

wwPDB X-ray Structure Validation Summary Report (i)

Oct 10, 2023 – 01:50 AM EDT

PDB ID : 7TDL

Title : M379A mutant tyrosine phenol-lyase complexed with 3-bromo-DL-phenylala

nine

Authors : Phillips, R.S. Deposited on : 2022-01-01

Resolution : 1.60 Å(reported)

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

We welcome your comments at 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 (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) 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 : 2.35.1

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

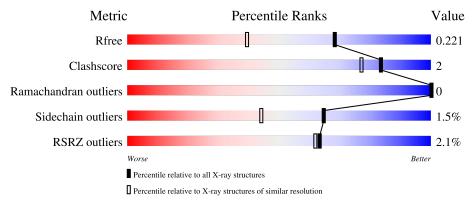
Validation Pipeline (wwPDB-VP) : 2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 1.60 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
R_{free}	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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 >=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 <=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	456	91%	9%
1	В	456	93%	7%
1	С	456	94%	6%
1	D	456	96%	



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 31090 atoms, of which 14679 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 Tyrosine phenol-lyase.

Mol	Chain	Residues		Atoms						AltConf	Trace
1	A	456	Total	С	Н	N	О	S	0	10	0
1	A	450	7289	2321	3628	633	680	27	0	10	
1	В	456	Total	С	Н	N	О	S	0	12	0
1	Ъ	450	7306	2323	3636	639	681	27			
1	С	456	Total	С	Н	N	О	S	0	15	0
1		450	7358	2340	3667	640	684	27		10	
1	1 D	456	Total	С	Н	N	О	S	0	14	0
1			7321	2330	3640	638	684	29	U	14	

There are 8 discrepancies between the modelled and reference sequences:

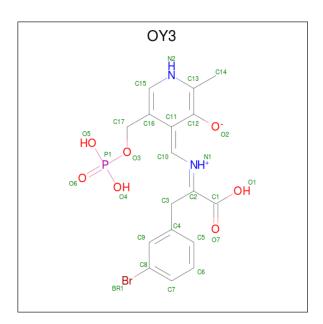
Chain	Residue	Modelled	Actual	Comment	Reference
A	205	ALA	GLU	$\operatorname{conflict}$	UNP P31013
A	379	ALA	MET	engineered mutation	UNP P31013
В	205	ALA	GLU	$\operatorname{conflict}$	UNP P31013
В	379	ALA	MET	engineered mutation	UNP P31013
С	205	ALA	GLU	$\operatorname{conflict}$	UNP P31013
С	379	ALA	MET	engineered mutation	UNP P31013
D	205	ALA	GLU	conflict	UNP P31013
D	379	ALA	MET	engineered mutation	UNP P31013

• Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total K 2 2	0	0
2	С	2	Total K 2 2	0	0

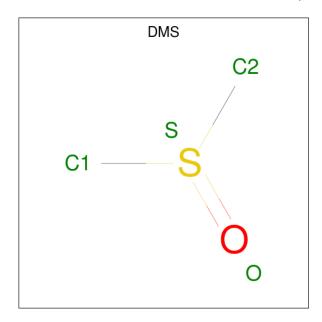
• Molecule 3 is (4Z)-4-($\{[(1E)-2-(3-bromophenyl)-1-carboxyethylidene]$ azaniumyl $\}$ methylidene)-2-methyl-5-[(phosphonooxy)methyl]-1,4-dihydropyridin-3-olate (three-letter code: OY3) (formula: $C_{17}H_{18}BrN_2O_7P$) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
9	9 Л	1	Total	Br	С	Н	N	О	Р	0	1
$\begin{array}{ c c c c } \hline 3 & A & \end{array}$	A		86	2	34	30	4	14	2	U	1
9	D	1	Total	Br	С	Н	N	О	Р	0	1
3 B	Б	1	86	2	34	30	4	14	2	U	1

 \bullet Molecule 4 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: $\mathrm{C_2H_6OS}).$



