



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 10:00 am BST

PDB ID : 2Y4O  
Title : Crystal Structure of PaaK2 in complex with phenylacetyl adenylate  
Authors : Law, A.; Boulanger, M.J.  
Deposited on : 2011-01-07  
Resolution : 1.90 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.

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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

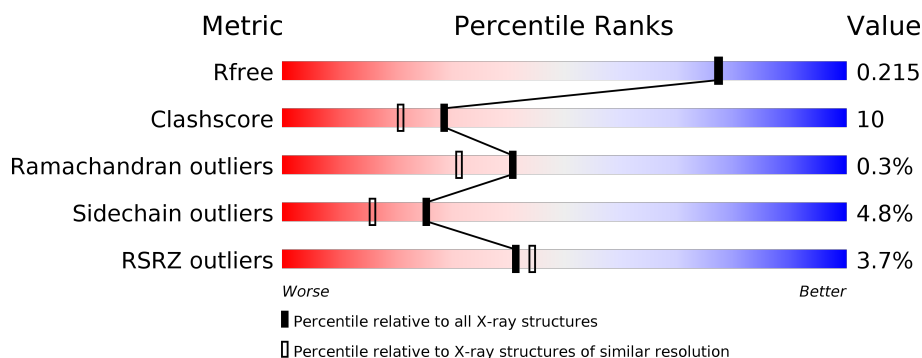
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	443	<div> <div>4%</div> <div> <div></div> <div>84%</div> <div>12%</div> <div>••</div> </div> </div>
1	B	443	<div> <div>4%</div> <div> <div></div> <div>84%</div> <div>12%</div> <div>••</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	PEG	B	1440	-	-	X	-
5	PEG	B	1442	-	-	X	-
5	PEG	B	1443	-	-	X	-

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7630 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 PHENYLACETATE-COENZYME A LIGASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	433	Total	C	N	O	S	0	2	0
			3339	2098	600	627	14			
1	B	432	Total	C	N	O	S	0	2	0
			3327	2091	596	626	14			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP B4EL89
A	-1	SER	-	expression tag	UNP B4EL89
A	0	HIS	-	expression tag	UNP B4EL89
B	-2	GLY	-	expression tag	UNP B4EL89
B	-1	SER	-	expression tag	UNP B4EL89
B	0	HIS	-	expression tag	UNP B4EL89

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

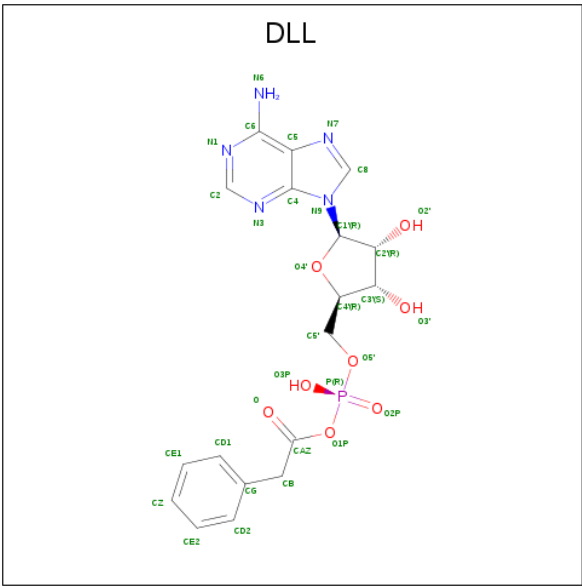
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	K	0	0
			1	1		
3	A	1	Total	K	0	0
			1	1		

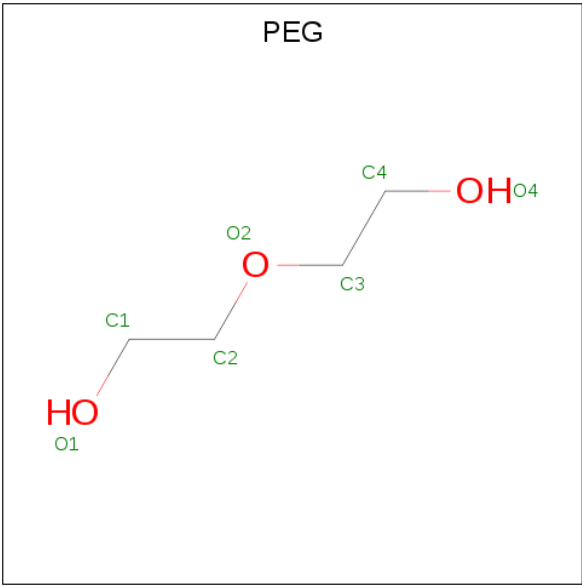
- Molecule 4 is 5'-O-[HYDROXY(PHENYLACETYL)PHOSPHORYL]ADENOSINE

(three-letter code: DLL) (formula: C<sub>18</sub>H<sub>20</sub>N<sub>5</sub>O<sub>8</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			32	18	5	8	1		
4	B	1	Total	C	N	O	P	0	0
			32	18	5	8	1		

