



# Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 07:03 am BST

PDB ID : 4TVW  
Title : Resorufin ligase with bound resorufin-AMP analog  
Authors : Goldman, P.J.; Drennan, C.L.  
Deposited on : 2014-06-28  
Resolution : 3.50 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

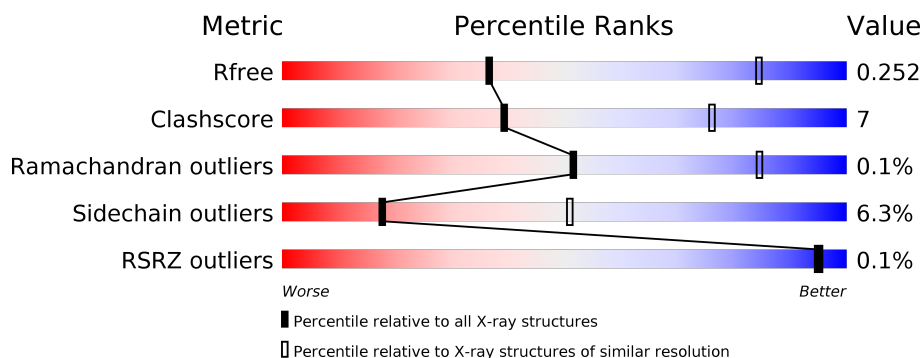
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (3.60-3.40)

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	341	<div> <div style="width: 81%;"></div> <div style="width: 16%;"></div> <div style="width: 3%;"></div> <div style="width: 2%;"></div> </div> <div>81% 16% ..</div>
1	B	341	<div> <div style="width: 81%;"></div> <div style="width: 17%;"></div> <div style="width: 3%;"></div> <div style="width: 2%;"></div> </div> <div>81% 17% ..</div>
1	C	341	<div> <div style="width: 76%;"></div> <div style="width: 21%;"></div> <div style="width: 3%;"></div> <div style="width: 2%;"></div> </div> <div>76% 21% ..</div>
1	D	341	<div> <div style="width: 75%;"></div> <div style="width: 21%;"></div> <div style="width: 3%;"></div> <div style="width: 2%;"></div> </div> <div>75% 21% ..</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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	37P	A	401	X	-	-	-
2	37P	B	401	X	-	-	-
2	37P	D	401	X	-	-	-

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10768 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 Lipoate-protein ligase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2647	1661	474	501	11			
1	B	337	Total	C	N	O	S	0	0	0
			2647	1661	474	501	11			
1	C	337	Total	C	N	O	S	0	0	0
			2647	1661	474	501	11			
1	D	337	Total	C	N	O	S	0	0	0
			2647	1661	474	501	11			

There are 24 discrepancies between the modelled and reference sequences:

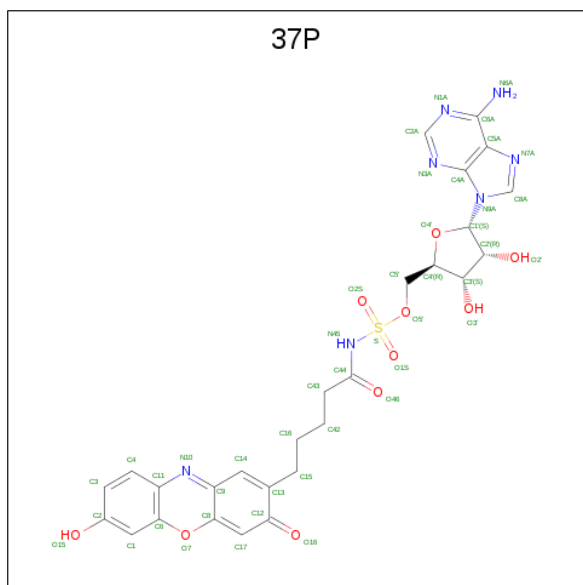
Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P32099
A	-2	SER	-	expression tag	UNP P32099
A	-1	HIS	-	expression tag	UNP P32099
A	20	ALA	GLU	engineered mutation	UNP P32099
A	147	ALA	PHE	engineered mutation	UNP P32099
A	149	GLY	HIS	engineered mutation	UNP P32099
B	-3	GLY	-	expression tag	UNP P32099
B	-2	SER	-	expression tag	UNP P32099
B	-1	HIS	-	expression tag	UNP P32099
B	20	ALA	GLU	engineered mutation	UNP P32099
B	147	ALA	PHE	engineered mutation	UNP P32099
B	149	GLY	HIS	engineered mutation	UNP P32099
C	-3	GLY	-	expression tag	UNP P32099
C	-2	SER	-	expression tag	UNP P32099
C	-1	HIS	-	expression tag	UNP P32099
C	20	ALA	GLU	engineered mutation	UNP P32099
C	147	ALA	PHE	engineered mutation	UNP P32099
C	149	GLY	HIS	engineered mutation	UNP P32099
D	-3	GLY	-	expression tag	UNP P32099
D	-2	SER	-	expression tag	UNP P32099
D	-1	HIS	-	expression tag	UNP P32099

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Chain	Residue	Modelled	Actual	Comment	Reference
D	20	ALA	GLU	engineered mutation	UNP P32099
D	147	ALA	PHE	engineered mutation	UNP P32099
D	149	GLY	HIS	engineered mutation	UNP P32099

- Molecule 2 is 5'-O-[[5-(7-hydroxy-3-oxo-3H-phenoxazin-2-yl)pentanoyl]sulfamoyl]adenosine (three-letter code: 37P) (formula: C<sub>27</sub>H<sub>27</sub>N<sub>7</sub>O<sub>10</sub>S).