N	/Iol	Chain	Residues	Atoms					ZeroOcc	AltConf	
	4	Δ	1	Total	С	Н	О	S	0	0	
	T 1	11	1	10	2	6	1	1			
	1	0	\sim	1	Total	С	Η	О	S	0	0
	4 C	1	10	2	6	1	1	U	0		

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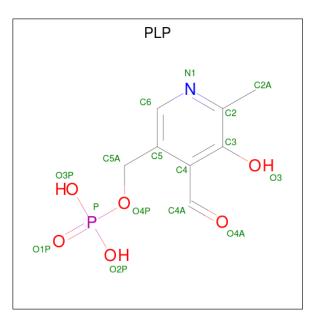
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	С	1	Total	С	Н	О	S	0	0
			10	2	6	1	1		
1	D	1	Total	С	Η	О	S	0	0
4 D	1	10	2	6	1	1	0		

• Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Cl 1 1	0	0
5	С	1	Total Cl 1 1	0	0
5	D	1	Total Cl 1 1	0	0

 $\bullet \ \ Molecule\ 6\ is\ PYRIDOXAL-5'-PHOSPHATE\ (three-letter\ code:\ PLP)\ (formula:\ C_8H_{10}NO_6P).$



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	В	1	Total C H N O P 23 8 8 1 5 1	0	1
6	С	1	Total C H N O P 23 8 8 1 5 1	0	0
6	D	1	Total C H N O P 23 8 8 1 5 1	0	0

• Molecule 7 is water.



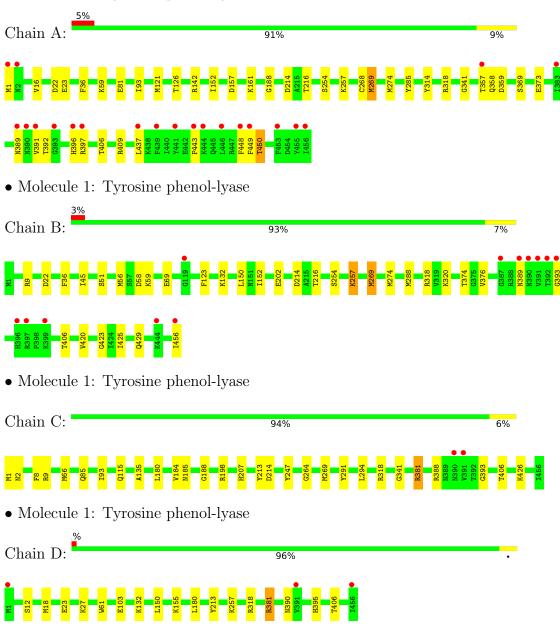
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	341	Total O 341 341	0	27
7	В	286	Total O 286 286	0	10
7	С	465	Total O 465 465	0	29
7	D	436	Total O 436 436	0	25



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Tyrosine phenol-lyase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants	61.25Å 82.87Å 94.33Å	Donositor
a, b, c, α , β , γ	113.16° 96.54° 102.21°	Depositor
Resolution (Å)	58.41 - 1.60	Depositor
Resolution (A)	58.41 - 1.60	EDS
% Data completeness	98.7 (58.41-1.60)	Depositor
(in resolution range)	98.6 (58.41-1.60)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.04 (at 1.60Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
D D.	0.177 , 0.221	Depositor
R, R_{free}	0.178 , 0.221	DCC
R_{free} test set	1985 reflections (0.94%)	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.38, 39.4	EDS
L-test for twinning ²	$ < L >=0.51, < L^2>=0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	31090	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.15% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

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: PLP, OY3, DMS, K, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with |Z| > 5 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Во	nd lengths	Bond angles		
IVIOI	Moi Chain		# Z > 5	RMSZ	# Z >5	
1	A	0.69	1/3767~(0.0%)	0.75	0/5073	
1	В	0.61	$1/3776 \ (0.0\%)$	0.71	1/5081 (0.0%)	
1	С	0.75	1/3809 (0.0%)	0.79	0/5126	
1	D	0.74	0/3793	0.78	0/5104	
All	All	0.70	3/15145 (0.0%)	0.76	1/20384 (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 maintenain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	С	0	1
1	D	0	2
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	$\operatorname{Ideal}(\text{\AA})$
1	A	268	CYS	CB-SG	-6.96	1.70	1.82
1	В	69	GLU	CB-CG	6.52	1.64	1.52
1	С	291	TYR	CE1-CZ	5.56	1.45	1.38