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



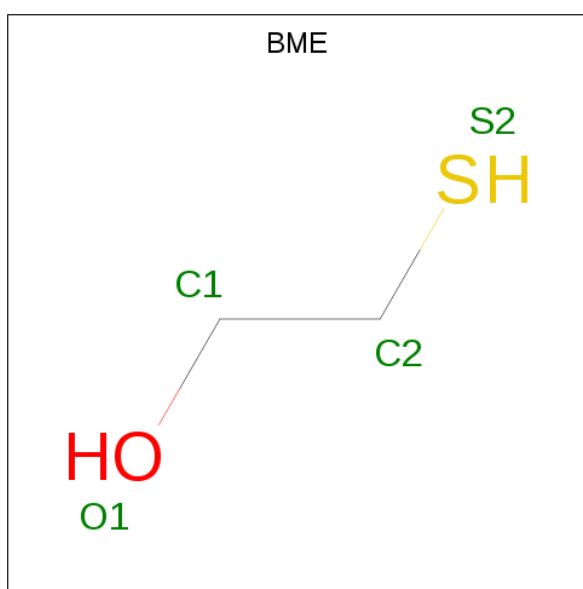
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 6 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C<sub>2</sub>H<sub>6</sub>OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	O	S	0	0
			4	2	1	1		

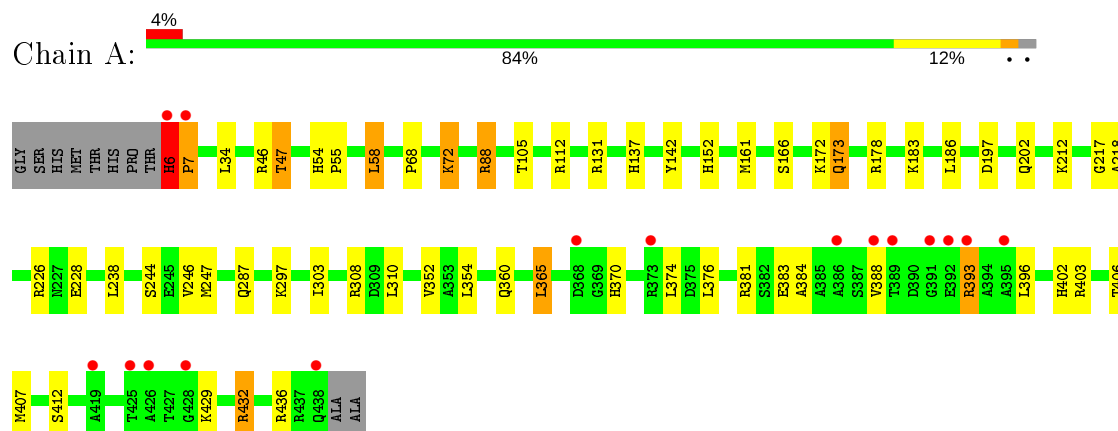
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	383	Total	O	0	0
			383	383		
7	B	477	Total	O	0	0
			477	477		

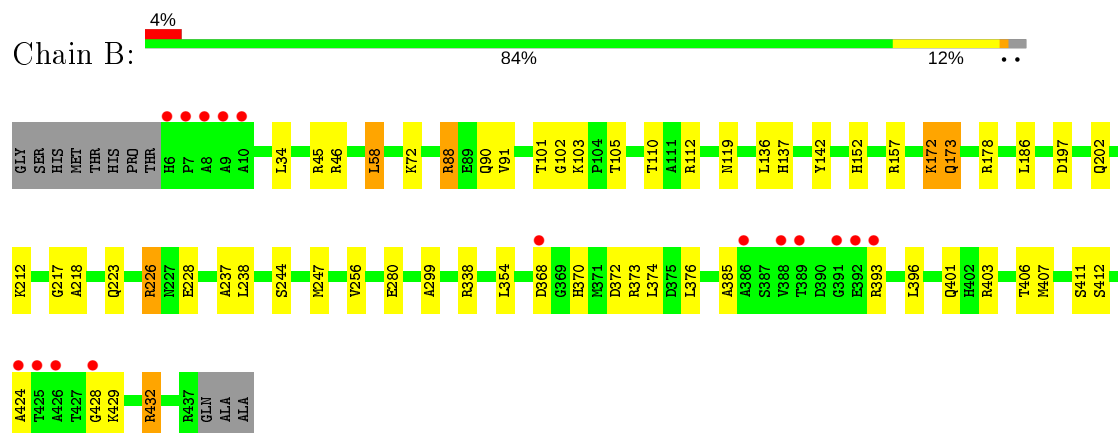
### 3 Residue-property plots [i](#)

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: PHENYLACETATE-COENZYME A LIGASE



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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.15Å 81.97Å 80.89Å 90.00° 97.27° 90.00°	Depositor
Resolution (Å)	40.99 – 1.90 40.98 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.9 (40.99-1.90) 98.9 (40.98-1.90)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.49 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.166 , 0.214 0.166 , 0.215	Depositor DCC
$R_{free}$ test set	3532 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	17.6	Xtriage
Anisotropy	0.094	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 53.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7630	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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: PEG, MG, DLL, K, BME

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	0.36	0/3410	0.49	1/4630 (0.0%)
1	B	0.37	0/3398	0.51	0/4614
All	All	0.36	0/6808	0.50	1/9244 (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	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	186	LEU	CA-CB-CG	-5.05	103.69	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	166	SER	Peptide
1	A	6	HIS	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	3339	0	3357	67	0
1	B	3327	0	3342	76	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	32	0	18	0	0
4	B	32	0	18	0	0
5	B	32	0	45	16	0
6	B	4	0	6	0	0
7	A	383	0	0	13	1
7	B	477	0	0	15	1
All	All	7630	0	6786	132	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 10.

