117:GLY:O	1:A:127:ARG:HD3	2.16	0.45
1:B:109:GLY:HA3	1:B:164:ILE:CD1	2.46	0.45
1:C:85:LYS:HD2	1:C:85:LYS:HA	1.54	0.45
1:D:117:GLY:HA3	1:D:128:PHE:CE2	2.51	0.45
1:A:131:MET:HB3	1:A:131:MET:HE3	1.70	0.45
1:C:185:LEU:HD11	1:C:341:LEU:CD1	2.47	0.45
1:D:87:TYR:HA	1:D:91:ALA:HB3	1.99	0.44
1:A:85:LYS:HA	1:A:85:LYS:HD2	1.65	0.44
1:B:227:PHE:O	1:B:246:ILE:HG12	2.18	0.44
1:A:77:ASN:ND2	1:A:79:PRO:HD2	2.33	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:189:ASN:HA	1:C:190:PRO:HD2	1.93	0.43
1:B:169:HIS:CG	1:B:240:ARG:HG2	2.53	0.43
1:B:251:MET:HE3	1:B:251:MET:HB2	1.92	0.43
1:C:52:TYR:HB2	1:C:70:VAL:HG13	2.01	0.42
1:A:78:GLY:N	1:A:79:PRO:CD	2.83	0.42
1:C:155:LEU:HD21	1:C:298:GLU:HG3	2.02	0.42
1:A:205:CYS:O	1:A:205:CYS:SG	2.78	0.41
1:D:42:PRO:HA	1:D:52:TYR:HA	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	320/363 (88%)	315 (98%)	5 (2%)	0	100	100
1	B	257/363 (71%)	253 (98%)	4 (2%)	0	100	100
1	C	313/363 (86%)	307 (98%)	6 (2%)	0	100	100
1	D	258/363 (71%)	253 (98%)	5 (2%)	0	100	100
All	All	1148/1452 (79%)	1128 (98%)	20 (2%)	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	244/303 (80%)	238 (98%)	6 (2%)	47	67
1	B	206/303 (68%)	202 (98%)	4 (2%)	57	75
1	C	253/303 (84%)	243 (96%)	10 (4%)	31	49
1	D	215/303 (71%)	205 (95%)	10 (5%)	26	42
All	All	918/1212 (76%)	888 (97%)	30 (3%)	38	57

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	SER
1	A	76	ASP
1	A	80	LEU
1	A	125	SER
1	A	140	LYS
1	A	274	SER
1	B	113	TYR
1	B	132	ASP
1	B	157	LEU
1	B	225	ILE
1	C	38	LYS
1	C	48	PHE
1	C	51	ILE
1	C	80	LEU
1	C	84	LEU
1	C	112	LYS
1	C	133	ARG
1	C	187	TYR
1	C	221	HIS
1	C	228	THR
1	D	41	LEU
1	D	52	TYR
1	D	85	LYS
1	D	113	TYR
1	D	116	SER
1	D	205	CYS
1	D	222	ASP
1	D	277	ARG
1	D	318	GLU
1	D	340	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 ⓘ

Of 19 ligands modelled in this entry, 7 are monoatomic - leaving 12 for Mogul analysis.

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	8E1	A	401	-	24,28,28	0.88	2 (8%)	33,42,42	1.11	2 (6%)
2	8E1	B	401	-	24,28,28	0.82	1 (4%)	33,42,42	0.69	1 (3%)
3	PO4	C	403	-	4,4,4	2.49	1 (25%)	6,6,6	0.49	0
2	8E1	D	402	-	24,28,28	0.85	1 (4%)	33,42,42	0.81	1 (3%)
3	PO4	B	402	-	4,4,4	2.54	2 (50%)	6,6,6	0.45	0
5	PEG	D	401	-	6,6,6	0.33	0	5,5,5	0.51	0
3	PO4	B	405	-	4,4,4	1.18	0	6,6,6	0.45	0
3	PO4	C	402	-	4,4,4	2.43	1 (25%)	6,6,6	0.60	0
3	PO4	B	403	-	4,4,4	2.48	1 (25%)	6,6,6	0.49	0
3	PO4	B	404	-	4,4,4	2.49	1 (25%)	6,6,6	0.57	0
3	PO4	A	402	-	4,4,4	2.49	1 (25%)	6,6,6	0.58	0
2	8E1	C	401	-	24,28,28	0.85	1 (4%)	33,42,42	1.04	2 (6%)