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	В	58	ASP	CB-CG-OD1	6.96	124.56	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	С	381	ARG	Sidechain
1	D	381	ARG	Sidechain
1	D	395	HIS	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	3661	3628	3630	32	0
1	В	3670	3636	3634	20	0
1	С	3691	3667	3659	17	0
1	D	3681	3640	3634	9	0
2	A	2	0	0	0	0
2	С	2	0	0	0	0
3	A	56	30	0	6	0
3	В	56	30	0	1	0
4	A	4	6	6	0	0
4	С	8	12	12	0	0
4	D	4	6	6	1	0
5	A	1	0	0	0	0
5	С	1	0	0	0	0
5	D	1	0	0	0	0
6	В	15	8	7	1	0
6	С	15	8	7	1	0
6	D	15	8	7	2	0
7	A	341	0	0	6	0
7	В	286	0	0	2	0
7	С	465	0	0	7	0
7	D	436	0	0	1	0
All	All	16411	14679	14602	72	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

The worst 5 of 72 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.



Atom-1	Atom-2	$egin{array}{ll} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:A:391:VAL:HG13	1:A:392:THR:HG23	1.66	0.77
1:A:214:ASP:OD1	1:A:216[B]:THR:HG23	1.88	0.74
1:B:123:PHE:HB2	6:B:502[A]:PLP:H2A3	1.82	0.61
1:C:184:VAL:HG13	7:C:840:HOH:O	2.04	0.57
1:A:16[A]:VAL:HG11	1:C:9:ARG:HG3	1.86	0.56

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	Perce	ntiles
1	A	$464/456 \; (102\%)$	451 (97%)	13 (3%)	0	100	100
1	В	466/456 (102%)	454 (97%)	12 (3%)	0	100	100
1	С	469/456 (103%)	458 (98%)	11 (2%)	0	100	100
1	D	468/456 (103%)	457 (98%)	11 (2%)	0	100	100
All	All	1867/1824 (102%)	1820 (98%)	47 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	$387/377 \ (103\%)$	380 (98%)	7 (2%)	59 36	
1	В	$387/377 \ (103\%)$	378 (98%)	9 (2%)	50 25	

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles		
1	С	391/377 (104%)	387 (99%)	4 (1%)	76 61		
1	D	389/377 (103%)	385 (99%)	4 (1%)	76 61		
All	All	1554/1508 (103%)	1530 (98%)	24 (2%)	65 44		

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	318	ARG
1	С	318	ARG
1	С	269	MET
1	С	381	ARG
1	A	450	THR

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

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)

Of 18 ligands modelled in this entry, 7 are monoatomic - leaving 11 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	Trino	Chain	Res	Link	Во	ond leng	ths	В	ond ang	les
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
4	DMS	С	503	-	3,3,3	0.42	0	3,3,3	0.80	0
4	DMS	С	504	-	3,3,3	0.75	0	3,3,3	0.85	0
3	OY3	A	501[B]	-	27,29,29	1.21	4 (14%)	33,41,41	1.24	4 (12%)
3	OY3	В	501[B]	-	27,29,29	1.46	4 (14%)	33,41,41	1.21	4 (12%)
6	PLP	С	502	1	15,15,16	2.01	4 (26%)	20,22,23	1.42	3 (15%)
4	DMS	A	504	-	3,3,3	0.80	0	3,3,3	1.08	0
6	PLP	В	502[A]	1	15,15,16	2.24	5 (33%)	20,22,23	1.31	2 (10%)
3	OY3	A	501[A]	-	27,29,29	1.39	4 (14%)	33,41,41	1.17	4 (12%)
4	DMS	D	501	-	3,3,3	0.41	0	3,3,3	1.69	1 (33%)
6	PLP	D	502	1	15,15,16	2.22	5 (33%)	20,22,23	1.16	1 (5%)
3	OY3	В	501[C]	-	27,29,29	1.50	4 (14%)	33,41,41	1.21	5 (15%)