All (132) 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:90:GLN:HG3	7:B:2168:HOH:O	1.42	1.17
1:A:406:THR:HG22	1:A:407:MET:CE	1.73	1.16
1:B:406:THR:HG22	1:B:407:MET:HE2	1.16	1.13
1:B:406:THR:HG22	1:B:407:MET:CE	1.80	1.10
1:A:393:ARG:CB	1:A:393:ARG:HH11	1.65	1.09
1:A:432:ARG:HH11	1:A:432:ARG:HG3	0.91	1.08
1:B:432:ARG:HH11	1:B:432:ARG:HG3	0.92	1.07
1:A:47:THR:HG21	7:A:2073:HOH:O	1.53	1.06
1:A:161:MET:HE2	1:B:91:VAL:HB	1.39	1.04
1:A:393:ARG:HB3	1:A:393:ARG:HH11	1.18	1.02
1:A:406:THR:HG22	1:A:407:MET:HE3	1.02	1.01
1:A:432:ARG:HH11	1:A:432:ARG:CG	1.72	1.00
1:A:432:ARG:NH1	1:A:432:ARG:HG3	1.67	0.98
1:A:161:MET:CE	1:B:91:VAL:HB	1.94	0.97
1:A:406:THR:CG2	1:A:407:MET:HE3	1.93	0.97
1:B:432:ARG:CG	1:B:432:ARG:HH11	1.76	0.96
1:B:432:ARG:NH1	1:B:432:ARG:HG3	1.67	0.95
1:B:172:LYS:CE	7:B:2262:HOH:O	2.19	0.90
1:A:212:LYS:HD3	7:A:2245:HOH:O	1.71	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:385:ALA:HB2	7:B:2454:HOH:O	1.72	0.88
1:B:45:ARG:HH21	5:B:1440:PEG:C4	1.91	0.84
1:B:101:THR:HG22	1:B:103:LYS:H	1.44	0.83
1:B:406:THR:CG2	1:B:407:MET:HE2	2.05	0.83
1:A:287:GLN:HE22	1:A:310:LEU:HB3	1.44	0.82
1:B:46:ARG:HE	5:B:1442:PEG:C2	1.94	0.81
1:B:101:THR:HG22	1:B:102:GLY:N	1.95	0.81
1:B:101:THR:CG2	1:B:102:GLY:N	2.44	0.80
1:A:393:ARG:HB3	1:A:393:ARG:NH1	1.97	0.80
1:B:223:GLN:HA	5:B:1443:PEG:H41	1.66	0.78
1:B:212:LYS:HE2	7:B:2306:HOH:O	1.85	0.77
1:B:368:ASP:HB2	1:B:373:ARG:HD2	1.66	0.77
1:A:393:ARG:HH11	1:A:393:ARG:CG	1.99	0.76
1:A:403:ARG:HD2	7:A:2357:HOH:O	1.85	0.75
1:B:244[B]:SER:OG	7:B:2334:HOH:O	2.05	0.74
1:B:172:LYS:HE2	7:B:2262:HOH:O	1.86	0.74
1:B:45:ARG:NH2	5:B:1440:PEG:O4	2.22	0.73
1:A:137:HIS:HE1	1:A:173:GLN:NE2	1.87	0.72
1:A:407:MET:HE2	1:B:172:LYS:HB2	1.70	0.72
1:B:101:THR:CG2	1:B:102:GLY:H	2.03	0.71
1:B:137:HIS:HE1	1:B:173:GLN:NE2	1.90	0.69
1:A:161:MET:HE3	1:B:88:ARG:O	1.91	0.69
1:A:161:MET:CE	1:B:88:ARG:O	2.41	0.69
1:A:197:ASP:OD2	1:A:370:HIS:HE1	1.77	0.67
1:B:223:GLN:HG2	5:B:1443:PEG:O4	1.95	0.67
1:A:6:HIS:N	1:A:7:PRO:CA	2.58	0.67
1:B:197:ASP:OD2	1:B:370:HIS:HE1	1.79	0.65
1:A:172:LYS:HB2	1:B:407:MET:CE	2.27	0.65
1:A:172:LYS:HB2	1:B:407:MET:HE1	1.78	0.64
1:B:112:ARG:HG3	7:B:2198:HOH:O	1.97	0.63
1:B:34:LEU:HD11	1:B:58:LEU:HD13	1.78	0.63
1:A:393:ARG:CB	1:A:393:ARG:NH1	2.50	0.63
1:A:244[B]:SER:OG	7:A:2269:HOH:O	2.15	0.63
1:B:45:ARG:NH2	5:B:1440:PEG:C4	2.61	0.63
1:A:407:MET:CE	1:B:172:LYS:HB2	2.29	0.62
1:B:101:THR:HG23	1:B:102:GLY:H	1.65	0.61
1:A:161:MET:HE1	1:B:91:VAL:HB	1.80	0.60
1:A:406:THR:CG2	1:A:407:MET:CE	2.65	0.59
1:B:101:THR:HG21	7:B:2183:HOH:O	2.01	0.59
1:B:142:TYR:OH	1:B:152:HIS:HD2	1.86	0.58
1:B:172:LYS:HB2	1:B:172:LYS:HZ3	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:TYR:OH	1:A:152:HIS:HD2	1.88	0.57
1:A:6:HIS:N	1:A:7:PRO:HA	2.20	0.56
1:A:365:LEU:HD22	1:A:365:LEU:N	2.21	0.56
1:A:183:LYS:HD2	7:A:2178:HOH:O	2.06	0.56
1:A:34:LEU:HD23	1:A:303:ILE:HD12	1.87	0.55
1:B:137:HIS:HD2	7:B:2263:HOH:O	1.88	0.55
