



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 13, 2023 – 12:03 am GMT

PDB ID : 2XUO  
Title : CRYSTAL STRUCTURE OF MACHE-Y337A mutant in complex with soaked TZ2PA6 ANTI inhibitor  
Authors : Bourne, Y.; Radic, Z.; Taylor, P.; Marchot, P.  
Deposited on : 2010-10-19  
Resolution : 2.80 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

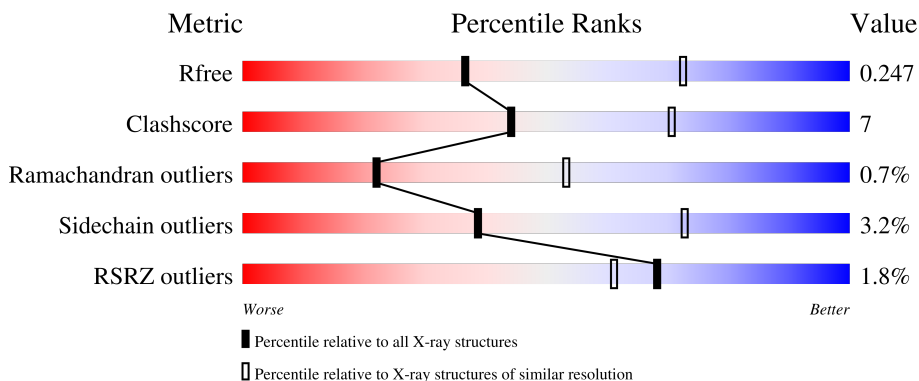
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


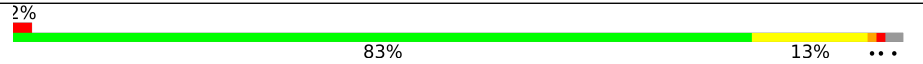
The reported resolution of this entry is 2.80 Å.

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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 of 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	543	 2% 82% 16% ..
1	B	543	 2% 83% 13% ...

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8515 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.

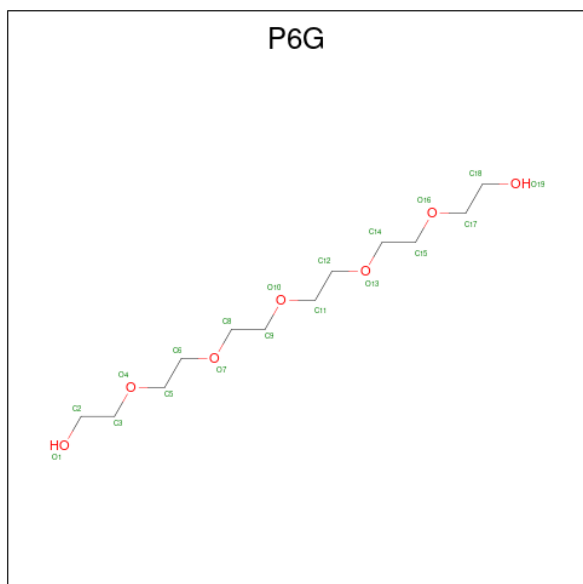
- Molecule 1 is a protein called ACETYLCHOLINESTERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	536	4203	2699	730	760	14	0	3	0
1	B	532	4156	2668	718	756	14	0	2	0

There are 2 discrepancies between the modelled and reference sequences:

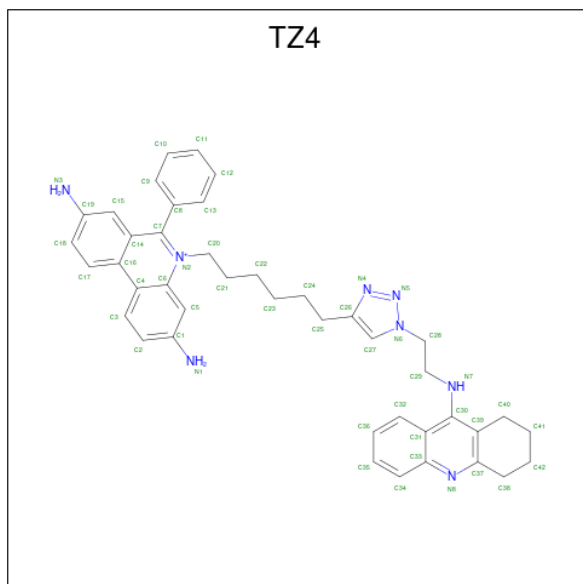
Chain	Residue	Modelled	Actual	Comment	Reference
A	337	ALA	TYR	engineered mutation	UNP P21836
B	337	ALA	TYR	engineered mutation	UNP P21836

- Molecule 2 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: C<sub>12</sub>H<sub>26</sub>O<sub>7</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	19	12	7	0	0