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	OY3	A	501[B]	-	-	1/16/19/19	0/2/2/2
3	OY3	В	501[B]	-	-	1/16/19/19	0/2/2/2
6	PLP	С	502	1	-	0/6/6/8	0/1/1/1
6	PLP	В	502[A]	1	-	5/6/6/8	0/1/1/1
3	OY3	A	501[A]	-	-	1/16/19/19	0/2/2/2
6	PLP	D	502	1	-	0/6/6/8	0/1/1/1
3	OY3	В	501[C]	-	-	1/16/19/19	0/2/2/2

The worst 5 of 30 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	$\operatorname{Ideal}(\text{\AA})$
3	В	501[C]	OY3	O3-C17	5.17	1.46	1.43
3	В	501[B]	OY3	O3-C17	5.03	1.46	1.43
6	В	502[A]	PLP	C4A-C4	4.93	1.61	1.51
6	С	502	PLP	C4A-C4	4.69	1.61	1.51
3	A	501[A]	OY3	O3-C17	4.50	1.46	1.43

The worst 5 of 24 bond angle outliers are listed below:

N	Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\mathrm{Ideal}(^{o})$
	6	С	502	PLP	O4P-C5A-C5	3.46	115.95	109.35
	3	A	501[A]	OY3	C3-C2-N1	-3.23	117.77	125.81

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Mol	Chain	Res	Type	$\Gamma ext{ype} \qquad ext{Atoms}$		$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
3	В	501[B]	OY3	C3-C2-C1	3.16	123.69	118.28
3	A	501[A]	OY3	C3-C2-C1	3.09	123.56	118.28
3	В	501[B]	OY3	C3-C2-N1	-3.05	118.22	125.81

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501[A]	OY3	N1-C10-C11-C16
3	A	501[B]	OY3	N1-C10-C11-C16
3	В	501[B]	OY3	N1-C10-C11-C16
3	В	501[C]	OY3	N1-C10-C11-C16
6	В	502[A]	PLP	C4-C5-C5A-O4P

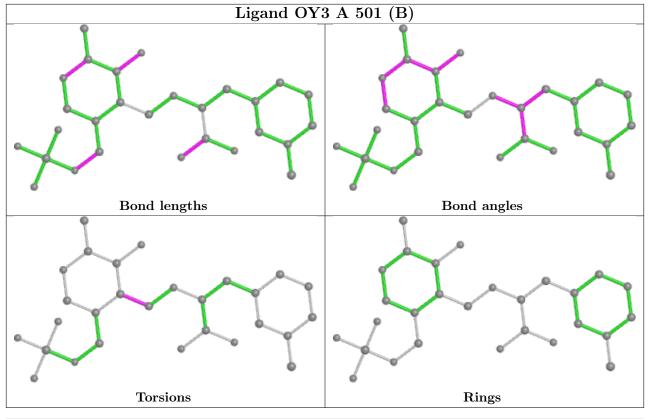
There are no ring outliers.

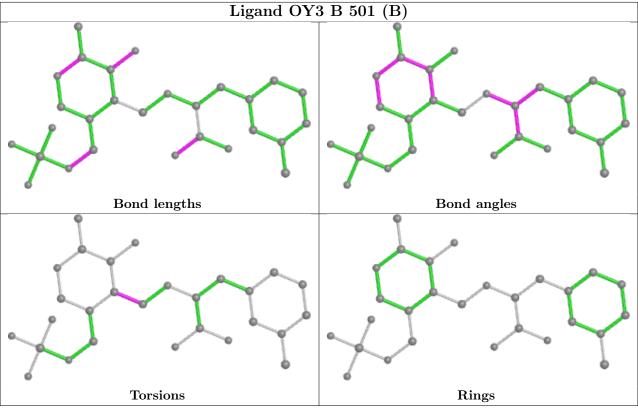
7 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501[B]	OY3	2	0
3	В	501[B]	OY3	1	0
6	С	502	PLP	1	0
6	В	502[A]	PLP	1	0
3	A	501[A]	OY3	4	0
4	D	501	DMS	1	0
6	D	502	PLP	2	0