1:B:46:ARG:HB2	5:B:1442:PEG:H21	1.88	0.54
1:A:402:HIS:HD2	7:A:2362:HOH:O	1.90	0.54
1:A:46:ARG:NH1	7:A:2070:HOH:O	2.39	0.53
1:A:112:ARG:NH2	7:A:2155:HOH:O	2.40	0.53
1:B:226:ARG:CZ	5:B:1443:PEG:H31	2.38	0.53
1:A:47:THR:HG23	1:A:68:PRO:HG3	1.90	0.53
1:B:403:ARG:HG2	1:B:403:ARG:HH11	1.74	0.52
1:A:393:ARG:NH1	1:A:393:ARG:CG	2.65	0.51
1:B:46:ARG:HH21	5:B:1442:PEG:H42	1.75	0.51
1:B:338:ARG:NH2	1:B:372:ASP:O	2.42	0.51
1:B:197:ASP:OD2	1:B:370:HIS:CE1	2.63	0.51
1:B:172:LYS:CD	7:B:2262:HOH:O	2.56	0.50
1:B:403:ARG:HG2	1:B:403:ARG:NH1	2.27	0.50
1:A:197:ASP:OD2	1:A:370:HIS:CE1	2.62	0.50
1:A:34:LEU:HD11	1:A:58:LEU:HD13	1.94	0.50
1:A:217:GLY:O	1:A:218:ALA:HB3	2.12	0.49
1:B:110:THR:HG23	1:B:299:ALA:HB2	1.94	0.49
1:B:46:ARG:CB	5:B:1442:PEG:H21	2.42	0.49
1:A:54:HIS:CG	1:A:55:PRO:HD2	2.48	0.48
1:B:46:ARG:HH21	5:B:1442:PEG:C3	2.26	0.48
1:A:161:MET:HE2	1:B:91:VAL:CB	2.27	0.48
1:A:287:GLN:NE2	1:A:310:LEU:HB3	2.20	0.48
1:A:6:HIS:N	1:A:7:PRO:C	2.67	0.48
1:A:178:ARG:HH11	1:A:202:GLN:HE21	1.61	0.48
1:B:45:ARG:HD3	7:B:2089:HOH:O	2.13	0.48
1:A:88:ARG:HD3	7:B:2227:HOH:O	2.14	0.48
5:B:1440:PEG:O1	5:B:1440:PEG:H32	2.14	0.48
1:A:137:HIS:HD2	7:A:2210:HOH:O	1.96	0.48
1:B:119[B]:ASN:OD1	1:B:157:ARG:NH1	2.46	0.48
1:B:46:ARG:HH21	5:B:1442:PEG:C4	2.27	0.48
1:A:352:VAL:HG23	7:A:2336:HOH:O	2.14	0.47
1:A:381:ARG:HD3	1:A:383:GLU:OE2	2.14	0.47
1:B:45:ARG:HH21	5:B:1440:PEG:H41	1.74	0.47
1:A:142:TYR:HB3	1:B:142:TYR:HB3	1.96	0.47
1:B:34:LEU:CD1	1:B:58:LEU:HD13	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:LYS:HE3	7:B:2262:HOH:O	2.02	0.46
1:B:178:ARG:HH11	1:B:202:GLN:HE21	1.63	0.46
1:A:246:VAL:O	1:A:297:LYS:HD2	2.15	0.46
1:B:217:GLY:O	1:B:218:ALA:HB3	2.17	0.45
1:B:407:MET:HA	1:B:407:MET:CE	2.46	0.45
1:B:407:MET:HE1	1:B:407:MET:HA	1.98	0.45
1:A:403:ARG:HD3	7:A:2364:HOH:O	2.17	0.44
1:A:47:THR:HG23	1:A:68:PRO:CG	2.47	0.44
1:B:137:HIS:HE1	1:B:173:GLN:HE22	1.65	0.44
1:B:424:ALA:HB1	1:B:428:GLY:HA2	2.00	0.44
1:A:142:TYR:OH	1:A:152:HIS:CD2	2.69	0.44
1:B:72:LYS:HE3	1:B:105:THR:OG1	2.17	0.43
1:A:374:LEU:HD22	1:A:412:SER:OG	2.19	0.43
1:A:88:ARG:CD	7:B:2227:HOH:O	2.67	0.43
1:B:46:ARG:NE	5:B:1442:PEG:C2	2.72	0.43
1:A:432:ARG:NH1	1:A:432:ARG:CG	2.43	0.43
5:B:1443:PEG:H42	5:B:1443:PEG:H22	1.49	0.42
1:A:172:LYS:HB2	1:B:407:MET:HE3	2.01	0.42
1:B:401:GLN:HG3	1:B:412:SER:O	2.19	0.41
1:A:137:HIS:HE1	1:A:173:GLN:HE22	1.67	0.41
1:A:365:LEU:CD2	1:A:365:LEU:N	2.84	0.41
1:B:406:THR:HG22	1:B:407:MET:HE3	1.88	0.41
1:A:384:ALA:O	1:A:388:VAL:HG23	2.20	0.41
1:B:101:THR:HB	7:B:2188:HOH:O	2.20	0.41
1:B:238:LEU:HD21	1:B:256:VAL:HA	2.03	0.41
1:A:407:MET:HE1	1:B:172:LYS:N	2.36	0.41
1:B:136:LEU:HD11	1:B:186:LEU:HD13	2.02	0.41
1:A:72:LYS:HE3	1:A:105:THR:OG1	2.21	0.41
1:A:436:ARG:NE	7:A:2378:HOH:O	2.27	0.41
1:A:436:ARG:NH2	7:A:2377:HOH:O	2.54	0.40
1:B:237:ALA:C	1:B:238:LEU:HG	2.42	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)
7:A:2326:HOH:O	7:B:2293:HOH:O[2_757]	2.17	0.03