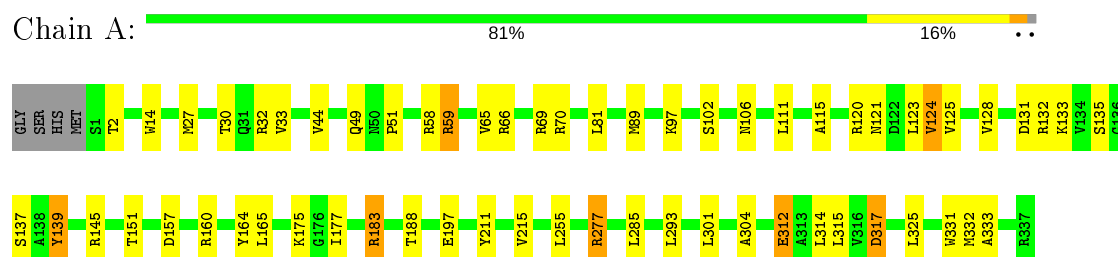


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 45	C 27	N 7	O 10	S 1	0	0
2	B	1	Total 45	C 27	N 7	O 10	S 1	0	0
2	C	1	Total 45	C 27	N 7	O 10	S 1	0	0
2	D	1	Total 45	C 27	N 7	O 10	S 1	0	0

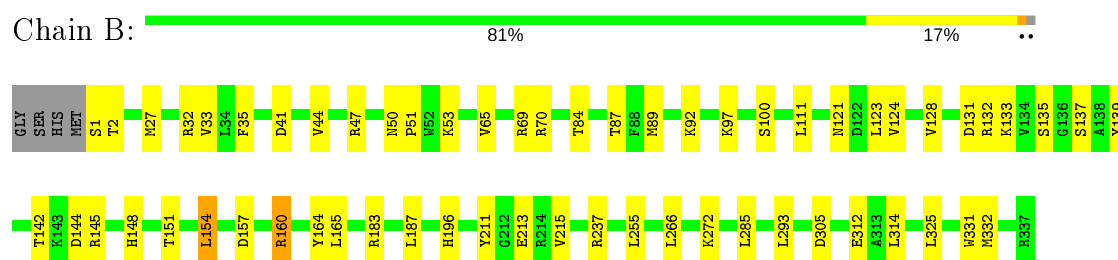
### 3 Residue-property plots

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.

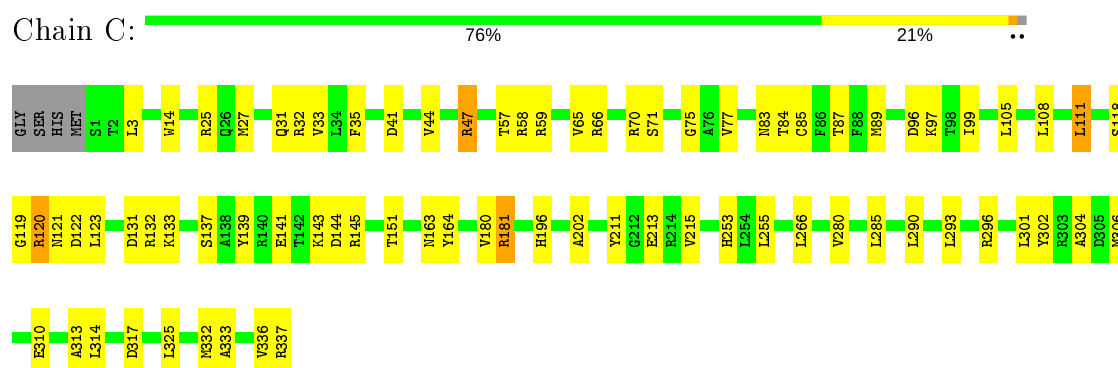
#### • Molecule 1: Lipoate-protein ligase A



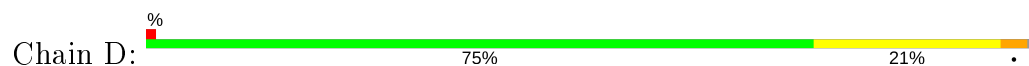
#### • Molecule 1: Lipoate-protein ligase A

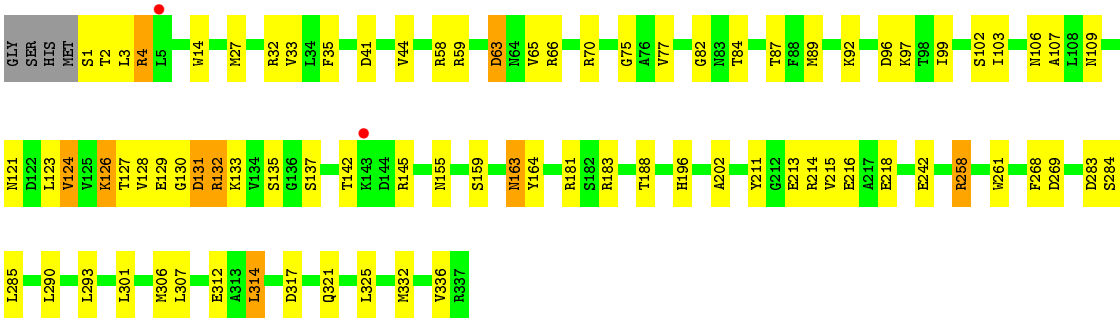


#### • Molecule 1: Lipoate-protein ligase A



#### • Molecule 1: Lipoate-protein ligase A





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.35Å 99.93Å 107.28Å 90.00° 105.79° 90.00°	Depositor
Resolution (Å)	49.97 – 3.50 49.97 – 3.51	Depositor EDS
% Data completeness (in resolution range)	87.1 (49.97-3.50) 87.2 (49.97-3.51)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.14	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.64 (at 3.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.201 , 0.248 0.206 , 0.252	Depositor DCC
$R_{free}$ test set	1049 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	54.0	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 28.6	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	0.90	EDS
Total number of atoms	10768	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% 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: 37P

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	0.22	0/2702	0.43	0/3661
1	B	0.23	0/2702	0.44	1/3661 (0.0%)
1	C	0.24	0/2702	0.48	1/3661 (0.0%)
1	D	0.22	0/2702	0.46	0/3661
All	All	0.23	0/10808	0.45	2/14644 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	141	GLU	CB-CA-C	5.21	120.83	110.40
1	B	142	THR	N-CA-C	-5.04	97.41	111.00

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	2647	0	2596	32	0
1	B	2647	0	2596	32	0
1	C	2647	0	2596	47	0
1	D	2647	0	2596	50	0
2	A	45	0	26	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	45	0	26	4	0
2	C	45	0	26	9	0
2	D	45	0	26	2	0
All	All	10768	0	10488	158	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 7.