- Molecule 3 is 3,8-DIAMINO-6-PHENYL-5-[6-[1-[2-[(1,2,3,4-TETRAHYDRO-9-ACRIDINYL)AMINO]ETHYL]-1H-1,2,3-TRIAZOL-4-YL]HEXYL]-PHENANTHRIDINIUM (three-letter code: TZ4) (formula: C<sub>42</sub>H<sub>45</sub>N<sub>8</sub>).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C N	0	0
			50	42 8		
3	B	1	Total	C N	0	0
			50	42 8		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	24	Total	O	0	0
			24	24		
4	B	13	Total	O	0	0
			13	13		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.91Å 110.41Å 227.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 19.92 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.00-2.80) 99.9 (19.92-2.80)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.52 (at 2.79Å)	Xtrriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.199 , 0.247 0.202 , 0.247	Depositor DCC
$R_{free}$ test set	1020 reflections (2.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	63.6	Xtrriage
Anisotropy	0.071	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 42.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8515	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% 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: P6G, TZ4

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	0.57	0/4339	0.67	1/5931 (0.0%)
1	B	0.54	0/4287	0.83	6/5860 (0.1%)
All	All	0.56	0/8626	0.75	7/11791 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	4
All	All	0	5

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	322	GLN	CA-C-N	24.25	170.54	117.20
1	B	322	GLN	O-C-N	-20.19	90.40	122.70
1	B	322	GLN	CA-C-O	-18.20	81.87	120.10
1	B	525	ARG	N-CA-C	-10.42	82.87	111.00
1	B	524	LEU	CA-CB-CG	6.47	130.19	115.30
1	B	387	HIS	C-N-CD	-5.59	108.31	120.60
1	A	322	GLN	N-CA-C	5.51	125.88	111.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	322	GLN	Peptide
1	B	322	GLN	Mainchain
1	B	387	HIS	Peptide
1	B	524	LEU	Peptide
1	B	525	ARG	Peptide