## 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	433/443 (98%)	419 (97%)	12 (3%)	2 (0%)	29	18
1	B	432/443 (98%)	421 (98%)	10 (2%)	1 (0%)	47	38
All	All	865/886 (98%)	840 (97%)	22 (2%)	3 (0%)	41	31

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	247	MET
1	B	247	MET
1	A	7	PRO

### 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	355/360 (99%)	335 (94%)	20 (6%)	21	11
1	B	354/360 (98%)	339 (96%)	15 (4%)	30	20
All	All	709/720 (98%)	674 (95%)	35 (5%)	25	15

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	HIS
1	A	47	THR
1	A	58	LEU

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Mol	Chain	Res	Type
1	A	72	LYS
1	A	88	ARG
1	A	131[A]	ARG
1	A	131[B]	ARG
1	A	173	GLN
1	A	226	ARG
1	A	228	GLU
1	A	238	LEU
1	A	308	ARG
1	A	354	LEU
1	A	360	GLN
1	A	365	LEU
1	A	376	LEU
1	A	393	ARG
1	A	396	LEU
1	A	429	LYS
1	A	432	ARG
1	B	58	LEU
1	B	88	ARG
1	B	172	LYS
1	B	173	GLN
1	B	226	ARG
1	B	228	GLU
1	B	280	GLU
1	B	354	LEU
1	B	374	LEU
1	B	376	LEU
1	B	393	ARG
1	B	396	LEU
1	B	411	SER
1	B	429	LYS
1	B	432	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	137	HIS
1	A	152	HIS
1	A	173	GLN
1	A	202	GLN
1	A	227	ASN

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Mol	Chain	Res	Type
1	A	287	GLN
1	A	360	GLN
1	A	370	HIS
1	B	137	HIS
1	B	152	HIS
1	B	173	GLN
1	B	202	GLN
1	B	227	ASN
1	B	370	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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	$\# Z  > 2$
4	DLL	B	1438	3	31,35,35	1.65	4 (12%)	35,51,51	1.26	3 (8%)
5	PEG	B	1439	-	3,3,6	0.54	0	2,2,5	0.27	0
4	DLL	A	1441	3	31,35,35	1.63	4 (12%)	35,51,51	1.45	4 (11%)
5	PEG	B	1441	-	6,6,6	0.46	0	5,5,5	0.55	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	PEG	B	1440	-	6,6,6	0.49	0	5,5,5	0.52	0
5	PEG	B	1442	-	6,6,6	0.51	0	5,5,5	0.56	0
5	PEG	B	1443	-	6,6,6	0.53	0	5,5,5	0.46	0
6	BME	B	1444	-	3,3,3	0.39	0	1,2,2	0.37	0

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
4	DLL	B	1438	3	-	1/13/35/35	0/4/4/4
5	PEG	B	1439	-	-	0/1/1/4	-
4	DLL	A	1441	3	-	1/13/35/35	0/4/4/4
5	PEG	B	1441	-	-	2/4/4/4	-
5	PEG	B	1440	-	-	2/4/4/4	-
5	PEG	B	1442	-	-	1/4/4/4	-
5	PEG	B	1443	-	-	1/4/4/4	-
6	BME	B	1444	-	-	0/1/1/1	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1441	DLL	C8-N7	6.80	1.46	1.34
4	B	1438	DLL	C8-N7	6.44	1.46	1.34
4	B	1438	DLL	O4'-C4'	-3.37	1.37	1.45
4	A	1441	DLL	O4'-C4'	-3.26	1.37	1.45
4	B	1438	DLL	O4'-C1'	3.08	1.45	1.41
4	A	1441	DLL	O4'-C1'	2.34	1.44	1.41
4	B	1438	DLL	O5'-C5'	-2.24	1.36	1.44
4	A	1441	DLL	O5'-C5'	-2.11	1.36	1.44