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	8E1	A	401	-	-	2/11/23/23	0/4/4/4
2	8E1	B	401	-	-	5/11/23/23	0/4/4/4
2	8E1	D	402	-	-	3/11/23/23	0/4/4/4
5	PEG	D	401	-	-	0/4/4/4	-
2	8E1	C	401	-	-	5/11/23/23	0/4/4/4

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	PO4	P-O1	4.14	1.60	1.50
3	B	403	PO4	P-O1	4.11	1.60	1.50
3	A	402	PO4	P-O1	4.11	1.60	1.50
3	C	403	PO4	P-O1	4.09	1.60	1.50
3	B	404	PO4	P-O1	4.09	1.60	1.50
3	C	402	PO4	P-O1	4.02	1.60	1.50
2	C	401	8E1	C11-C10	2.99	1.42	1.36
2	B	401	8E1	C11-C10	2.97	1.42	1.36
2	A	401	8E1	C11-C10	2.92	1.42	1.36
2	D	402	8E1	C11-C10	2.91	1.42	1.36
2	A	401	8E1	C7-C15	-2.01	1.47	1.50
3	B	402	PO4	P-O2	2.01	1.60	1.54

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	8E1	C7-C6-N	-4.74	115.27	123.92
2	C	401	8E1	C7-C6-N	-4.30	116.08	123.92
2	D	402	8E1	C7-C6-N	-2.37	119.60	123.92
2	B	401	8E1	C14-S-C13	-2.24	88.44	90.88
2	C	401	8E1	C14-S-C13	-2.19	88.49	90.88
2	A	401	8E1	C6-C7-C15	-2.16	121.52	124.74

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	402	8E1	N-C6-C7-C8
2	D	402	8E1	N-C6-C7-C15
2	C	401	8E1	C7-C6-N-C5