## 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	4203	0	4095	61	0
1	B	4156	0	4044	46	0
2	A	19	0	26	1	0
3	A	50	0	45	6	0
3	B	50	0	45	5	0
4	A	24	0	0	0	0
4	B	13	0	0	1	0
All	All	8515	0	8255	110	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 (110) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286[B]:TRP:CZ2	3:A:1543:TZ4:H211	1.89	1.06
1:A:322:GLN:N	1:A:323:ASP:HB3	1.68	1.05
1:A:376:GLU:HG3	4:B:2013:HOH:O	1.60	1.02
1:A:4:GLU:OE1	1:A:18:ARG:HD3	1.69	0.91
1:A:496:LYS:HA	1:A:497:SER:O	1.72	0.88
1:A:113:PRO:HG2	1:A:485:ARG:HG2	1.64	0.79
1:B:497:SER:HB2	1:B:498:PRO:C	2.07	0.75
1:A:496:LYS:HA	1:A:497:SER:C	2.06	0.73
1:A:437:LEU:HD11	1:A:449:TYR:CD2	2.24	0.73
3:A:1543:TZ4:H402	3:A:1543:TZ4:H292	1.71	0.71
1:A:459:LEU:HD23	1:A:470:ARG:HG2	1.71	0.70
1:B:458:PRO:HA	1:B:465:TYR:CD2	2.26	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:THR:HG21	1:B:143:GLY:O	1.93	0.68
1:B:211:MET:HG2	1:B:308:LEU:HD21	1.77	0.67
1:B:113:PRO:HG2	1:B:485:ARG:HG2	1.78	0.65
1:A:385:TRP:HD1	1:B:527:GLN:OE1	1.80	0.64
1:B:286:TRP:HB2	3:B:1542:TZ4:C3	2.28	0.64
1:B:285:GLU:HG3	3:B:1542:TZ4:H2	1.80	0.63
1:A:458:PRO:HA	1:A:465:TYR:CD2	2.34	0.63
1:A:322:GLN:N	1:A:323:ASP:CB	2.56	0.62
1:B:224:ARG:HG2	1:B:325:GLN:HB2	1.81	0.60
1:B:103:THR:HG22	1:B:145:VAL:HG22	1.84	0.59
1:A:293:SER:HA	1:A:365:ILE:HG23	1.86	0.57
1:A:322:GLN:CA	1:A:323:ASP:HB3	2.34	0.57
1:B:329:GLY:HA3	1:B:428:TYR:CZ	2.40	0.57
1:A:197:VAL:H	1:A:223[A]:HIS:HD2	1.53	0.57
1:A:285:GLU:HG3	3:A:1543:TZ4:H2	1.87	0.56
1:B:213:ILE:O	1:B:219:ARG:HD3	2.05	0.56
1:A:226:VAL:HG13	1:A:327:LEU:HD23	1.88	0.56
1:B:113:PRO:HG2	1:B:485:ARG:CG	2.36	0.54
1:B:497:SER:HB2	1:B:498:PRO:O	2.08	0.53
1:B:328:VAL:O	1:B:427:ALA:HA	2.08	0.53
1:A:103:THR:HG22	1:A:145:VAL:HG13	1.91	0.53
1:A:413:GLN:O	1:A:414:LEU:C	2.47	0.53
1:B:329:GLY:HA3	1:B:428:TYR:CE2	2.44	0.53
1:A:321:PHE:C	1:A:323:ASP:HB3	2.29	0.52
1:B:353:LEU:HB3	1:B:391:PRO:HB2	1.92	0.52
1:A:294:ILE:HD11	1:A:402:VAL:HG21	1.92	0.51
3:A:1543:TZ4:H292	3:A:1543:TZ4:C40	2.41	0.51
1:B:495:SER:OG	1:B:496:LYS:N	2.43	0.50
1:B:66:GLN:HG3	1:B:98:TYR:CG	2.46	0.50
3:B:1542:TZ4:H402	3:B:1542:TZ4:H292	1.95	0.49
1:A:286[B]:TRP:CE3	1:A:286[B]:TRP:HA	2.47	0.49
1:A:459:LEU:HD21	1:A:474:GLN:HG2	1.95	0.49
1:B:528:THR:O	1:B:531:PHE:HB3	2.13	0.48
1:A:328:VAL:O	1:A:427:ALA:HA	2.13	0.48
1:B:265:ASN:HD22	1:B:266:ASP:N	2.12	0.48
3:B:1542:TZ4:H5	3:B:1542:TZ4:C21	2.43	0.48
1:A:286[B]:TRP:CH2	3:A:1543:TZ4:H211	2.45	0.47
1:A:527:GLN:HE21	2:A:1542:P6G:H181	1.79	0.47
1:A:364:ARG:HH22	1:A:372:ASP:CG	2.18	0.47
1:A:43:GLY:C	1:A:45:ARG:H	2.16	0.47
1:A:528:THR:O	1:A:531:PHE:HB3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:ARG:HH21	1:B:107:ARG:HG2	1.79	0.47
1:A:522:ARG:HB3	1:B:386:LEU:HD21	1.97	0.47
1:A:29:VAL:CG1	1:A:140:GLN:HB2	2.45	0.47
1:A:112:THR:HG21	1:A:143:GLY:O	2.14	0.47
1:A:470:ARG:O	1:A:474:GLN:HG3	2.16	0.46
1:B:66:GLN:HG3	1:B:98:TYR:CD1	2.50	0.46
1:B:384:ASP:O	1:B:388:PRO:HA	2.15	0.46
1:A:223[B]:HIS:HD2	1:A:224:ARG:NH2	2.13	0.46
1:A:29:VAL:HG21	1:A:136:ARG:HB2	1.97	0.46
1:A:286[A]:TRP:HB2	3:A:1543:TZ4:C3	2.46	0.46
1:B:357:ALA:HA	1:B:360:LEU:HD12	1.97	0.46
1:A:122:GLY:O	1:A:123:PHE:HB2	2.16	0.46
1:B:177:ARG:CZ	1:B:217:PRO:HB2	2.46	0.45
1:B:379:VAL:O	1:B:383:THR:OG1	2.34	0.45
1:B:304:ASP:OD2	1:B:306:ASP:HB3	2.16	0.45
1:A:322:GLN:CA	1:A:323:ASP:CB	2.95	0.45
1:B:80:PHE:CE2	1:B:82:GLY:HA3	2.52	0.45
1:A:276:ARG:HA	1:A:276:ARG:HD3	1.81	0.45
1:A:161:LEU:HD11	1:A:269:LEU:HD22	1.98	0.45
1:B:524:LEU:HD13	1:B:524:LEU:O	2.17	0.45
1:B:336:SER:HB2	1:B:443:MET:HG2	1.98	0.44
1:A:38:ALA:HB2	1:A:178:LEU:HD23	2.00	0.44
1:A:434:ALA:HB3	1:A:437:LEU:HD13	2.00	0.44
1:B:32:PHE:HB3	1:B:35:ILE:HD11	1.99	0.44
1:A:103:THR:HG21	1:A:190:PHE:HB3	2.00	0.44
1:A:327:LEU:HD11	1:A:500:TRP:CH2	2.53	0.44
1:A:258:PRO:HA	1:A:259:PRO:HD2	1.84	0.43
1:B:24:ALA:HB3	1:B:27:GLY:O	2.18	0.43
1:A:75:THR:O	1:A:78:PRO:HD3	2.17	0.43
1:B:353:LEU:HB3	1:B:391:PRO:CB	2.47	0.43
1:A:33:LEU:HD23	1:A:100:ASN:HB3	1.99	0.43
1:A:112:THR:HB	1:A:145:VAL:HG23	2.00	0.43
1:B:360:LEU:HD23	1:B:379:VAL:HG11	1.99	0.43
1:A:322:GLN:H	1:A:323:ASP:HB3	1.70	0.42
1:B:197:VAL:H	1:B:223:HIS:HD2	1.67	0.42
1:B:265:ASN:HD22	1:B:265:ASN:C	2.22	0.42
1:B:84:GLU:HG3	1:B:87:ASN:HD22	1.83	0.42
1:A:224:ARG:HG2	1:A:325:GLN:HB2	2.02	0.42
1:B:77:TYR:CZ	1:B:348:LYS:HG2	2.54	0.42
1:A:128:ALA:HB1	1:A:148:SER:OG	2.20	0.42
1:A:183:VAL:HG13	1:A:187:ILE:HB	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:257:CYS:HA	1:B:258:PRO:HD3	1.89	0.42
1:B:300:VAL:HB	1:B:301:PRO:HD2	2.02	0.42
1:B:460:ASP:HB3	1:B:463:LEU:HD12	2.01	0.42
1:A:226:VAL:HG22	1:A:327:LEU:HB3	2.02	0.42
1:B:197:VAL:H	1:B:223:HIS:CD2	2.37	0.42
1:A:8:LEU:HD11	1:A:21:ARG:HB2	2.02	0.41
1:A:36:PRO:HB2	1:A:53:LYS:HD3	2.01	0.41
3:B:1542:TZ4:H252	3:B:1542:TZ4:H421	2.03	0.41
1:A:33:LEU:HD22	1:A:65:PHE:CE1	2.56	0.41
1:B:511:VAL:HG11	1:B:518:LEU:HD13	2.01	0.41
1:A:540:LEU:O	1:A:541:SER:C	2.58	0.41
1:A:231:THR:HB	1:A:233:ASN:OD1	2.21	0.41
1:B:24:ALA:O	1:B:25:PRO:C	2.59	0.41
1:A:213:ILE:HD11	1:A:225:ALA:CB	2.51	0.40
1:A:491:ASP:HA	1:A:492:PRO:HD3	1.99	0.40
1:A:524:LEU:O	1:A:525:ARG:C	2.60	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	535/543 (98%)	506 (95%)	26 (5%)	3 (1%)	25 56
1	B	530/543 (98%)	500 (94%)	26 (5%)	4 (1%)	19 49
All	All	1065/1086 (98%)	1006 (94%)	52 (5%)	7 (1%)	22 53

