



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 05:09 am BST

PDB ID : 3UVV  
Title : Crystal Structure of the ligand binding domains of the thyroid receptor:retinoid X receptor complexed with 3,3',5 triiodo-L-thyronine and 9-cis retinoic acid  
Authors : Fernandez, E.J.; Putcha, B.-D.K.; Wright, E.; Brunzelle, J.S.  
Deposited on : 2011-11-30  
Resolution : 2.95 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

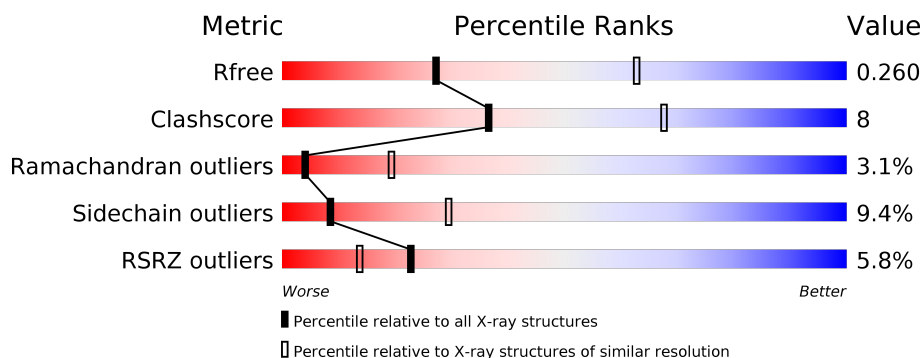
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.95 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	3104 (3.00-2.92)
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)
RSRZ outliers	127900	2986 (3.00-2.92)

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	265	<div> <div>6%</div> <div> <div></div> <div>67%</div> <div>25%</div> <div>•• 6%</div> </div> </div>
2	B	244	<div> <div>5%</div> <div> <div></div> <div>69%</div> <div>18%</div> <div>• 11%</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
3	T3	A	501	-	-	X	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3755 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 Thyroid hormone receptor alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	250	1971	1263	332	360	16	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	144	GLY	-	EXPRESSION TAG	UNP P04625
A	145	SER	-	EXPRESSION TAG	UNP P04625
A	146	HIS	-	EXPRESSION TAG	UNP P04625
A	147	GLU	-	EXPRESSION TAG	UNP P04625

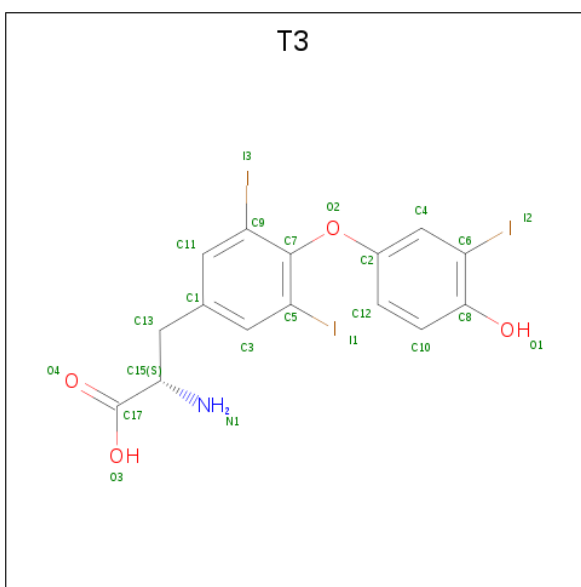
- Molecule 2 is a protein called Retinoic acid receptor RXR-alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	218	1720	1103	295	312	10	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

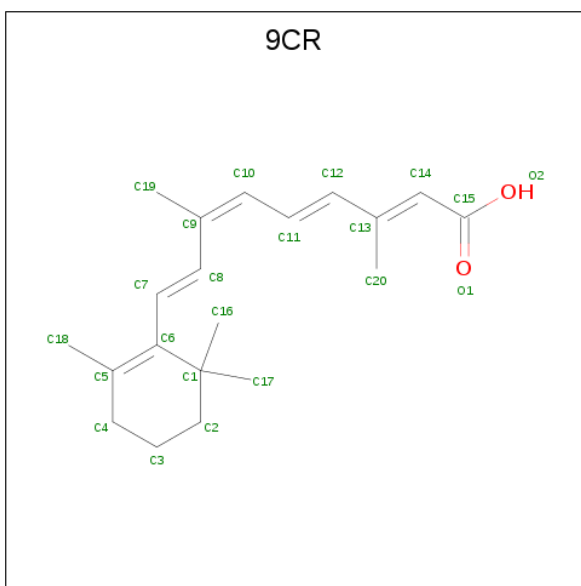
Chain	Residue	Modelled	Actual	Comment	Reference
B	221	MET	-	EXPRESSION TAG	UNP P19793
B	222	LYS	-	EXPRESSION TAG	UNP P19793
B	223	LYS	-	EXPRESSION TAG	UNP P19793
B	224	GLY	-	EXPRESSION TAG	UNP P19793
B	463	MET	-	EXPRESSION TAG	UNP P19793
B	464	THR	-	EXPRESSION TAG	UNP P19793