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1441	DLL	N3-C2-N1	-5.10	120.70	128.68
4	B	1438	DLL	N3-C2-N1	-4.73	121.28	128.68
4	A	1441	DLL	O4'-C4'-C3'	3.26	111.56	105.11
4	A	1441	DLL	O4'-C4'-C5'	2.98	119.16	109.37
4	B	1438	DLL	O4'-C4'-C5'	2.83	118.69	109.37
4	A	1441	DLL	C5'-C4'-C3'	2.34	123.94	115.18
4	B	1438	DLL	O4'-C4'-C3'	2.15	109.36	105.11

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1441	DLL	O4'-C4'-C5'-O5'
5	B	1443	PEG	C4-C3-O2-C2
5	B	1442	PEG	O2-C3-C4-O4
4	B	1438	DLL	O4'-C4'-C5'-O5'
5	B	1441	PEG	O1-C1-C2-O2
5	B	1440	PEG	O2-C3-C4-O4
5	B	1440	PEG	C4-C3-O2-C2
5	B	1441	PEG	C4-C3-O2-C2

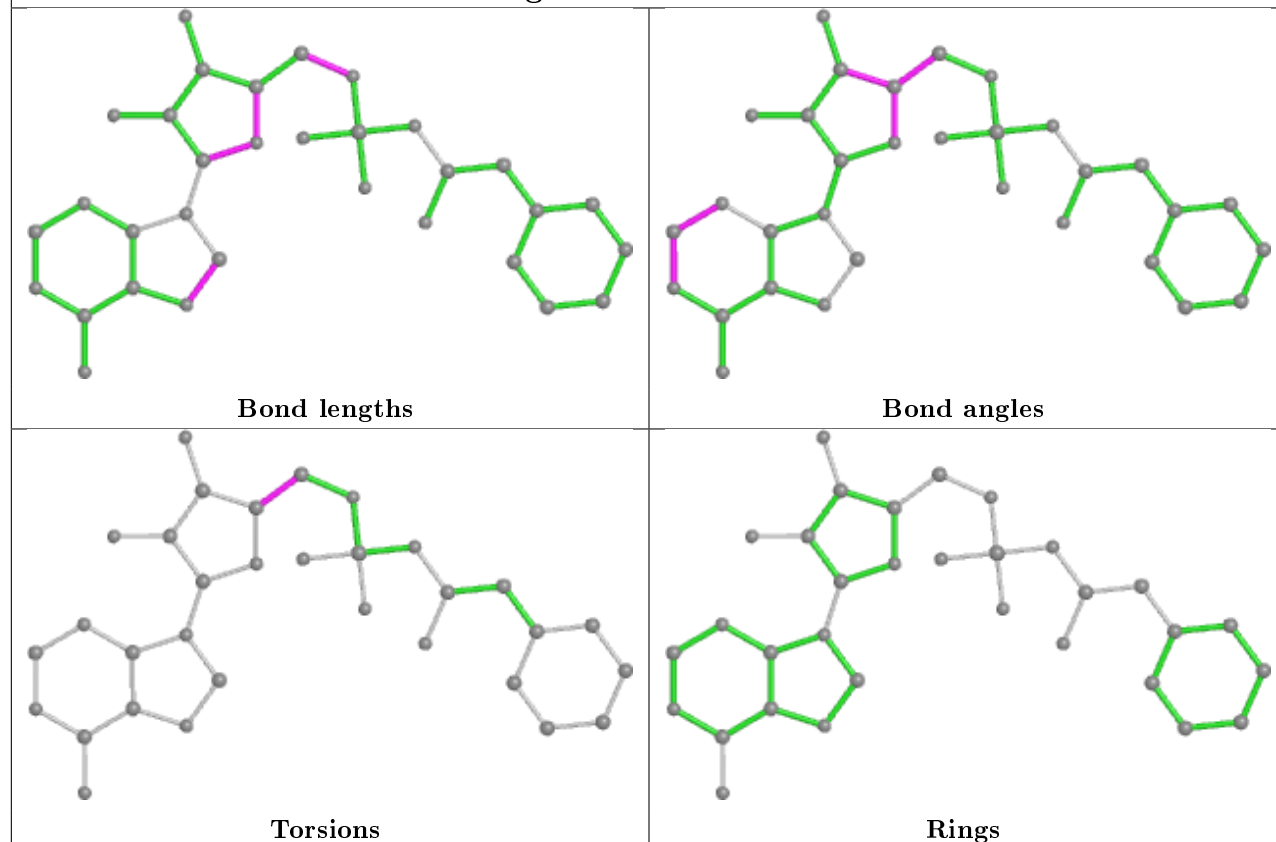
There are no ring outliers.