*Continued on next page...*



*Continued from previous page...*

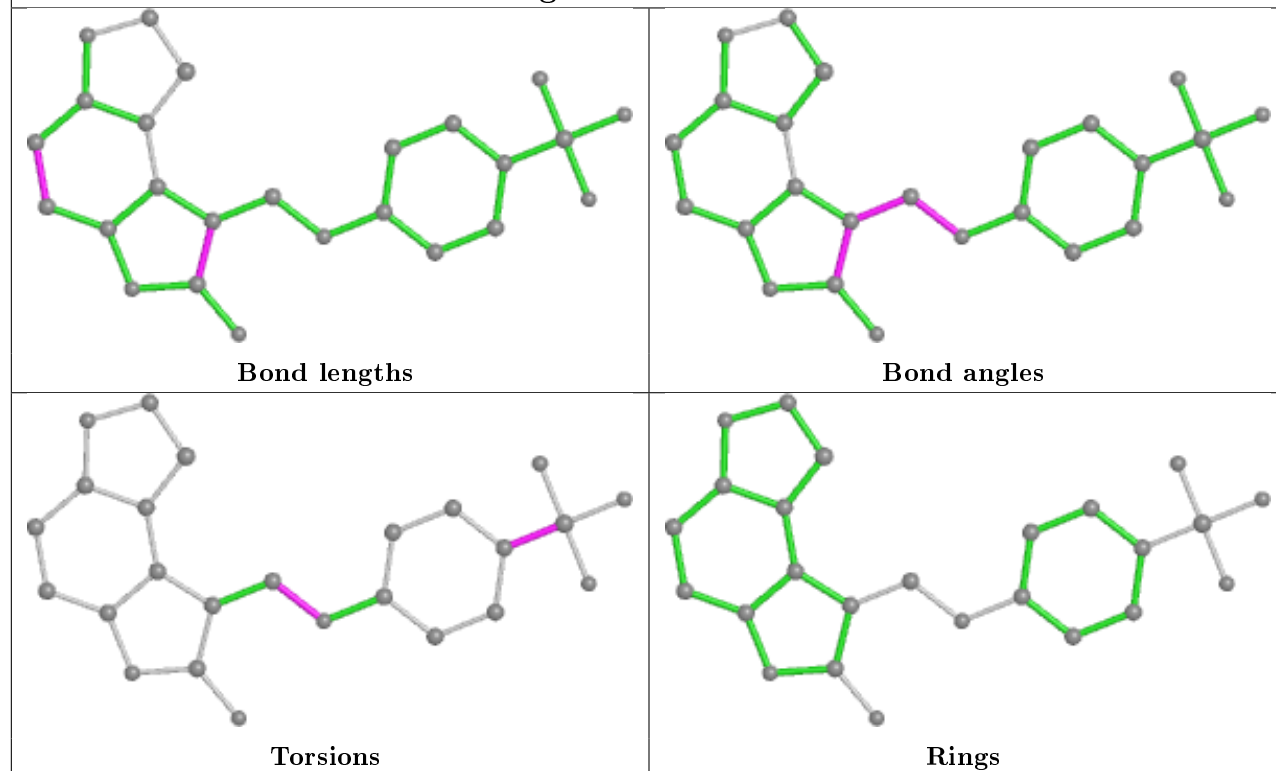
Mol	Chain	Res	Type	Atoms
2	D	402	8E1	C7-C6-N-C5
2	C	401	8E1	C1-C2-S1-N3
2	C	401	8E1	C1-C2-S1-O1
2	C	401	8E1	C3-C2-S1-N3
2	A	401	8E1	C7-C6-N-C5
2	B	401	8E1	C1-C2-S1-N3
2	B	401	8E1	C3-C2-S1-N3
2	C	401	8E1	C3-C2-S1-O1
2	B	401	8E1	C7-C6-N-C5
2	B	401	8E1	C3-C2-S1-O2
2	B	401	8E1	C1-C2-S1-O2
2	A	401	8E1	C1-C2-S1-O1

There are no ring outliers.