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	525	ARG
1	A	323	ASP

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Mol	Chain	Res	Type
1	B	495	SER
1	B	497	SER
1	A	497	SER
1	A	523	GLY
1	B	523	GLY

### 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	443/442 (100%)	428 (97%)	15 (3%)	37	71
1	B	439/442 (99%)	425 (97%)	14 (3%)	39	73
All	All	882/884 (100%)	853 (97%)	29 (3%)	39	72

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	70	TYR
1	A	75	THR
1	A	81	GLU
1	A	84	GLU
1	A	200	PHE
1	A	216	LEU
1	A	286[A]	TRP
1	A	286[B]	TRP
1	A	295	PHE
1	A	322	GLN
1	A	326	VAL
1	A	327	LEU
1	A	372	ASP
1	A	524	LEU
1	B	84	GLU
1	B	100	ASN
1	B	161	LEU

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Mol	Chain	Res	Type
1	B	200	PHE
1	B	265	ASN
1	B	295	PHE
1	B	352	SER
1	B	360	LEU
1	B	386	LEU
1	B	413	GLN
1	B	437	LEU
1	B	497	SER
1	B	524	LEU
1	B	536	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	287	HIS
1	A	509	GLN
1	A	527	GLN
1	B	223	HIS
1	B	265	ASN

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

3 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
3	TZ4	A	1543	-	56,57,57	1.74	13 (23%)	65,80,80	1.64	12 (18%)
2	P6G	A	1542	-	18,18,18	2.17	6 (33%)	17,17,17	1.28	2 (11%)
3	TZ4	B	1542	-	56,57,57	1.70	11 (19%)	65,80,80	1.79	13 (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
3	TZ4	A	1543	-	-	4/15/26/26	0/8/8/8
2	P6G	A	1542	-	-	10/16/16/16	-
3	TZ4	B	1542	-	-	4/15/26/26	0/8/8/8