All (158) 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:C:120:ARG:HH21	1:C:180:VAL:HG21	1.16	1.08
1:C:120:ARG:NH2	1:C:180:VAL:HG21	1.81	0.94
1:B:41:ASP:O	1:B:65:VAL:HG13	1.69	0.92
1:C:120:ARG:HH11	1:C:120:ARG:HG3	1.36	0.90
1:D:124:VAL:HG21	1:D:131:ASP:OD2	1.91	0.70
1:C:33:VAL:HB	1:C:89:MET:HB2	1.73	0.70
1:D:44:VAL:HG12	1:D:70:ARG:HH11	1.56	0.69
1:C:122:ASP:OD1	1:C:133:LYS:NZ	2.21	0.69
1:D:126:LYS:HE3	1:D:131:ASP:OD1	1.93	0.69
1:D:84:THR:HG1	1:D:196:HIS:HD1	1.39	0.69
1:C:35:PHE:HB3	1:C:87:THR:HB	1.75	0.68
1:C:120:ARG:NH1	1:C:120:ARG:HG3	1.98	0.68
1:C:151:THR:OG1	2:C:401:37P:N1A	2.28	0.65
1:C:120:ARG:NH2	1:C:180:VAL:CG2	2.59	0.65
1:A:123:LEU:HB2	1:A:135:SER:HB3	1.79	0.64
1:A:115:ALA:HB2	1:A:125:VAL:HG12	1.81	0.63
1:D:126:LYS:CE	1:D:131:ASP:OD1	2.47	0.63
1:B:44:VAL:HG12	1:B:70:ARG:HH11	1.63	0.63
1:C:120:ARG:HH11	1:C:120:ARG:CG	2.08	0.63
1:C:27:MET:O	1:C:145:ARG:NH1	2.25	0.63
1:C:44:VAL:HG12	1:C:70:ARG:HH11	1.63	0.63
1:A:277:ARG:NH2	1:D:269:ASP:OD1	2.29	0.63
1:C:118:SER:OG	1:C:119:GLY:N	2.32	0.62
1:D:41:ASP:O	1:D:65:VAL:HG13	1.99	0.62
1:B:133:LYS:NZ	2:B:401:37P:H2	2.15	0.61
1:C:133:LYS:HZ3	2:C:401:37P:H1	1.65	0.61
1:A:27:MET:O	1:A:145:ARG:NH1	2.33	0.61
1:B:157:ASP:OD2	1:B:160:ARG:NH1	2.34	0.60
1:D:132:ARG:HH12	1:D:183:ARG:HH11	1.48	0.60
1:B:35:PHE:HB3	1:B:87:THR:HB	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:ASP:O	1:C:65:VAL:HG13	2.01	0.60
1:C:293:LEU:HD21	1:C:332:MET:HE1	1.83	0.60
1:C:181:ARG:HA	2:C:401:37P:O3'	2.02	0.60
1:D:27:MET:O	1:D:145:ARG:NH1	2.29	0.59
1:A:44:VAL:HG12	1:A:70:ARG:HH11	1.67	0.59
1:D:92:LYS:NZ	1:D:142:THR:O	2.27	0.59
1:B:84:THR:OG1	1:B:196:HIS:ND1	2.36	0.58
1:A:157:ASP:OD2	1:A:160:ARG:NH1	2.37	0.58
1:B:92:LYS:N	1:B:144:ASP:O	2.24	0.57
1:C:65:VAL:HG21	1:C:164:TYR:CE2	2.40	0.57
1:B:123:LEU:HB2	1:B:135:SER:HB3	1.86	0.57
1:D:102:SER:O	1:D:106:ASN:ND2	2.30	0.57
1:B:33:VAL:HB	1:B:89:MET:HB2	1.86	0.56
1:C:105:LEU:HA	1:C:108:LEU:HD12	1.87	0.56
1:D:133:LYS:HE3	1:D:135:SER:O	2.05	0.56
1:D:4:ARG:NH2	1:D:218:GLU:OE1	2.37	0.56
1:B:124:VAL:HG21	1:B:131:ASP:HB3	1.88	0.56
1:D:65:VAL:HG21	1:D:164:TYR:CE2	2.41	0.56
1:B:84:THR:HG1	1:B:196:HIS:HD1	1.53	0.55
1:A:58:ARG:HH22	1:A:59:ARG:HH21	1.53	0.54
1:C:121:ASN:HB2	1:C:137:SER:OG	2.07	0.54
1:C:296:ARG:NH2	1:C:317:ASP:OD2	2.37	0.54
1:D:96:ASP:HB3	1:D:99:ILE:HG13	1.89	0.54
1:D:124:VAL:CG2	1:D:131:ASP:OD2	2.56	0.54
1:C:121:ASN:ND2	1:C:137:SER:O	2.41	0.53
1:A:133:LYS:NZ	2:A:401:37P:H2	2.24	0.53
1:C:293:LEU:HB2	1:C:314:LEU:HD11	1.90	0.53
1:D:133:LYS:CE	1:D:135:SER:O	2.57	0.53
1:A:33:VAL:HB	1:A:89:MET:HB2	1.90	0.52
1:C:301:LEU:HB2	1:C:306:MET:HG3	1.91	0.52
1:D:84:THR:OG1	1:D:196:HIS:ND1	2.32	0.52
1:A:293:LEU:HD21	1:A:332:MET:HE1	1.91	0.52
1:D:293:LEU:HD21	1:D:332:MET:HE1	1.92	0.52
1:B:133:LYS:HZ1	2:B:401:37P:H2	1.74	0.52
1:D:132:ARG:NH1	1:D:183:ARG:HD2	2.25	0.51
1:B:121:ASN:HB2	1:B:137:SER:OG	2.09	0.51
1:B:132:ARG:NH1	1:B:183:ARG:HH11	2.08	0.51
1:A:151:THR:HG23	2:A:401:37P:H21	1.93	0.51
1:B:27:MET:O	1:B:145:ARG:NH1	2.29	0.51
1:D:70:ARG:NH2	1:D:75:GLY:O	2.37	0.51
1:D:301:LEU:HB2	1:D:306:MET:HG3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:TRP:CD2	1:C:66:ARG:HD3	2.46	0.50
1:B:160:ARG:HG2	1:B:164:TYR:CE2	2.46	0.50
1:A:131:ASP:O	1:A:132:ARG:NH1	2.40	0.50
1:B:65:VAL:HG21	1:B:164:TYR:CE2	2.46	0.49
1:B:255:LEU:HB2	1:B:266:LEU:HB2	1.94	0.49
1:A:165:LEU:CD1	2:A:401:37P:H12	2.42	0.49
1:D:3:LEU:HB2	1:D:215:VAL:HG21	1.94	0.49
1:D:307:LEU:HD12	1:D:336:VAL:HG21	1.94	0.49
1:A:32:ARG:HG2	1:A:211:TYR:HE1	1.77	0.49
1:B:293:LEU:HD21	1:B:332:MET:HE1	1.93	0.49
1:D:121:ASN:HB2	1:D:137:SER:OG	2.12	0.49