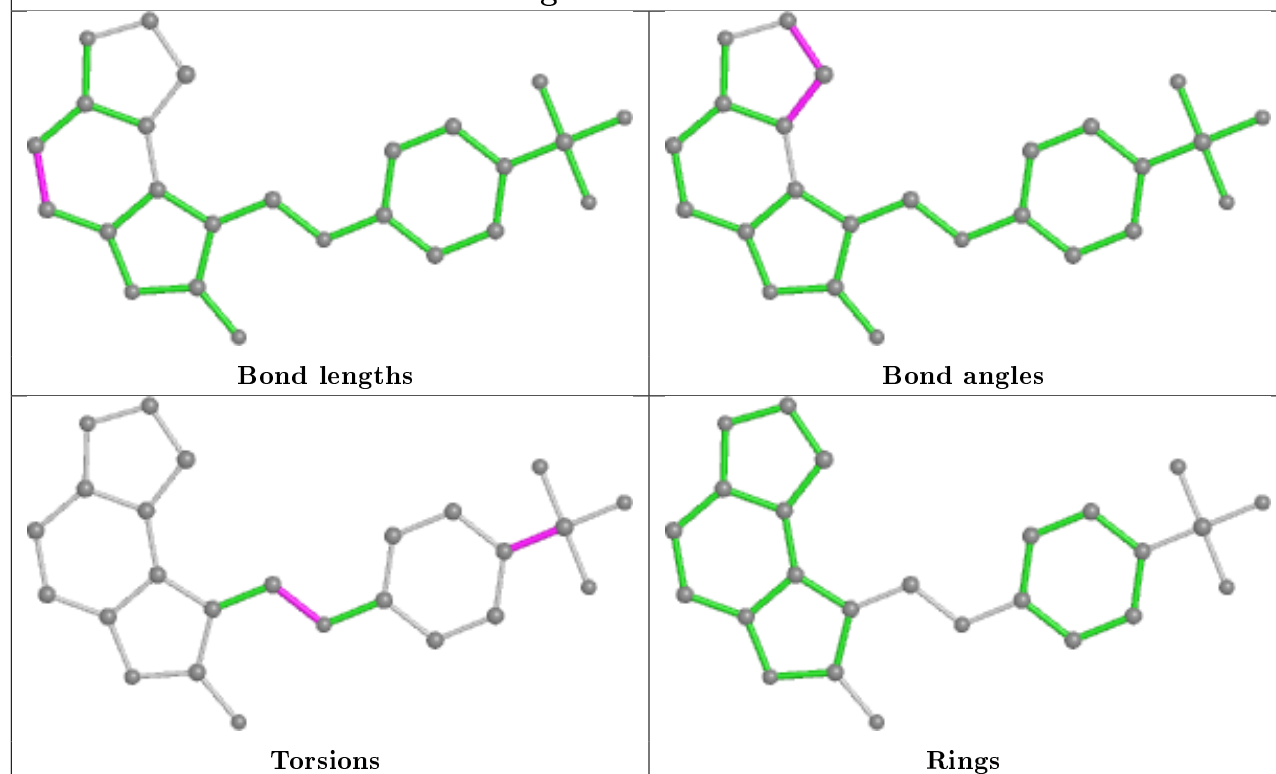
No monomer is involved in short contacts.

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.