- Molecule 3 is 3,5,3'-TRIiodothyronine (three-letter code: T3) (formula: C<sub>15</sub>H<sub>12</sub>I<sub>3</sub>NO<sub>4</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	I	N	O	0	0
			23	15	3	1	4		

- Molecule 4 is (9cis)-retinoic acid (three-letter code: 9CR) (formula: C<sub>20</sub>H<sub>28</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			22	20	2		

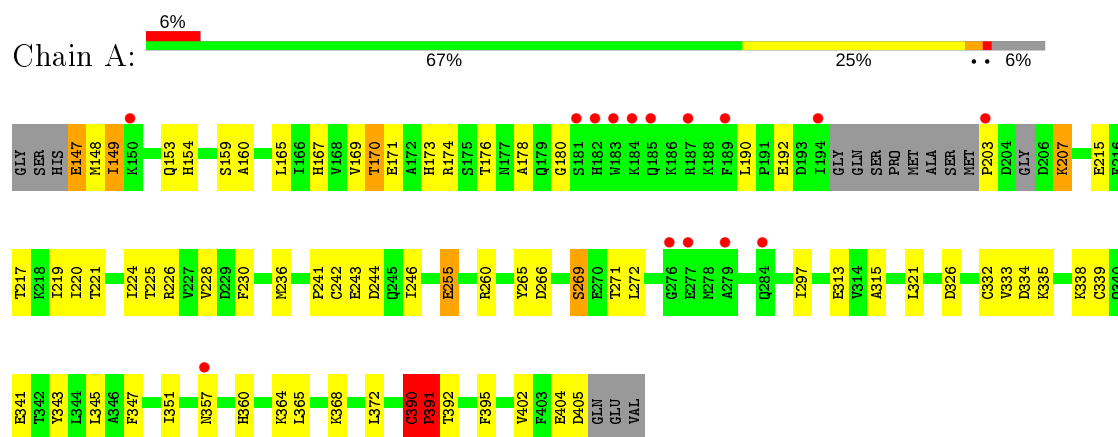
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	11	Total 11	O 11	0	0
5	B	8	Total 8	O 8	0	0

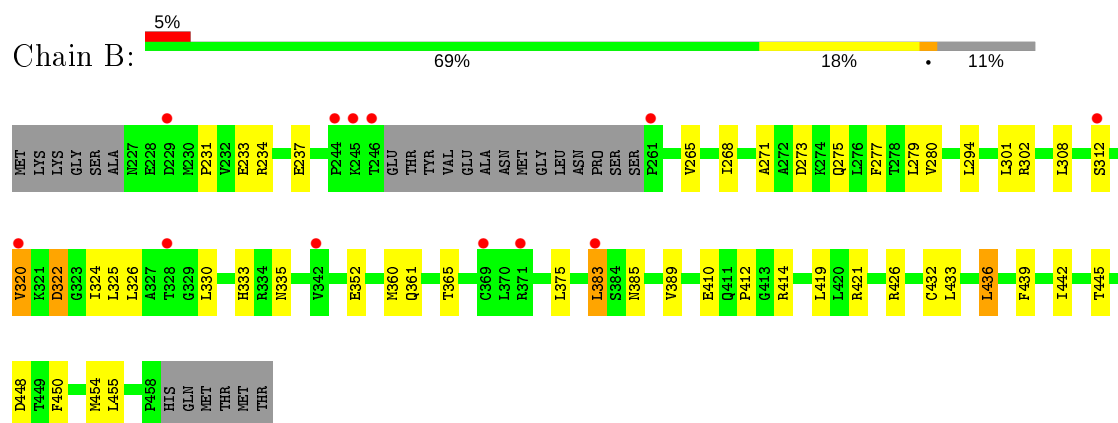
### 3 Residue-property plots [i](#)