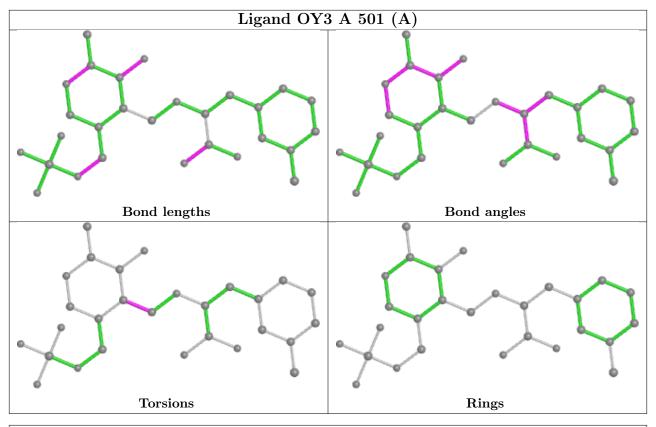
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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.

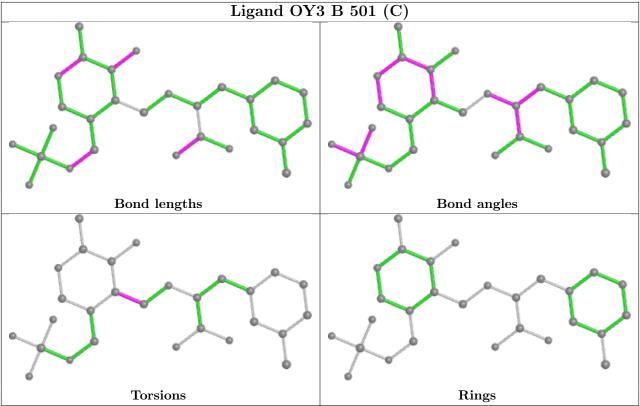














5.7 Other polymers (i)

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^{th} 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>	# RSRZ > 2	$OWAB(Å^2)$	Q<0.9
1	A	456/456 (100%)	-0.02	21 (4%) 32 29	19, 32, 67, 107	0
1	В	$456/456 \; (100\%)$	0.05	12 (2%) 56 53	21, 38, 68, 130	0
1	С	$456/456 \; (100\%)$	-0.32	2 (0%) 92 92	18, 27, 42, 88	0
1	D	$456/456 \; (100\%)$	-0.36	3 (0%) 87 87	19, 30, 47, 103	0
All	All	1824/1824 (100%)	-0.16	38 (2%) 63 62	18, 30, 61, 130	0

The worst 5 of 38 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	391	VAL	9.2
1	В	390	ASN	8.9
1	D	1	MET	8.0
1	A	391	VAL	5.2
1	A	456	ILE	5.2

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)

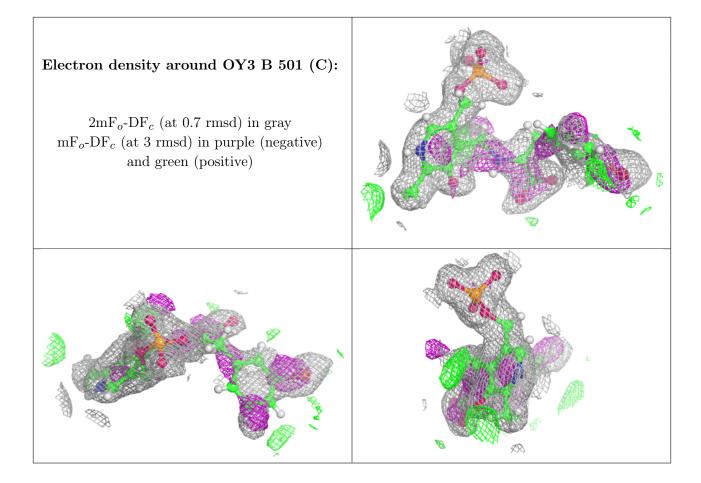
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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