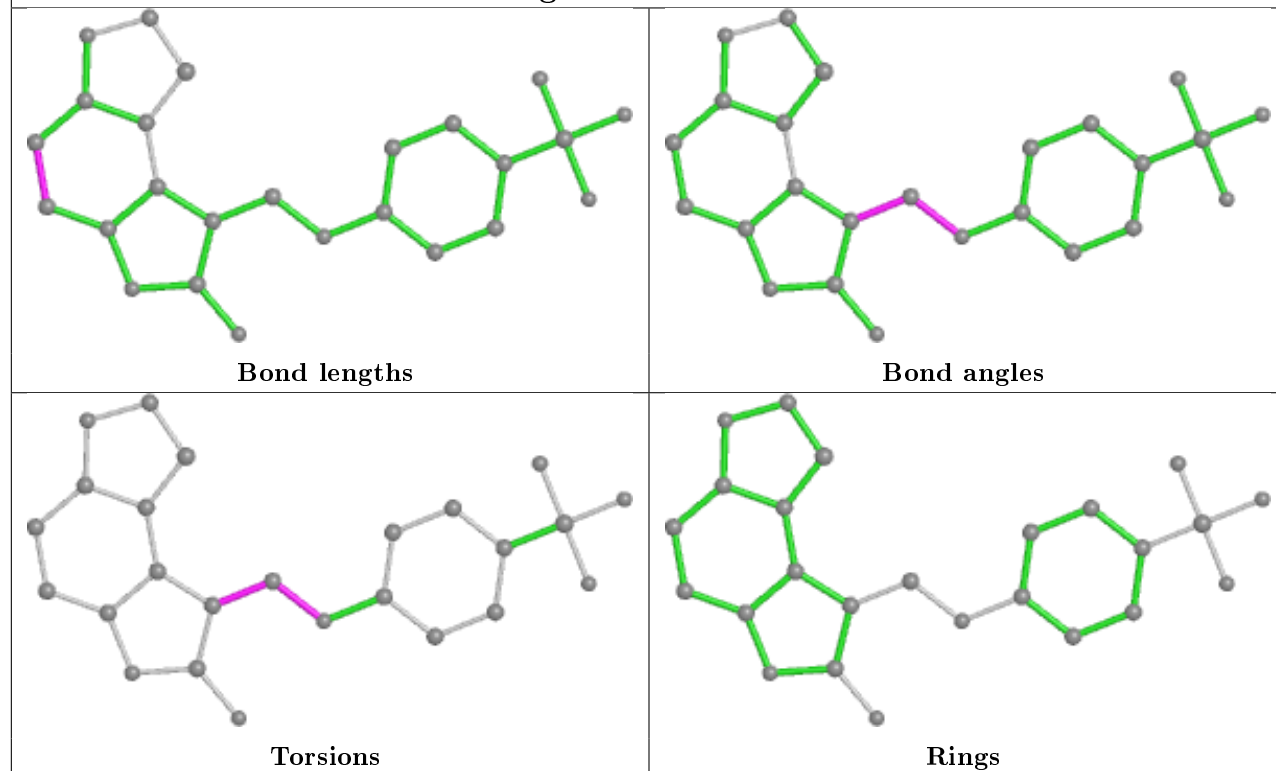
## Ligand 8E1 A 401



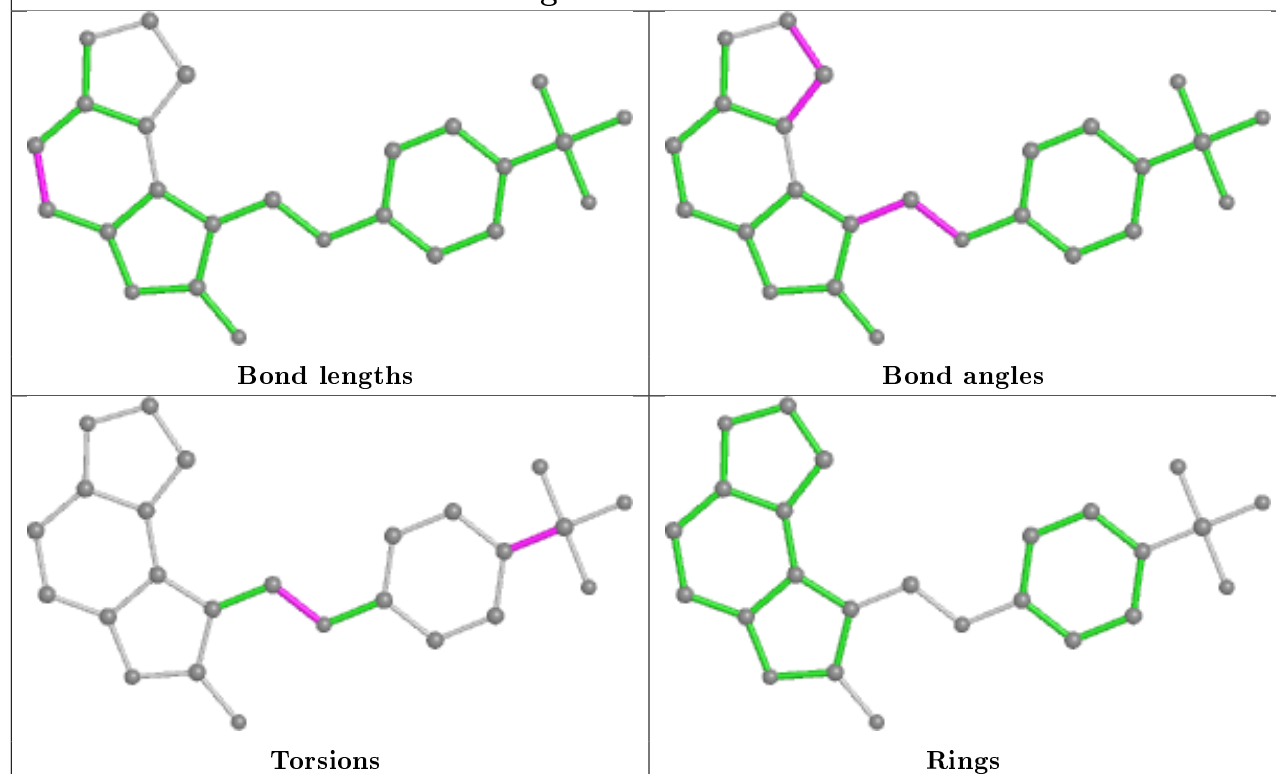
## Ligand 8E1 B 401



## Ligand 8E1 D 402



## Ligand 8E1 C 401



## 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	322/363 (88%)	0.11	24 (7%)	14 13	37, 57, 90, 119	0
1	B	265/363 (73%)	0.12	15 (5%)	23 22	36, 58, 109, 153	0
1	C	317/363 (87%)	0.15	15 (4%)	31 30	35, 58, 100, 119	0
1	D	268/363 (73%)	0.07	19 (7%)	16 14	33, 50, 84, 138	0
All	All	1172/1452 (80%)	0.11	73 (6%)	20 19	33, 56, 99, 153	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	64	SER	5.9
1	A	217	PRO	5.4
1	C	49	GLY	5.4
1	C	48	PHE	5.0
1	D	41	LEU	4.9
1	B	131	MET	4.8
1	A	20	ALA	4.8
1	C	218	LYS	4.4
1	C	217	PRO	4.3
1	A	209	VAL	4.3
1	D	54	ALA	4.2
1	B	52	TYR	4.0
1	A	63	GLY	3.7
1	D	79	PRO	3.6
1	B	114	TRP	3.6
1	D	161	ILE	3.6
1	B	117	GLY	3.5
1	A	218	LYS	3.5
1	C	63	GLY	3.5
1	B	135	GLY	3.4
1	A	292	ALA	3.4

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	C	41	LEU	3.3
1	A	76	ASP	3.2
1	D	52	TYR	3.1
1	D	40	GLY	3.1
1	C	76	ASP	3.1
1	A	25	VAL	3.1
1	D	187	TYR	3.1
1	B	118	LEU	3.1
1	C	50	CYS	3.0
1	C	187	TYR	3.0
1	D	53	LEU	3.0
1	D	78	GLY	3.0
1	A	215	ALA	3.0
1	D	39	VAL	2.9
1	D	246	ILE	2.9
1	A	47	GLY	2.8
1	C	247	LEU	2.7
1	B	128	PHE	2.7
1	D	42	PRO	2.7