1:A:124:VAL:HG21	1:A:131:ASP:HB3	1.93	0.48
1:B:154:LEU:HD12	1:B:187:LEU:HD12	1.95	0.48
1:C:180:VAL:O	2:C:401:37P:H2	2.13	0.48
1:A:121:ASN:HB3	1:A:139:TYR:OH	2.12	0.48
1:C:296:ARG:NH1	1:C:313:ALA:O	2.46	0.48
1:D:32:ARG:HG2	1:D:211:TYR:HE1	1.78	0.48
1:A:102:SER:O	1:A:106:ASN:ND2	2.30	0.48
1:D:35:PHE:HB3	1:D:87:THR:HB	1.94	0.48
1:C:133:LYS:NZ	2:C:401:37P:H1	2.28	0.48
1:D:77:VAL:HG11	2:D:401:37P:H15	1.95	0.47
1:C:3:LEU:HD22	1:C:213:GLU:HB2	1.97	0.47
1:C:253:HIS:ND1	1:C:302:TYR:OH	2.38	0.47
1:C:32:ARG:HG2	1:C:211:TYR:HE1	1.80	0.47
1:B:255:LEU:HB3	1:B:331:TRP:CZ2	2.50	0.47
1:B:51:PRO:HD3	1:B:69:ARG:HH22	1.80	0.47
1:D:127:THR:N	1:D:130:GLY:O	2.46	0.47
1:A:255:LEU:HB3	1:A:331:TRP:CZ2	2.50	0.46
1:D:293:LEU:HB2	1:D:314:LEU:HD11	1.96	0.46
1:D:33:VAL:HB	1:D:89:MET:HB2	1.97	0.46
1:D:261:TRP:HZ3	1:D:290:LEU:HD11	1.81	0.46
1:B:121:ASN:ND2	1:B:137:SER:O	2.48	0.46
1:B:132:ARG:HH11	1:B:183:ARG:HH11	1.63	0.46
1:C:77:VAL:HG12	2:C:401:37P:H8	1.97	0.46
1:D:123:LEU:HB2	1:D:135:SER:HB3	1.98	0.46
1:A:51:PRO:HD3	1:A:69:ARG:HH22	1.81	0.46
1:B:132:ARG:HH11	1:B:183:ARG:HD2	1.80	0.46
1:C:58:ARG:HH22	1:C:59:ARG:HH21	1.64	0.46
1:D:107:ALA:HB1	1:D:202:ALA:HB1	1.97	0.46
1:B:32:ARG:HG2	1:B:211:TYR:HE1	1.82	0.45
1:B:165:LEU:CD1	2:B:401:37P:H12	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:TRP:CE2	1:A:66:ARG:HD3	2.51	0.45
1:D:99:ILE:O	1:D:103:ILE:HG13	2.16	0.45
1:A:133:LYS:HZ3	2:A:401:37P:H2	1.82	0.45
1:C:131:ASP:O	1:C:132:ARG:NH1	2.44	0.45
1:D:1:SER:OG	1:D:213:GLU:OE1	2.31	0.45
1:B:100:SER:HB2	1:B:148:HIS:ND1	2.32	0.45
1:C:57:THR:OG1	1:C:337:ARG:OXT	2.28	0.45
1:C:84:THR:OG1	1:C:196:HIS:ND1	2.33	0.45
1:B:272:LYS:NZ	1:D:242:GLU:OE2	2.37	0.45
1:D:159:SER:O	1:D:163:ASN:HB2	2.17	0.45
1:A:317:ASP:N	1:A:317:ASP:OD1	2.50	0.44
1:D:58:ARG:HH22	1:D:59:ARG:HH21	1.65	0.44
1:D:121:ASN:ND2	1:D:137:SER:O	2.50	0.44
1:A:301:LEU:HA	1:A:301:LEU:HD23	1.86	0.44
1:D:132:ARG:HH12	1:D:183:ARG:HD2	1.82	0.44
1:C:77:VAL:CG1	2:C:401:37P:H16	2.48	0.43
1:D:128:VAL:HG23	1:D:129:GLU:N	2.33	0.43
1:C:280:VAL:HG21	1:C:290:LEU:HB3	2.00	0.43
1:B:151:THR:HG23	2:B:401:37P:H21	1.99	0.43
1:A:65:VAL:HG21	1:A:164:TYR:CE2	2.54	0.43
1:A:183:ARG:H	1:A:183:ARG:NE	2.17	0.42
1:D:14:TRP:CE2	1:D:66:ARG:HD3	2.54	0.42
1:A:81:LEU:HD23	1:A:81:LEU:HA	1.81	0.42
1:C:111:LEU:HD13	1:C:202:ALA:HB2	2.01	0.42
1:B:50:ASN:HB3	1:B:53:LYS:HB3	2.02	0.42
1:C:304:ALA:HB1	1:C:333:ALA:HB1	2.02	0.42
1:C:75:GLY:H	2:C:401:37P:H13	1.68	0.42
1:C:47:ARG:HG3	1:C:71:SER:HA	2.01	0.41
1:D:63:ASP:OD1	1:D:63:ASP:N	2.53	0.41
1:A:312:GLU:HA	1:A:315:LEU:HD12	2.01	0.41
1:A:14:TRP:CD2	1:A:66:ARG:HD3	2.55	0.41
1:D:82:GLY:HA3	1:D:155:ASN:O	2.21	0.41
1:D:126:LYS:CD	1:D:131:ASP:OD1	2.68	0.41
1:C:83:ASN:ND2	1:C:85:CYS:SG	2.94	0.41
1:A:49:GLN:HG2	1:A:177:ILE:O	2.21	0.41
1:B:1:SER:OG	1:B:213:GLU:OE1	2.38	0.41
1:C:255:LEU:HB2	1:C:266:LEU:HB2	2.02	0.41
2:C:401:37P:H24	2:C:401:37P:H18	1.73	0.41
1:A:304:ALA:HB1	1:A:333:ALA:HB1	2.02	0.41
1:A:175:LYS:NZ	1:C:310:GLU:OE2	2.43	0.41
1:A:177:ILE:HD11	1:C:313:ALA:HA	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:216:GLU:CD	1:D:216:GLU:H	2.23	0.40
1:D:258:ARG:NH2	1:D:283:ASP:OD1	2.54	0.40
1:C:96:ASP:HB3	1:C:99:ILE:HG13	2.03	0.40
1:D:133:LYS:HZ3	2:D:401:37P:H2	1.87	0.40
1:D:284:SER:OG	1:D:285:LEU:N	2.54	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	335/341 (98%)	323 (96%)	12 (4%)	0	100	100
1	B	335/341 (98%)	321 (96%)	14 (4%)	0	100	100
1	C	335/341 (98%)	323 (96%)	11 (3%)	1 (0%)	41	75
1	D	335/341 (98%)	324 (97%)	11 (3%)	0	100	100
All	All	1340/1364 (98%)	1291 (96%)	48 (4%)	1 (0%)	51	84