3 monomers are involved in 16 short contacts:

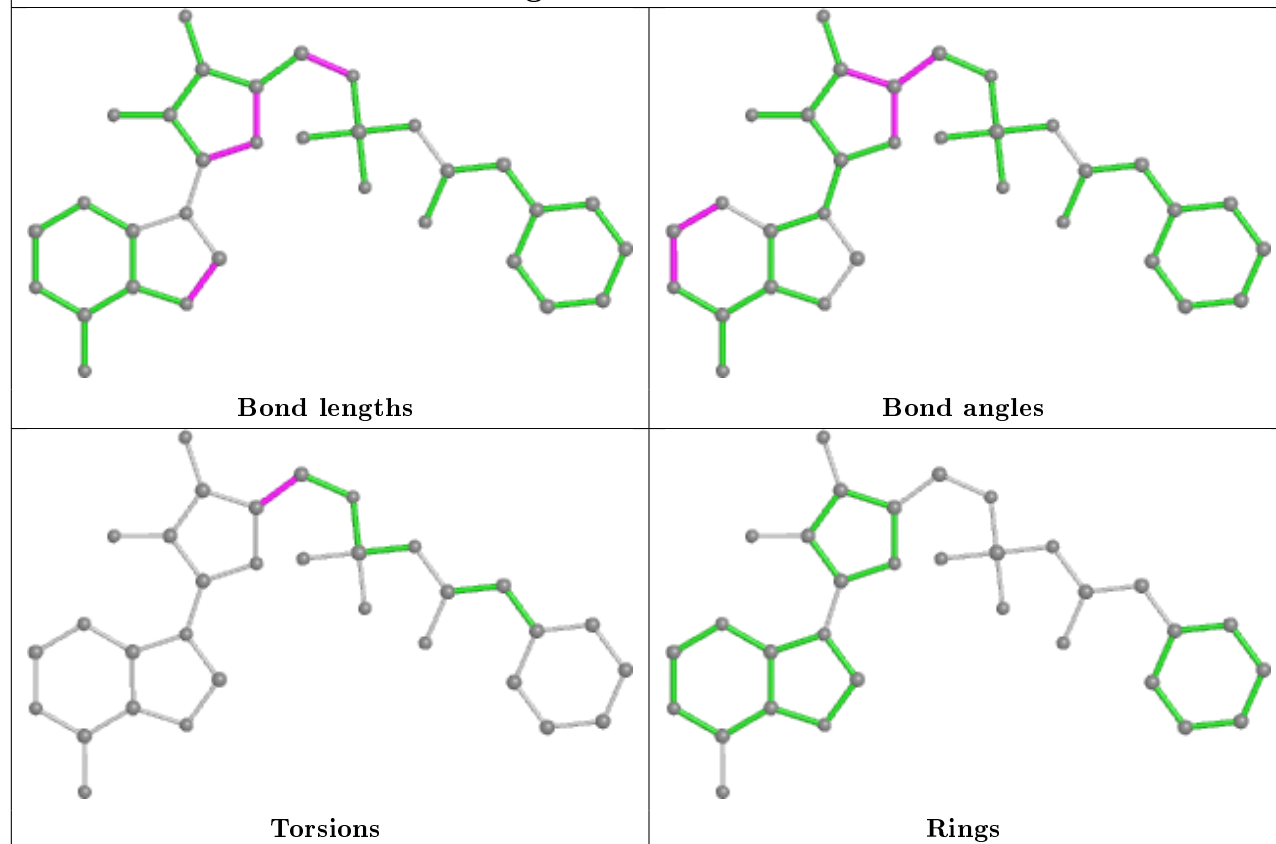
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1440	PEG	5	0
5	B	1442	PEG	7	0
5	B	1443	PEG	4	0

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.

## Ligand DLL B 1438



## Ligand DLL A 1441



## 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	433/443 (97%)	-0.02	16 (3%) 41 44	7, 17, 39, 51	0
1	B	432/443 (97%)	-0.01	16 (3%) 41 44	7, 16, 40, 50	0
All	All	865/886 (97%)	-0.02	32 (3%) 41 44	7, 17, 39, 51	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	426	ALA	5.2
1	B	7	PRO	4.8
1	A	426	ALA	4.5
1	B	428	GLY	4.3
1	A	7	PRO	3.7
1	B	391	GLY	3.7
1	A	6	HIS	3.4
1	B	9	ALA	3.4
1	B	386	ALA	3.3
1	B	6	HIS	3.2
1	A	425	THR	2.9
1	A	428	GLY	2.9
1	B	8	ALA	2.9
1	B	393	ARG	2.7
1	B	425	THR	2.7
1	B	392	GLU	2.7
1	A	392	GLU	2.6
1	A	368	ASP	2.6
1	A	393	ARG	2.5
1	B	389	THR	2.5
1	B	388	VAL	2.5
1	B	10	ALA	2.3
1	A	391	GLY	2.3
1	A	386	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	388	VAL	2.2
1	B	368	ASP	2.1
1	A	395	ALA	2.1
1	A	419	ALA	2.1
1	B	424	ALA	2.1
1	A	389	THR	2.1
1	A	438	GLN	2.1
1	A	373	ARG	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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.