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1542	TZ4	C38-C37	-4.41	1.43	1.50
2	A	1542	P6G	O10-C11	4.34	1.60	1.42
3	B	1542	TZ4	C40-C39	-4.26	1.44	1.51
3	B	1542	TZ4	C41-C40	-4.07	1.36	1.51
3	A	1543	TZ4	C38-C37	-4.03	1.44	1.50
3	A	1543	TZ4	C40-C39	-3.99	1.45	1.51
3	A	1543	TZ4	C41-C40	-3.93	1.37	1.51
3	B	1542	TZ4	C42-C38	-3.86	1.37	1.51
2	A	1542	P6G	O7-C8	3.81	1.58	1.42
3	A	1543	TZ4	C6-N2	3.69	1.43	1.38
3	A	1543	TZ4	C42-C38	-3.63	1.38	1.51
3	B	1542	TZ4	C6-N2	3.54	1.43	1.38
2	A	1542	P6G	O4-C5	3.51	1.57	1.42
2	A	1542	P6G	O1-C2	3.47	1.60	1.42
2	A	1542	P6G	O16-C17	3.39	1.56	1.42
3	A	1543	TZ4	C34-C33	2.73	1.46	1.41
3	B	1542	TZ4	C37-N8	2.71	1.35	1.32
2	A	1542	P6G	O13-C14	2.68	1.53	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1543	TZ4	C37-N8	2.56	1.35	1.32
3	A	1543	TZ4	C36-C35	2.47	1.44	1.38
3	B	1542	TZ4	C30-C39	2.35	1.42	1.38
3	A	1543	TZ4	C39-C37	2.34	1.44	1.40
3	B	1542	TZ4	C15-C14	2.26	1.45	1.41
3	A	1543	TZ4	C4-C6	2.17	1.45	1.41
3	B	1542	TZ4	C42-C41	-2.14	1.43	1.51
3	B	1542	TZ4	C27-C26	2.11	1.39	1.36
3	A	1543	TZ4	C15-C14	2.08	1.44	1.41
3	A	1543	TZ4	C30-C39	2.06	1.42	1.38
3	B	1542	TZ4	C17-C16	2.01	1.45	1.41
3	A	1543	TZ4	C17-C16	2.00	1.45	1.41

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1542	TZ4	N4-N5-N6	5.49	111.45	107.31
3	A	1543	TZ4	N4-N5-N6	5.08	111.14	107.31
3	B	1542	TZ4	C40-C39-C37	-4.57	116.80	121.08
3	B	1542	TZ4	C41-C40-C39	4.20	121.47	112.84
3	A	1543	TZ4	C40-C39-C37	-4.09	117.26	121.08
3	B	1542	TZ4	C28-N6-C27	-3.85	120.76	129.82
3	A	1543	TZ4	C42-C38-C37	3.77	119.73	113.53
3	A	1543	TZ4	C28-N6-C27	-3.76	120.97	129.82
3	A	1543	TZ4	C41-C40-C39	3.76	120.56	112.84
3	B	1542	TZ4	C39-C30-N7	3.64	126.20	119.54
3	B	1542	TZ4	C42-C38-C37	3.53	119.32	113.53
3	B	1542	TZ4	C20-N2-C6	-3.24	116.07	118.86
3	A	1543	TZ4	C19-C15-C14	3.21	123.68	120.90
3	B	1542	TZ4	C5-C6-C4	-3.20	116.55	120.61
2	A	1542	P6G	O19-C18-C17	3.14	130.01	111.81
3	A	1543	TZ4	C15-C14-C7	-3.06	118.56	124.00
3	B	1542	TZ4	C29-C28-N6	2.92	115.64	110.90
3	A	1543	TZ4	C5-C6-C4	-2.90	116.93	120.61
3	B	1542	TZ4	C7-N2-C6	2.77	125.60	121.90
3	A	1543	TZ4	C39-C30-N7	2.62	124.35	119.54
3	B	1542	TZ4	C19-C15-C14	2.36	122.95	120.90
3	A	1543	TZ4	C20-N2-C6	-2.33	116.86	118.86
3	B	1542	TZ4	C32-C31-C33	-2.20	115.86	118.33
3	B	1542	TZ4	C3-C4-C16	-2.17	117.87	122.39
2	A	1542	P6G	O7-C8-C9	2.16	120.13	110.39
3	A	1543	TZ4	C17-C16-C14	-2.10	115.63	118.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1543	TZ4	C29-C28-N6	2.01	114.16	110.90

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1542	TZ4	N2-C20-C21-C22
3	A	1543	TZ4	N2-C20-C21-C22
2	A	1542	P6G	O1-C2-C3-O4
2	A	1542	P6G	O4-C5-C6-O7
2	A	1542	P6G	O16-C17-C18-O19
2	A	1542	P6G	O10-C11-C12-O13
3	A	1543	TZ4	C31-C30-N7-C29
3	B	1542	TZ4	C31-C30-N7-C29
2	A	1542	P6G	O13-C14-C15-O16
2	A	1542	P6G	C15-C14-O13-C12
3	A	1543	TZ4	C20-C21-C22-C23
3	B	1542	TZ4	C39-C30-N7-C29
2	A	1542	P6G	C18-C17-O16-C15
3	B	1542	TZ4	C24-C25-C26-C27
3	A	1543	TZ4	C39-C30-N7-C29
2	A	1542	P6G	C11-C12-O13-C14
2	A	1542	P6G	C14-C15-O16-C17
2	A	1542	P6G	C2-C3-O4-C5

There are no ring outliers.