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: Thyroid hormone receptor alpha



#### • Molecule 2: Retinoic acid receptor RXR-alpha



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	129.13Å 165.32Å 85.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.55 – 2.95 27.32 – 2.95	Depositor EDS
% Data completeness (in resolution range)	(Not available) (27.55-2.95) 98.9 (27.32-2.95)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.24 (at 2.94Å)	Xtriage
Refinement program	BUSTER 2.8.0	Depositor
R, $R_{free}$	0.201 , 0.250 0.211 , 0.260	Depositor DCC
$R_{free}$ test set	952 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	97.4	Xtriage
Anisotropy	0.634	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 95.7	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	0.96	EDS
Total number of atoms	3755	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	114.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% 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: 9CR, T3

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	0.51	0/2012	0.79	4/2719 (0.1%)
2	B	0.51	0/1755	0.70	0/2374
All	All	0.51	0/3767	0.75	4/5093 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	391	PRO	N-CA-C	7.35	131.20	112.10
1	A	391	PRO	CA-N-CD	-6.35	102.61	111.50
1	A	203	PRO	N-CA-CB	6.05	110.56	103.30
1	A	390	CYS	CA-C-N	5.61	132.82	117.10

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	1971	0	1960	39	0
2	B	1720	0	1755	22	0
3	A	23	0	10	7	0
4	B	22	0	27	6	0
5	A	11	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	8	0	0	0	0
All	All	3755	0	3752	59	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (59) 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:169:VAL:HG21	1:A:343:TYR:HE2	1.30	0.93
2:B:320:VAL:HG21	2:B:325:LEU:HB2	1.63	0.77
1:A:219:ILE:HD11	3:A:501:T3:I1	2.57	0.75
1:A:220:ILE:CD1	3:A:501:T3:I1	3.11	0.69
1:A:176:THR:HG21	1:A:230:PHE:HB2	1.75	0.68
2:B:360:MET:O	2:B:414:ARG:NH2	2.27	0.66
1:A:269:SER:HB2	1:A:271:THR:HG22	1.78	0.66
1:A:220:ILE:HD13	3:A:501:T3:I1	2.67	0.65
1:A:220:ILE:HD12	3:A:501:T3:I1	2.68	0.63
1:A:169:VAL:HG21	1:A:343:TYR:CE2	2.23	0.62
1:A:147:GLU:O	1:A:149:ILE:N	2.32	0.61
1:A:326:ASP:HB2	2:B:352:GLU:HG2	1.82	0.61
1:A:219:ILE:CD1	3:A:501:T3:I1	3.20	0.60
1:A:297:ILE:HD13	3:A:501:T3:I3	2.72	0.59
2:B:432:CYS:SG	4:B:501:9CR:H16	2.43	0.58
2:B:279:LEU:HD11	2:B:308:LEU:HD13	1.87	0.56
1:A:220:ILE:HB	3:A:501:T3:I1	2.76	0.55
1:A:167:HIS:HA	1:A:170:THR:HG22	1.88	0.55
2:B:280:VAL:HG12	2:B:301:LEU:HD13	1.89	0.53
1:A:390:CYS:HB3	1:A:391:PRO:O	2.08	0.53
1:A:241:PRO:HD2	1:A:332:CYS:SG	2.50	0.52
1:A:224:ILE:HG21	1:A:402:VAL:HG21	1.92	0.51
1:A:345:LEU:HD11	2:B:412:PRO:HB2	1.93	0.51
1:A:332:CYS:HB3	1:A:335:LYS:HB2	1.93	0.50
2:B:302:ARG:HA	2:B:454:MET:CE	2.42	0.50
2:B:275:GLN:HG3	4:B:501:9CR:O2	2.12	0.49
1:A:392:THR:HA	1:A:395:PHE:HD1	1.77	0.49
2:B:436:LEU:HD22	2:B:455:LEU:HD11	1.95	0.49
2:B:439:PHE:HA	2:B:442:ILE:HG12	1.95	0.48
1:A:217:THR:HA	1:A:220:ILE:HG22	1.95	0.48
1:A:255:GLU:HB3	1:A:321:LEU:HD13	1.96	0.48
1:A:169:VAL:CG2	1:A:343:TYR:HE2	2.16	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:ILE:O	1:A:228:VAL:HG23	2.15	0.47
2:B:410:GLU:O	2:B:412:PRO:HD3	2.15	0.47
1:A:265:TYR:HB2	1:A:272:LEU:HD23	1.96	0.47
1:A:365:LEU:O	1:A:368:LYS:HB2	2.15	0.46
2:B:322:ASP:HB3	2:B:333:HIS:CE1	2.50	0.46
1:A:236:MET:HB2	1:A:339:CYS:SG	2.55	0.46
2:B:268:ILE:HD13	4:B:501:9CR:H1	1.98	0.46
2:B:265:VAL:HG12	2:B:445:THR:HG21	1.97	0.46
1:A:171:GLU:HA	1:A:174:ARG:HB2	1.99	0.44
1:A:169:VAL:HG12	1:A:315:ALA:HB1	2.00	0.44
1:A:190:LEU:HD11	1:A:215:GLU:HG3	1.99	0.44
1:A:241:PRO:HG2	1:A:244:ASP:HB2	2.00	0.44
2:B:268:ILE:HG23	4:B:501:9CR:H27	2.00	0.44
2:B:383:LEU:HD22	2:B:389:VAL:HG21	1.99	0.44
2:B:273:ASP:CG	2:B:448:ASP:HB3	2.38	0.44
1:A:404:GLU:O	1:A:405:ASP:HB2	2.17	0.43
1:A:321:LEU:HD11	1:A:372:LEU:HD13	2.00	0.43
1:A:176:THR:HG23	1:A:226:ARG:HB3	2.00	0.43
1:A:165:LEU:O	1:A:169:VAL:HG23	2.19	0.42
1:A:338:LYS:HA	1:A:341:GLU:HB2	2.02	0.42
2:B:436:LEU:HD11	4:B:501:9CR:H23	2.01	0.42
1:A:360:HIS:O	1:A:364:LYS:HB2	2.19	0.42
1:A:173:HIS:HD2	1:A:260:ARG:HH21	1.66	0.41
2:B:277:PHE:O	2:B:280:VAL:HG22	2.20	0.41
1:A:176:THR:HG22	1:A:260:ARG:NH2	2.36	0.41
2:B:436:LEU:HG	4:B:501:9CR:C18	2.50	0.41
2:B:271:ALA:HB2	2:B:326:LEU:HD23	2.03	0.41