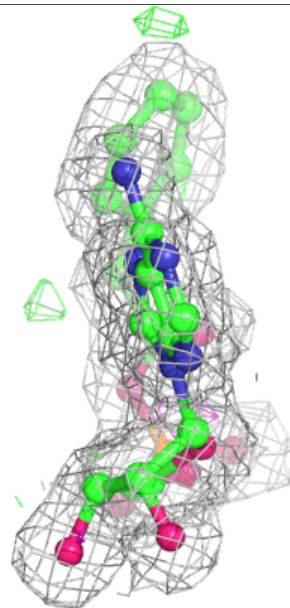
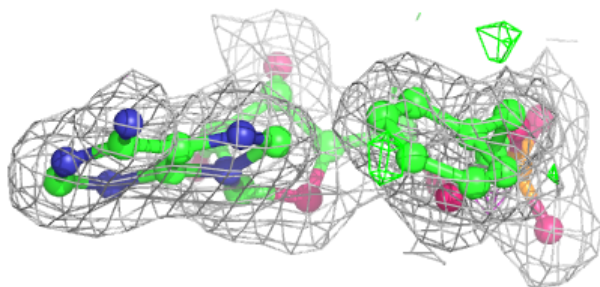
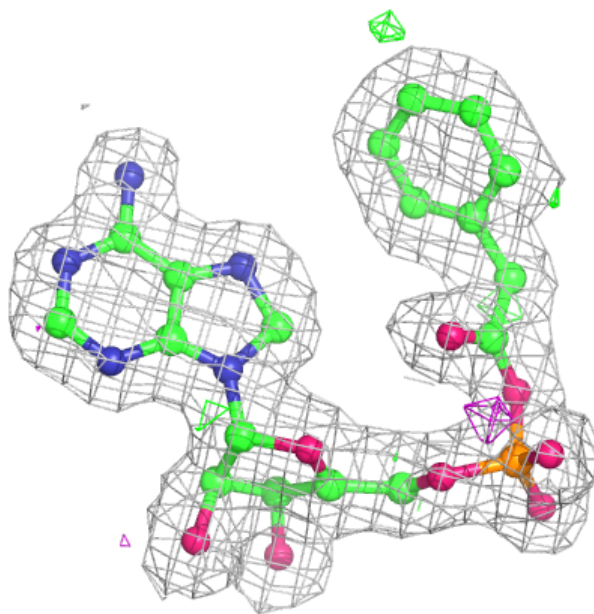
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	PEG	B	1439	4/7	0.79	0.26	38,40,43,44	0
5	PEG	B	1442	7/7	0.79	0.28	42,43,43,46	0
6	BME	B	1444	4/4	0.83	0.15	32,41,45,58	0
5	PEG	B	1440	7/7	0.87	0.20	52,54,56,58	0
5	PEG	B	1441	7/7	0.90	0.16	50,52,55,56	0
5	PEG	B	1443	7/7	0.91	0.13	39,40,48,51	0
4	DLL	A	1441	32/32	0.97	0.11	8,12,16,17	0
4	DLL	B	1438	32/32	0.97	0.12	8,11,14,20	0
2	MG	B	1445	1/1	0.99	0.09	24,24,24,24	0
3	K	B	1446	1/1	0.99	0.12	12,12,12,12	0
2	MG	A	1439	1/1	0.99	0.06	19,19,19,19	0
3	K	A	1440	1/1	1.00	0.09	12,12,12,12	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.

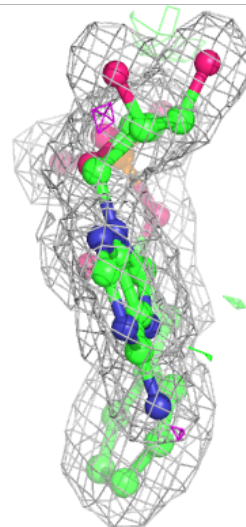
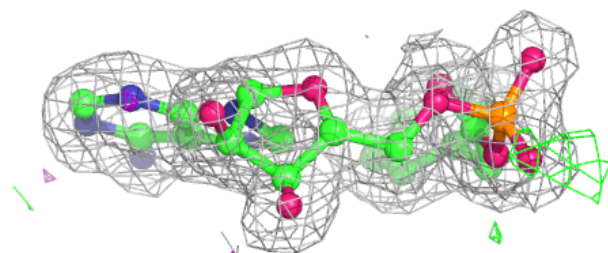
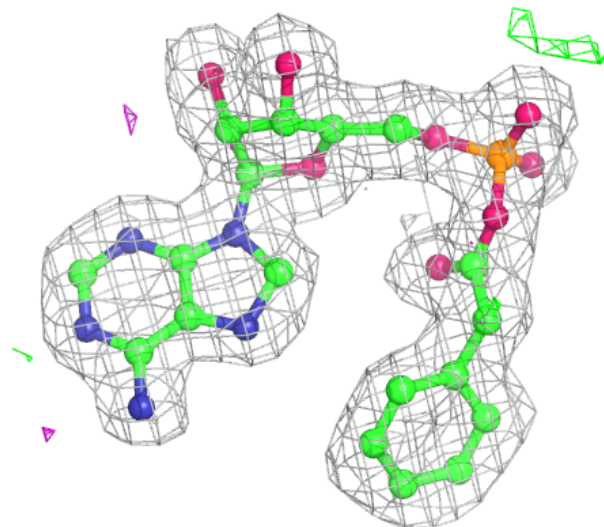
**Electron density around DLL A 1441:**

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



**Electron density around DLL B 1438:**

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



## 6.5 Other polymers ⓘ

There are no such residues in this entry.