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	336	VAL

### 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	279/282 (99%)	259 (93%)	20 (7%)	14	45
1	B	279/282 (99%)	264 (95%)	15 (5%)	22	55
1	C	279/282 (99%)	264 (95%)	15 (5%)	22	55
1	D	279/282 (99%)	259 (93%)	20 (7%)	14	45
All	All	1116/1128 (99%)	1046 (94%)	70 (6%)	18	51

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	30	THR
1	A	59	ARG
1	A	97	LYS
1	A	111	LEU
1	A	120	ARG
1	A	124	VAL
1	A	128	VAL
1	A	137	SER
1	A	139	TYR
1	A	183	ARG
1	A	188	THR
1	A	197	GLU
1	A	215	VAL
1	A	277	ARG
1	A	285	LEU
1	A	312	GLU
1	A	314	LEU
1	A	317	ASP
1	A	325	LEU
1	B	2	THR
1	B	47	ARG
1	B	97	LYS
1	B	111	LEU
1	B	128	VAL
1	B	139	TYR
1	B	154	LEU
1	B	160	ARG
1	B	215	VAL
1	B	237	ARG
1	B	285	LEU
1	B	305	ASP
1	B	312	GLU

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Mol	Chain	Res	Type
1	B	314	LEU
1	B	325	LEU
1	C	25	ARG
1	C	31	GLN
1	C	47	ARG
1	C	97	LYS
1	C	111	LEU
1	C	120	ARG
1	C	123	LEU
1	C	139	TYR
1	C	143	LYS
1	C	144	ASP
1	C	163	ASN
1	C	181	ARG
1	C	215	VAL
1	C	285	LEU
1	C	325	LEU
1	D	2	THR
1	D	4	ARG
1	D	63	ASP
1	D	97	LYS
1	D	109	ASN
1	D	124	VAL
1	D	126	LYS
1	D	131	ASP
1	D	132	ARG
1	D	163	ASN
1	D	181	ARG
1	D	188	THR
1	D	214	ARG
1	D	258	ARG
1	D	268	PHE
1	D	312	GLU
1	D	314	LEU
1	D	317	ASP
1	D	321	GLN
1	D	325	LEU

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 ⓘ

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	37P	A	401	-	45,50,50	1.22	5 (11%)	50,74,74	1.65	9 (18%)
2	37P	B	401	-	45,50,50	1.22	5 (11%)	50,74,74	1.67	10 (20%)
2	37P	C	401	-	45,50,50	1.20	5 (11%)	50,74,74	1.69	10 (20%)
2	37P	D	401	-	45,50,50	1.22	5 (11%)	50,74,74	1.60	10 (20%)

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	37P	A	401	-	1/1/6/7	1/17/42/42	0/5/6/6
2	37P	B	401	-	1/1/6/7	2/17/42/42	0/5/6/6
2	37P	C	401	-	-	9/17/42/42	0/5/6/6
2	37P	D	401	-	1/1/6/7	6/17/42/42	0/5/6/6

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	37P	O5'-S	-3.46	1.52	1.59
2	A	401	37P	O5'-S	-3.40	1.52	1.59
2	B	401	37P	O5'-S	-3.32	1.52	1.59
2	C	401	37P	O5'-S	-3.12	1.53	1.59
2	D	401	37P	S-N45	3.06	1.65	1.59
2	B	401	37P	C44-N45	-2.88	1.33	1.38
2	C	401	37P	C44-N45	-2.83	1.33	1.38
2	A	401	37P	C12-C13	2.79	1.49	1.41
2	D	401	37P	C12-C13	2.79	1.49	1.41
2	C	401	37P	C12-C13	2.76	1.49	1.41
2	A	401	37P	S-N45	2.75	1.64	1.59
2	A	401	37P	C44-N45	-2.70	1.33	1.38
2	C	401	37P	S-N45	2.69	1.64	1.59
2	B	401	37P	S-N45	2.61	1.64	1.59
2	B	401	37P	C12-C13	2.57	1.48	1.41
2	D	401	37P	C5A-C4A	2.55	1.47	1.40
2	B	401	37P	C5A-C4A	2.53	1.47	1.40
2	D	401	37P	C44-N45	-2.50	1.34	1.38
2	A	401	37P	C5A-C4A	2.45	1.47	1.40
2	C	401	37P	C5A-C4A	2.38	1.47	1.40