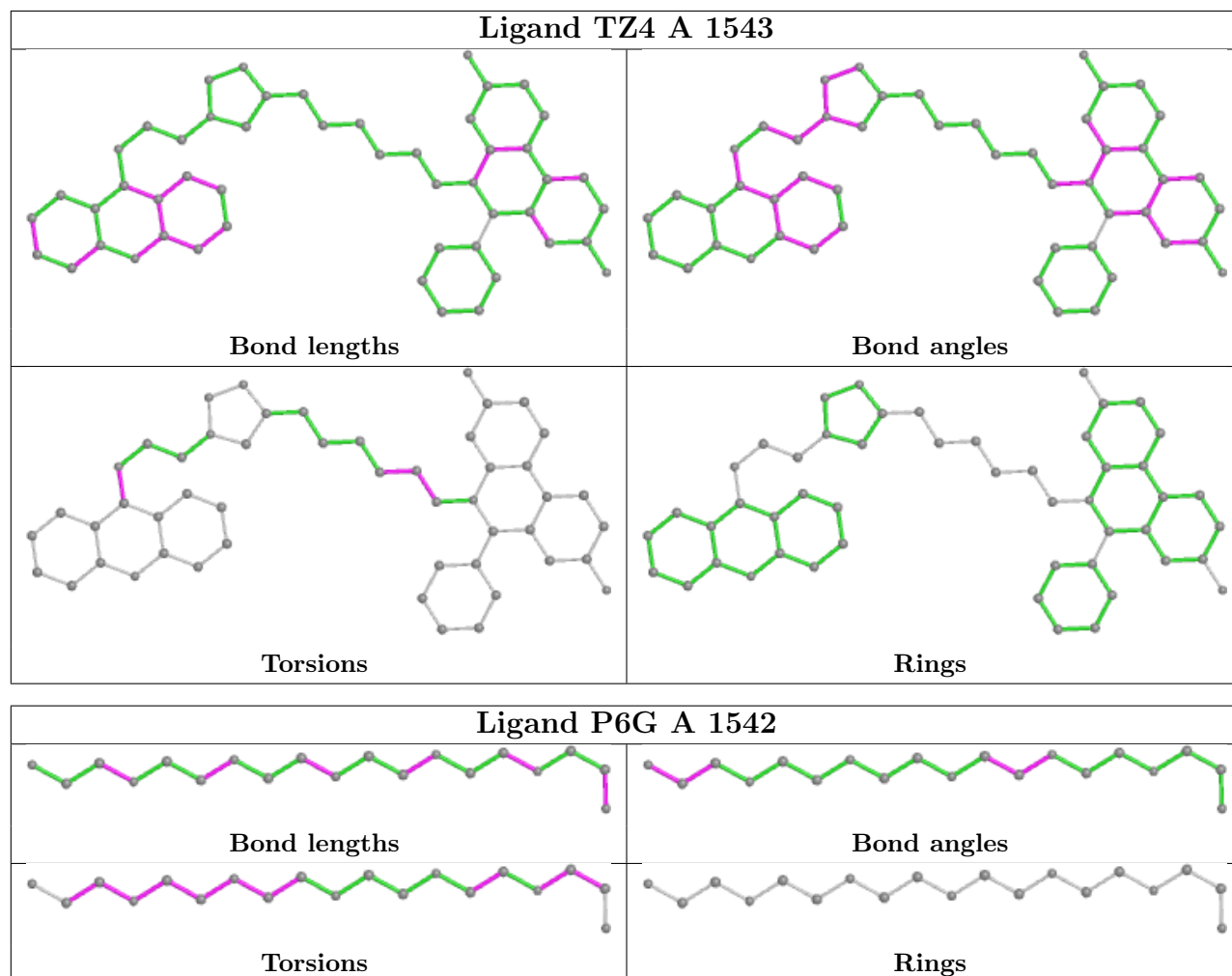
3 monomers are involved in 12 short contacts:

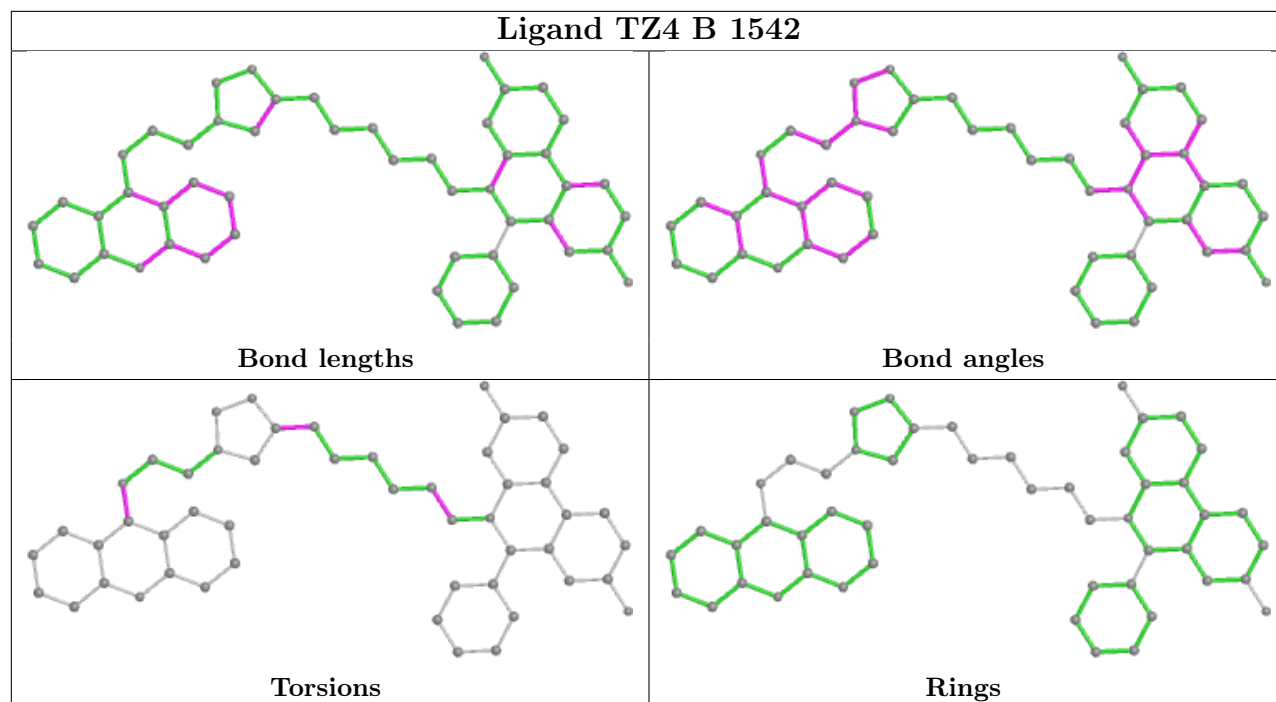
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1543	TZ4	6	0
2	A	1542	P6G	1	0
3	B	1542	TZ4	5	0

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(Å <sup>2</sup> )	Q<0.9
1	A	536/543 (98%)	-0.47	9 (1%) 70 63	35, 54, 78, 111	0
1	B	532/543 (97%)	-0.43	10 (1%) 66 59	42, 60, 87, 103	0
All	All	1068/1086 (98%)	-0.45	19 (1%) 68 61	35, 57, 85, 111	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	541	SER	5.8
1	A	259	PRO	5.4
1	A	495	SER	4.6
1	B	541	SER	3.8
1	B	540	LEU	3.4
1	A	1	GLU	3.2
1	B	323	ASP	3.2
1	A	540	LEU	3.1
1	B	495	SER	3.1
1	B	497	SER	2.9
1	A	494	ASP	2.9
1	B	268	GLU	2.7
1	A	493	ARG	2.7
1	A	496	LYS	2.5
1	B	109	ALA	2.3
1	A	165	ARG	2.2
1	B	110	SER	2.1
1	B	494	ASP	2.0
1	B	78	PRO	2.0