1	B	247	LEU	2.6
1	A	46	GLY	2.6
1	D	178	ILE	2.5
1	A	74	PRO	2.5
1	A	144	ALA	2.5
1	C	215	ALA	2.5
1	B	340	ASP	2.4
1	B	165	LEU	2.4
1	B	164	ILE	2.3
1	A	214	ALA	2.3
1	D	195	LEU	2.3
1	B	209	VAL	2.3
1	A	73	GLU	2.3
1	C	161	ILE	2.3
1	A	161	ILE	2.2
1	A	291	PRO	2.2
1	C	104	LYS	2.1
1	D	340	ASP	2.1
1	D	51	ILE	2.1
1	D	247	LEU	2.1
1	A	195	LEU	2.1
1	B	78	GLY	2.1
1	A	190	PRO	2.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	329	LYS	2.1
1	A	178	ILE	2.1
1	D	162	LEU	2.1
1	A	77	ASN	2.1
1	A	45	GLN	2.1
1	B	187	TYR	2.0
1	C	178	ILE	2.0
1	A	216	ASP	2.0
1	A	336	ASP	2.0
1	D	196	VAL	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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(Å <sup>2</sup> )	Q<0.9
2	8E1	C	401	25/25	0.71	0.28	85,90,122,123	0
3	PO4	C	403	5/5	0.72	0.29	133,135,135,136	0
5	PEG	D	401	7/7	0.73	0.38	61,71,73,74	0
4	CL	D	403	1/1	0.77	0.29	81,81,81,81	0
2	8E1	B	401	25/25	0.77	0.28	87,93,126,127	0
2	8E1	D	402	25/25	0.79	0.22	58,69,114,115	0
4	CL	C	404	1/1	0.82	0.21	102,102,102,102	0
3	PO4	B	403	5/5	0.83	0.21	134,135,135,136	0
2	8E1	A	401	25/25	0.87	0.18	50,54,93,95	0
3	PO4	B	404	5/5	0.89	0.22	103,104,106,106	0
4	CL	A	403	1/1	0.90	0.07	89,89,89,89	0
4	CL	D	406	1/1	0.91	0.37	82,82,82,82	0
3	PO4	A	402	5/5	0.92	0.15	92,96,97,99	0