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	37P	C6-C11-N10	-4.95	120.25	125.73
2	B	401	37P	C6-C11-N10	-4.93	120.27	125.73
2	A	401	37P	C6-C11-N10	-4.84	120.37	125.73
2	D	401	37P	C6-C11-N10	-4.73	120.50	125.73
2	A	401	37P	O7-C6-C1	4.51	121.39	116.11
2	B	401	37P	O7-C6-C1	4.44	121.31	116.11
2	D	401	37P	O7-C6-C1	4.10	120.91	116.11
2	C	401	37P	O7-C6-C1	4.07	120.88	116.11
2	B	401	37P	O7-C8-C17	3.85	120.60	116.03
2	A	401	37P	O7-C8-C17	3.68	120.40	116.03
2	B	401	37P	C1-C6-C11	-3.57	118.66	122.90
2	A	401	37P	C1-C6-C11	-3.51	118.74	122.90
2	C	401	37P	O7-C8-C17	3.46	120.13	116.03
2	D	401	37P	O7-C8-C17	3.24	119.87	116.03
2	C	401	37P	C4A-C5A-N7A	-3.20	106.07	109.40
2	C	401	37P	C1-C6-C11	-3.16	119.16	122.90
2	D	401	37P	C1-C6-C11	-3.16	119.16	122.90
2	A	401	37P	C9-N10-C11	3.14	120.40	118.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	37P	N3A-C2A-N1A	-3.11	123.81	128.68
2	C	401	37P	C9-N10-C11	3.09	120.36	118.06
2	C	401	37P	N3A-C2A-N1A	-3.09	123.85	128.68
2	D	401	37P	N3A-C2A-N1A	-3.07	123.89	128.68
2	B	401	37P	C9-N10-C11	3.05	120.33	118.06
2	B	401	37P	N3A-C2A-N1A	-3.02	123.96	128.68
2	D	401	37P	C9-N10-C11	2.84	120.17	118.06
2	B	401	37P	C4A-C5A-N7A	-2.79	106.50	109.40
2	D	401	37P	C4A-C5A-N7A	-2.71	106.58	109.40
2	C	401	37P	C44-N45-S	-2.59	120.09	124.30
2	C	401	37P	O5'-C5'-C4'	2.51	112.30	107.62
2	B	401	37P	O5'-C5'-C4'	2.49	112.26	107.62
2	A	401	37P	C4A-C5A-N7A	-2.48	106.82	109.40
2	D	401	37P	O1S-S-O2S	-2.40	117.02	120.76
2	A	401	37P	O5'-C5'-C4'	2.34	112.00	107.62
2	B	401	37P	C4-C11-C6	2.27	119.06	116.06
2	D	401	37P	O5'-C5'-C4'	2.17	111.66	107.62
2	A	401	37P	C4-C11-C6	2.15	118.91	116.06
2	D	401	37P	C4-C11-C6	2.14	118.90	116.06
2	C	401	37P	C4-C11-C6	2.06	118.80	116.06
2	B	401	37P	C16-C15-C13	-2.06	107.90	114.10

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	401	37P	C1'
2	B	401	37P	C1'
2	D	401	37P	C1'

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	37P	C44-N45-S-O2S
2	C	401	37P	O46-C44-N45-S
2	D	401	37P	C42-C43-C44-N45
2	D	401	37P	C42-C43-C44-O46
2	C	401	37P	C15-C16-C42-C43
2	C	401	37P	C13-C15-C16-C42
2	D	401	37P	C16-C42-C43-C44
2	D	401	37P	C14-C13-C15-C16
2	B	401	37P	C3'-C4'-C5'-O5'
2	C	401	37P	C5'-O5'-S-O2S

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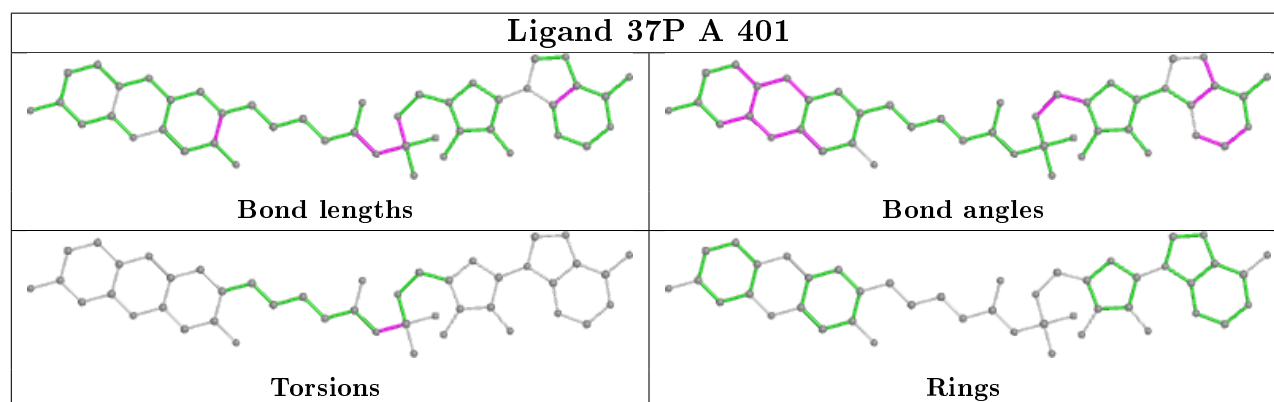
Mol	Chain	Res	Type	Atoms
2	C	401	37P	C5'-O5'-S-O1S
2	B	401	37P	C44-N45-S-O2S
2	D	401	37P	C13-C15-C16-C42
2	D	401	37P	C12-C13-C15-C16
2	C	401	37P	C43-C44-N45-S
2	C	401	37P	C16-C42-C43-C44
2	C	401	37P	C42-C43-C44-N45
2	C	401	37P	C42-C43-C44-O46