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	244/265 (92%)	215 (88%)	18 (7%)	11 (4%)	2	12
2	B	214/244 (88%)	197 (92%)	14 (6%)	3 (1%)	11	39
All	All	458/509 (90%)	412 (90%)	32 (7%)	14 (3%)	4	19

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	148	MET
1	A	160	ALA
1	A	192	GLU
1	A	390	CYS
1	A	391	PRO
2	B	320	VAL
1	A	159	SER
1	A	333	VAL
1	A	207	LYS
2	B	322	ASP
1	A	178	ALA
1	A	357	ASN
2	B	385	ASN
1	A	180	GLY

### 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	216/236 (92%)	197 (91%)	19 (9%)	10	33
2	B	188/210 (90%)	169 (90%)	19 (10%)	7	26
All	All	404/446 (91%)	366 (91%)	38 (9%)	8	29

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	147	GLU
1	A	149	ILE

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Mol	Chain	Res	Type
1	A	153	GLN
1	A	154	HIS
1	A	170	THR
1	A	207	LYS
1	A	221	THR
1	A	225	THR
1	A	242	CYS
1	A	243	GLU
1	A	246	ILE
1	A	255	GLU
1	A	266	ASP
1	A	269	SER
1	A	313	GLU
1	A	334	ASP
1	A	347	PHE
1	A	351	ILE
1	A	391	PRO
2	B	231	PRO
2	B	233	GLU
2	B	234	ARG
2	B	237	GLU
2	B	294	LEU
2	B	312	SER
2	B	324	ILE
2	B	330	LEU
2	B	335	ASN
2	B	361	GLN
2	B	365	THR
2	B	375	LEU
2	B	383	LEU
2	B	419	LEU
2	B	421	ARG
2	B	426	ARG
2	B	433	LEU
2	B	436	LEU
2	B	450	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	GLN
1	A	154	HIS

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Mol	Chain	Res	Type
1	A	167	HIS
1	A	173	HIS
1	A	245	GLN
1	A	318	GLN
1	A	352	ASN
2	B	297	GLN
2	B	385	ASN
2	B	411	GLN
2	B	435	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](#)