*Continued on next page...*

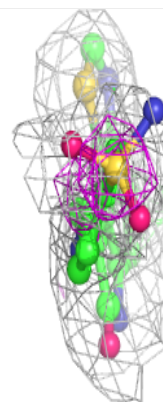
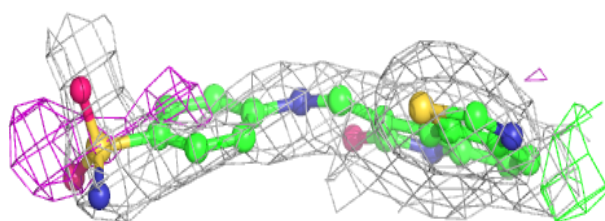
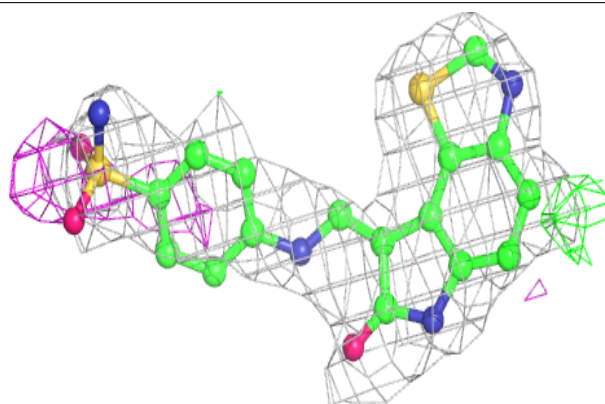
*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	CL	B	406	1/1	0.93	0.38	86,86,86,86	0
3	PO4	B	405	5/5	0.93	0.29	103,104,107,107	0
3	PO4	C	402	5/5	0.94	0.16	69,70,72,78	0
3	PO4	B	402	5/5	0.96	0.09	98,98,100,100	0
4	CL	D	405	1/1	0.97	0.23	86,86,86,86	0
4	CL	D	404	1/1	0.98	0.23	72,72,72,72	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 8E1 C 401:**

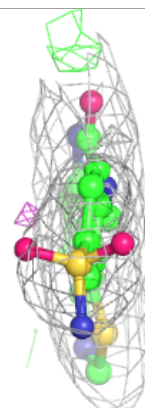
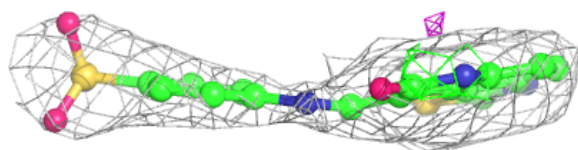
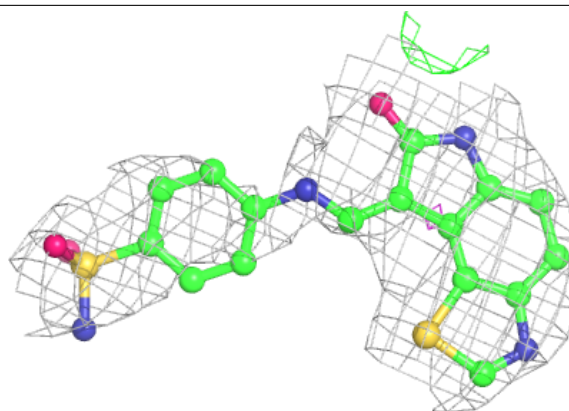
2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