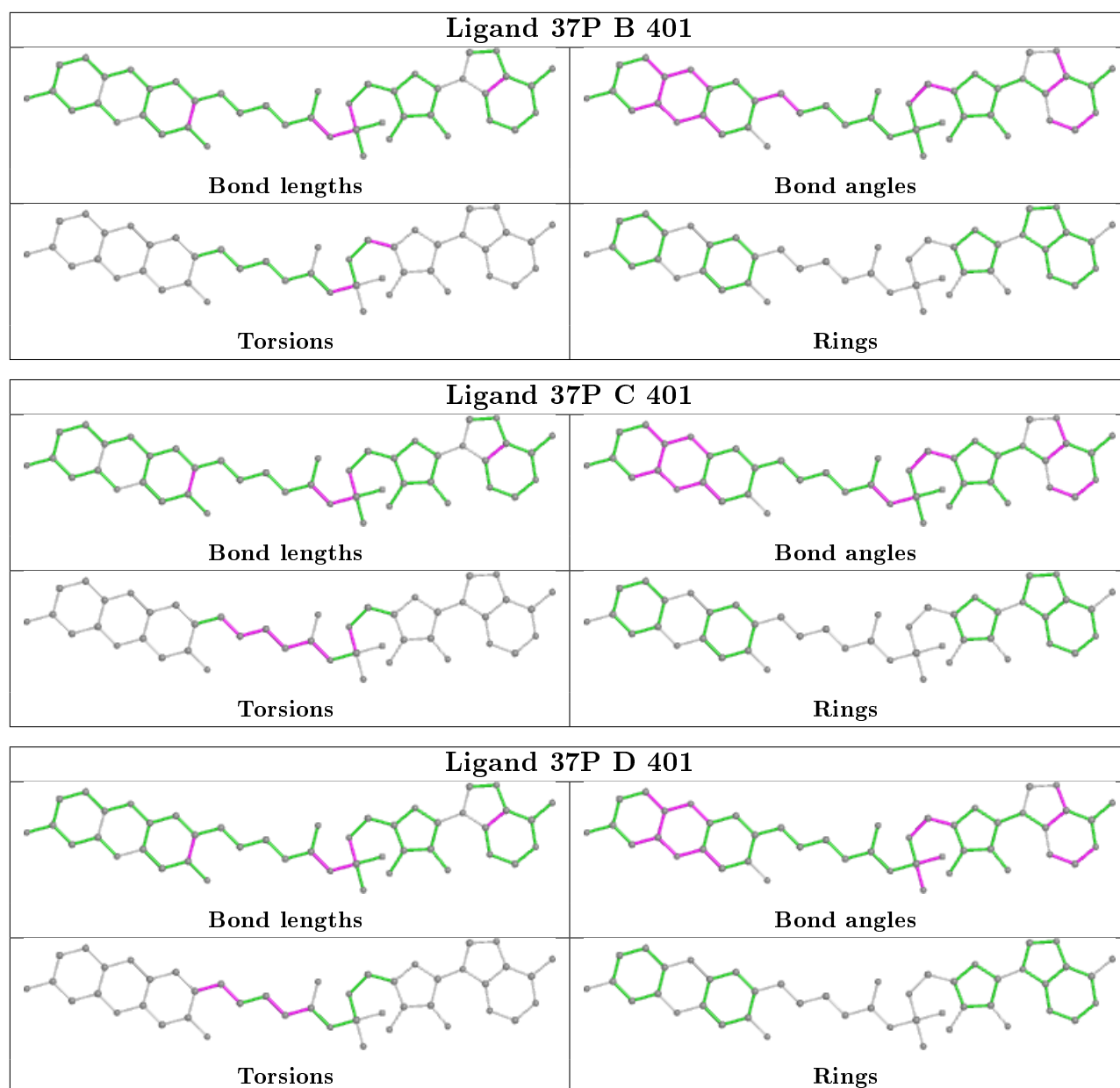
There are no ring outliers.

4 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	37P	4	0
2	B	401	37P	4	0
2	C	401	37P	9	0
2	D	401	37P	2	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, 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(Å²)	Q<0.9	
1	A	337/341 (98%)	-0.48	0	100	100	22, 41, 74, 99	0
1	B	337/341 (98%)	-0.43	0	100	100	20, 43, 77, 104	0
1	C	337/341 (98%)	-0.16	0	100	100	22, 63, 125, 158	0
1	D	337/341 (98%)	-0.07	2 (0%)	89	86	23, 72, 123, 154	0
All	All	1348/1364 (98%)	-0.28	2 (0%)	95	95	20, 51, 116, 158	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	5	LEU	2.4
1	D	143	LYS	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](#)

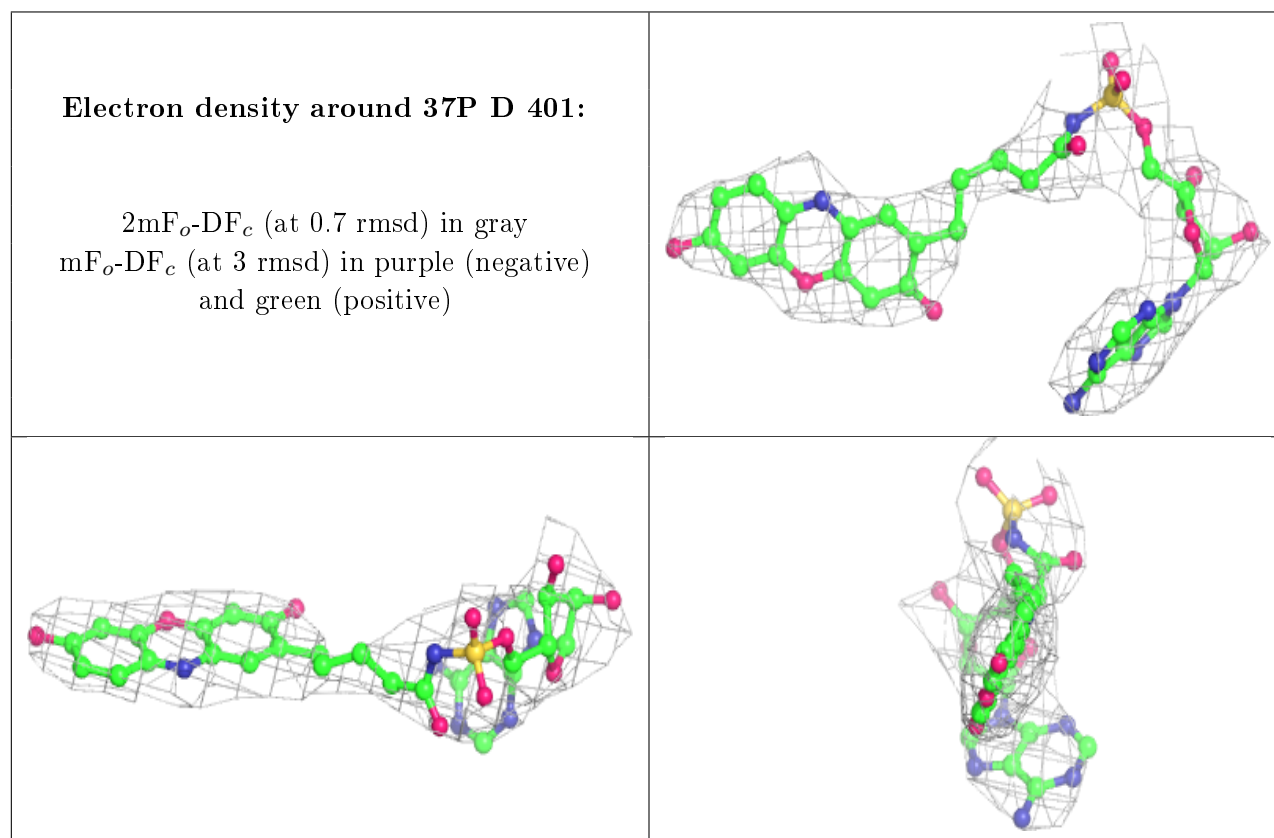
There are no carbohydrates in this entry.