2 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	$\# Z  > 2$
4	9CR	B	501	-	19,22,22	1.12	2 (10%)	26,30,30	2.28	9 (34%)
3	T3	A	501	-	21,24,24	1.33	3 (14%)	29,34,34	1.06	2 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	9CR	B	501	-	-	5/13/32/32	0/1/1/1
3	T3	A	501	-	-	1/8/12/12	0/2/2/2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	T3	C8-C6	4.42	1.48	1.39
3	A	501	T3	C7-C5	2.52	1.48	1.40
4	B	501	9CR	C12-C13	-2.30	1.41	1.45
3	A	501	T3	C7-C9	2.28	1.47	1.40
4	B	501	9CR	C1-C6	2.17	1.56	1.53

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	501	9CR	C19-C9-C10	-6.18	114.26	122.92
4	B	501	9CR	C16-C1-C6	5.07	118.53	110.30
4	B	501	9CR	C7-C8-C9	-3.15	121.47	126.23
4	B	501	9CR	C19-C9-C8	2.91	122.66	118.08
4	B	501	9CR	C8-C9-C10	2.69	123.06	118.94
4	B	501	9CR	C17-C1-C16	-2.42	101.12	108.53
4	B	501	9CR	C1-C6-C5	-2.31	119.35	122.61
4	B	501	9CR	C2-C3-C4	2.28	116.46	111.38
3	A	501	T3	C3-C5-I1	2.11	122.50	118.61
4	B	501	9CR	C10-C11-C12	-2.09	116.69	123.22
3	A	501	T3	C11-C1-C3	2.00	121.73	118.98

There are no chirality outliers.

All (6) torsion outliers are listed below:

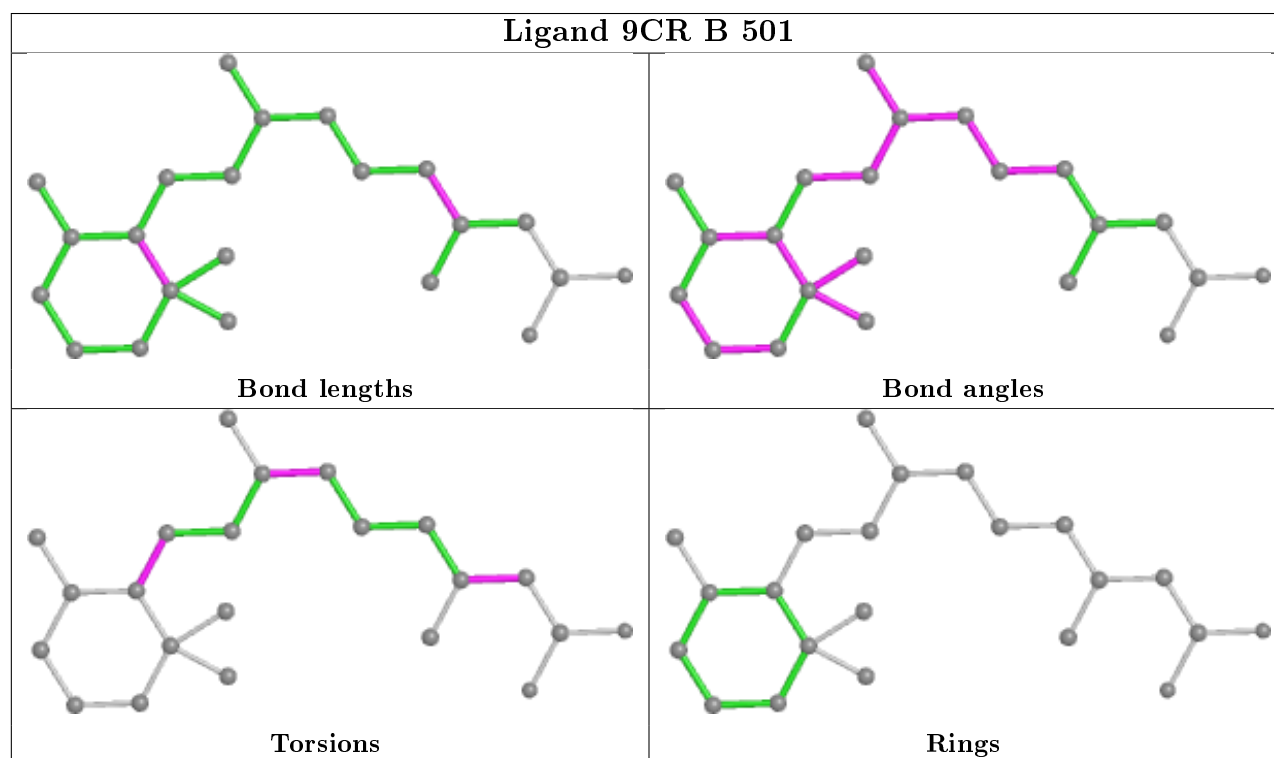
Mol	Chain	Res	Type	Atoms
4	B	501	9CR	C1-C6-C7-C8
4	B	501	9CR	C5-C6-C7-C8
4	B	501	9CR	C12-C13-C14-C15
3	A	501	T3	C1-C13-C15-N1
4	B	501	9CR	C11-C10-C9-C19
4	B	501	9CR	C11-C10-C9-C8