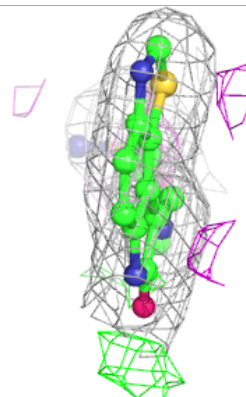
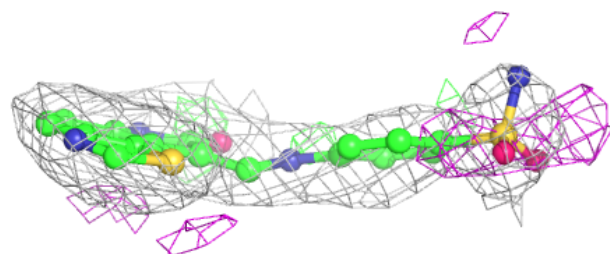
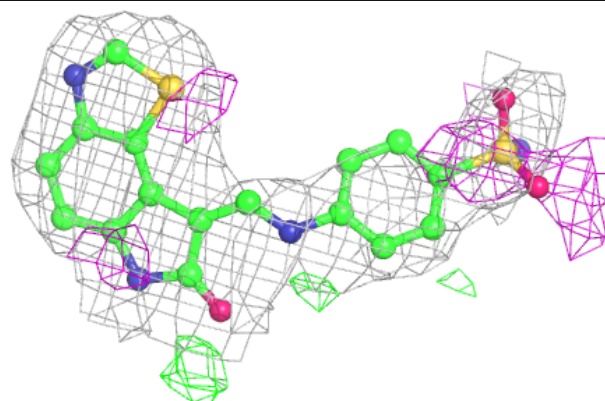


**Electron density around 8E1 B 401:**

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

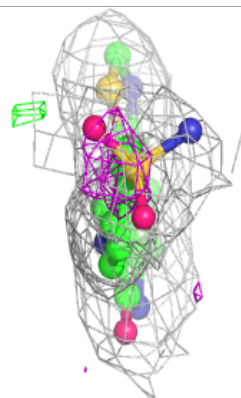
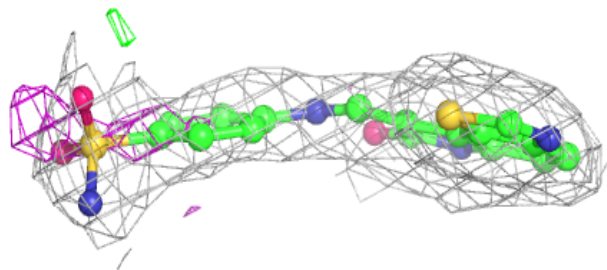
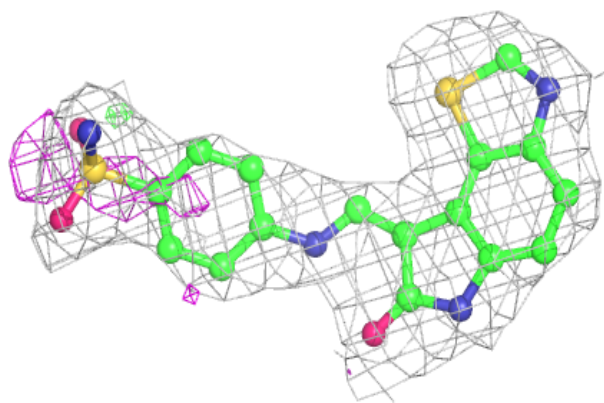
**Electron density around 8E1 D 402:**

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



**Electron density around 8E1 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.