### 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.

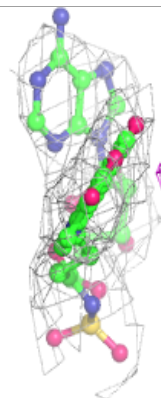
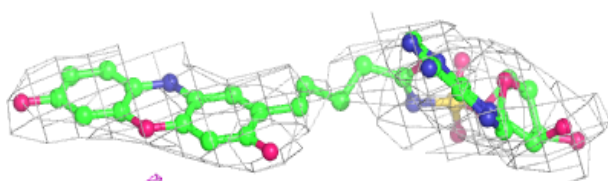
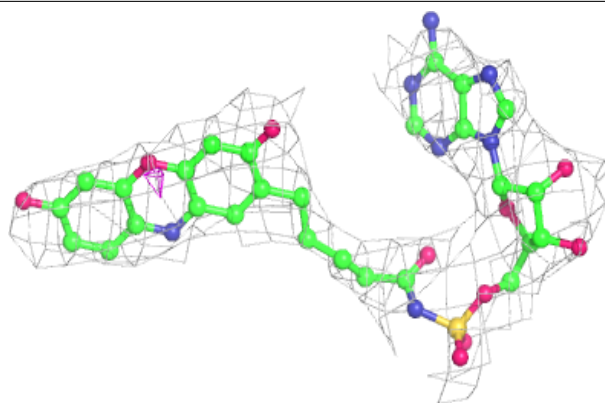
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	37P	D	401	45/45	0.93	0.27	56,66,78,82	0
2	37P	C	401	45/45	0.94	0.28	34,57,67,71	0
2	37P	A	401	45/45	0.97	0.23	22,31,43,45	0
2	37P	B	401	45/45	0.97	0.24	20,37,48,50	0

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.

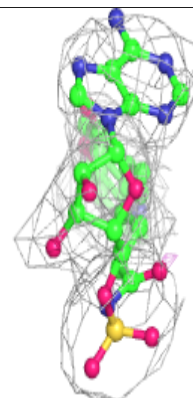
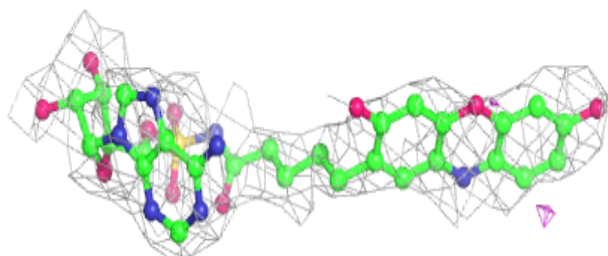
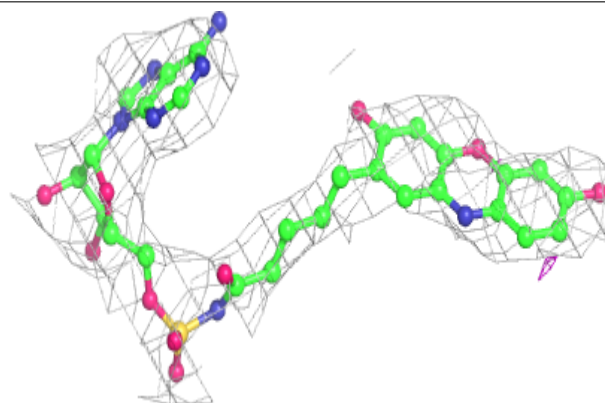


**Electron density around 37P C 401:**

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

**Electron density around 37P A 401:**

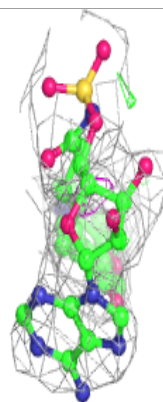
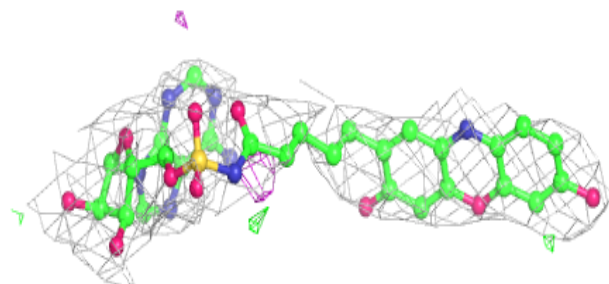
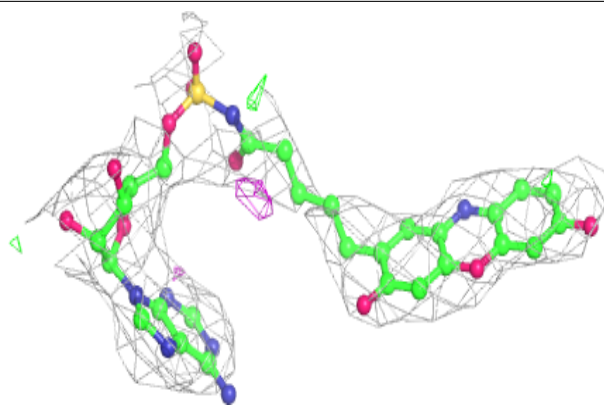
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around 37P 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)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.