There are no ring outliers.

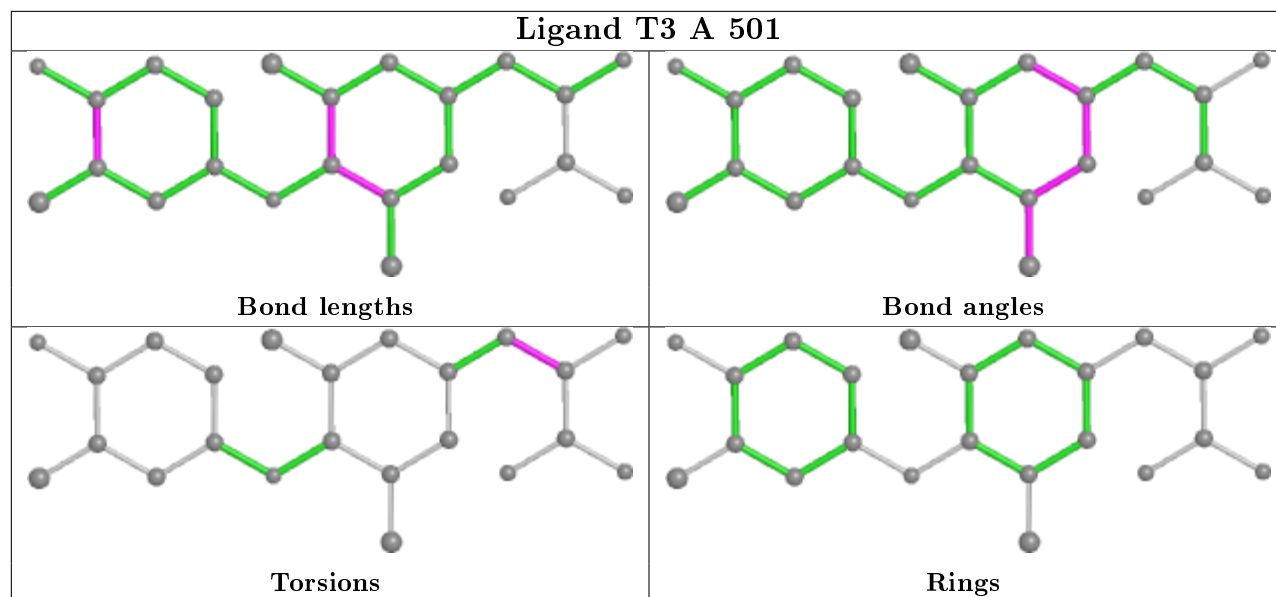
2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	501	9CR	6	0
3	A	501	T3	7	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

### 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	250/265 (94%)	0.28	15 (6%) 21 13	85, 114, 170, 195	8 (3%)
2	B	218/244 (89%)	0.25	12 (5%) 25 15	80, 110, 143, 173	0
All	All	468/509 (91%)	0.27	27 (5%) 23 14	80, 112, 146, 195	8 (1%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	183	TRP	5.8
1	A	182	HIS	5.6
2	B	261	PRO	5.0
1	A	185	GLN	4.5
1	A	181	SER	4.5
2	B	246	THR	4.2
1	A	184	LYS	4.2
2	B	245	LYS	4.0
1	A	187	ARG	3.5
1	A	276	GLY	3.3
2	B	312	SER	3.2
1	A	203	PRO	3.1
1	A	194	ILE	3.0
1	A	279	ALA	2.8
2	B	369	CYS	2.7
2	B	244	PRO	2.7
1	A	357	ASN	2.6
2	B	342	VAL	2.6
1	A	189	PHE	2.5
2	B	383	LEU	2.4
1	A	150	LYS	2.4
1	A	277	GLU	2.3
2	B	320	VAL	2.3
2	B	328	THR	2.2