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

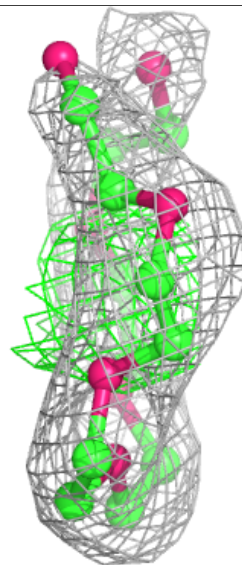
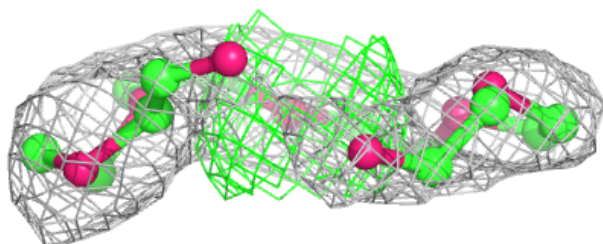
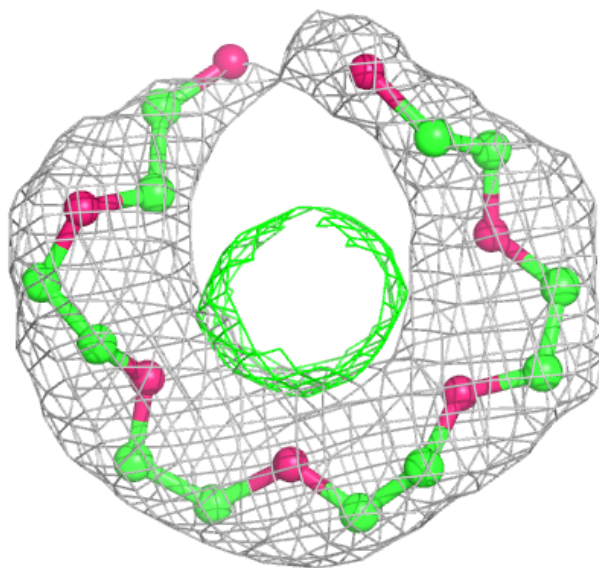
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	P6G	A	1542	19/19	0.89	0.26	59,63,74,76	0
3	TZ4	B	1542	50/50	0.90	0.20	48,70,91,91	0
3	TZ4	A	1543	50/50	0.91	0.20	42,55,82,82	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.

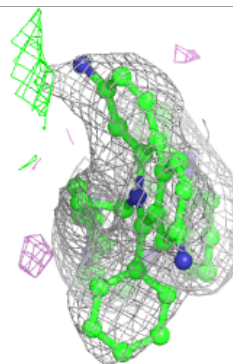
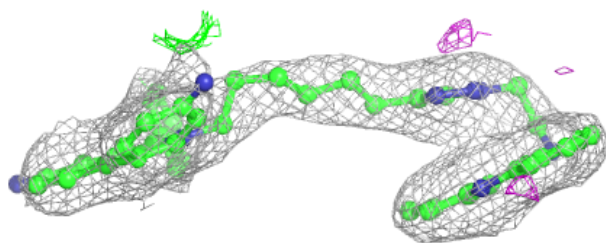
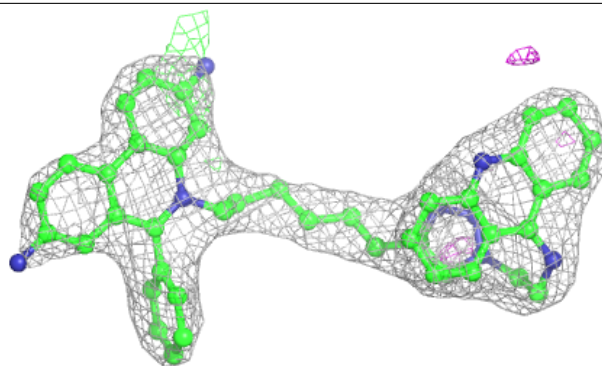
**Electron density around P6G A 1542:**

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

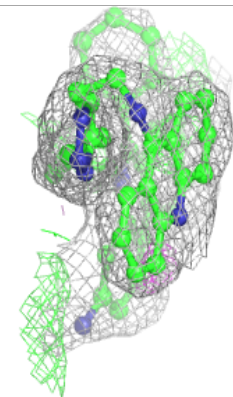
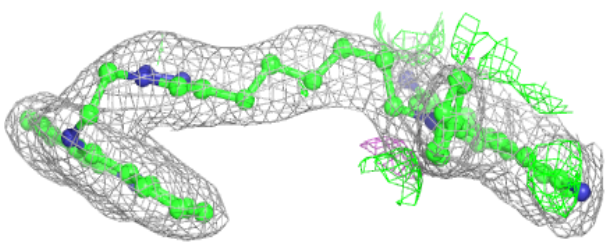
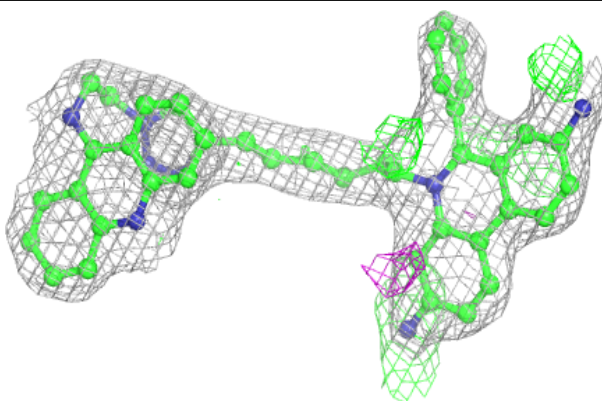


**Electron density around TZ4 B 1542:**

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

**Electron density around TZ4 A 1543:**

$2mF_o-DF_c$  (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.