*Continued on next page...*

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Mol	Chain	Res	Type	RSRZ
1	A	284	GLN	2.2
2	B	229	ASP	2.1
2	B	371	ARG	2.1

## 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](#)

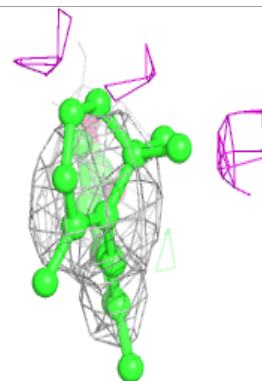
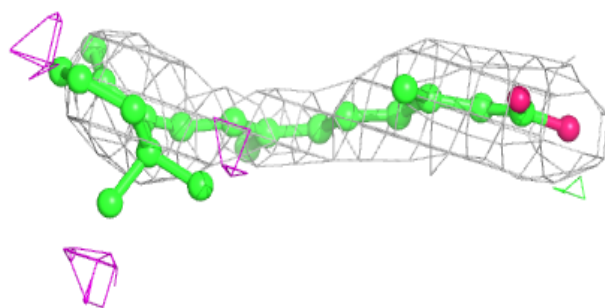
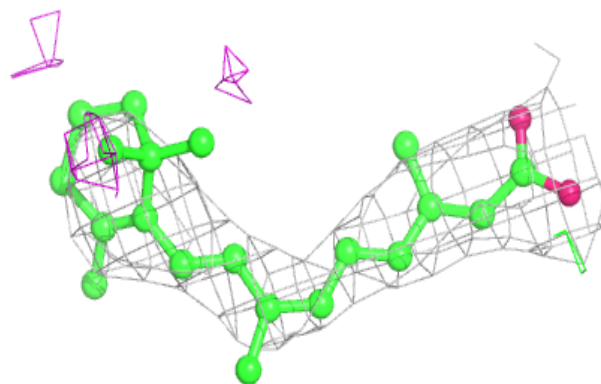
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
4	9CR	B	501	22/22	0.84	0.37	89,99,111,112	0
3	T3	A	501	23/23	0.97	0.35	88,98,109,121	23

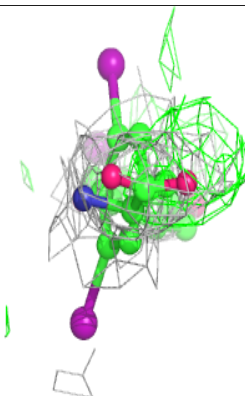
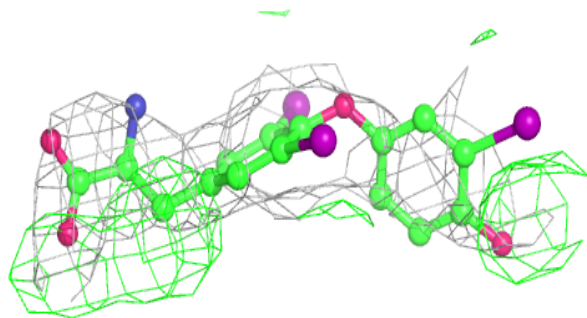
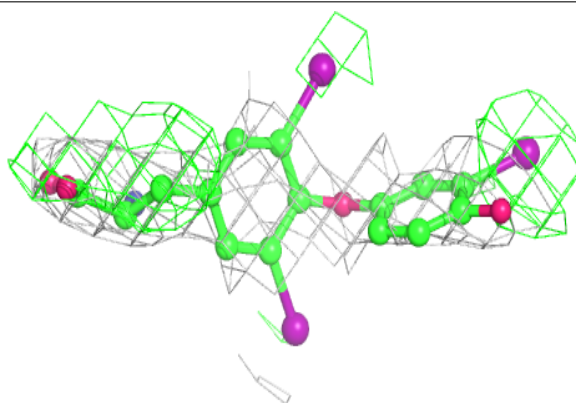
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around 9CR B 501:**

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

**Electron density around T3 A 501:**

$2mF_o-DF_c$  (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.