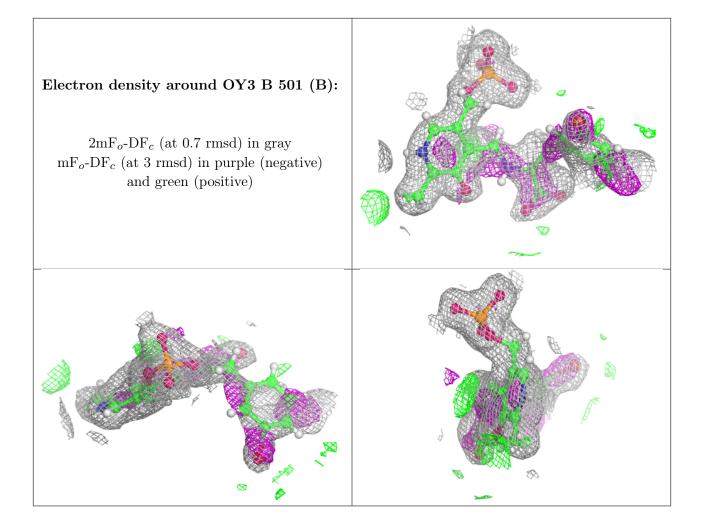
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
4	DMS	A	504	4/4	0.88	0.19	41,73,76,76	0
4	DMS	D	501	4/4	0.90	0.14	25,30,37,41	0
4	DMS	С	504	4/4	0.93	0.13	33,40,51,52	0
3	OY3	В	501[C]	28/28	0.94	0.18	28,42,56,77	43
3	OY3	В	501[B]	28/28	0.94	0.18	28,42,55,61	43
4	DMS	С	503	4/4	0.95	0.08	27,33,35,35	0
3	OY3	A	501[B]	28/28	0.95	0.11	24,35,50,56	43
3	OY3	A	501[A]	28/28	0.95	0.11	24,35,50,58	43
6	PLP	В	502[A]	15/16	0.97	0.10	28,36,44,44	23
6	PLP	С	502	15/16	0.97	0.09	21,24,30,32	0
5	CL	D	503	1/1	0.98	0.04	34,34,34,34	0
6	PLP	D	502	15/16	0.98	0.08	23,26,31,36	0
5	CL	A	505	1/1	0.99	0.03	41,41,41,41	0
5	CL	С	501	1/1	0.99	0.04	34,34,34,34	0
2	K	С	505	1/1	1.00	0.05	20,20,20,20	0
2	K	С	506	1/1	1.00	0.06	21,21,21,21	0
2	K	A	502	1/1	1.00	0.07	22,22,22,22	0
2	K	A	503	1/1	1.00	0.06	22,22,22,22	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.





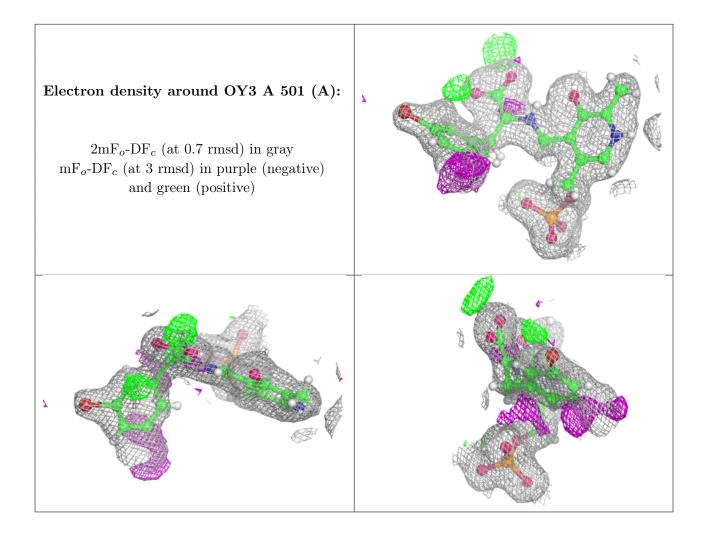






Electron density around OY3 A 501 (B): 2mF_o-DF_c (at 0.7 rmsd) in gray mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)





6.5 Other polymers (i)

There are no such residues in this entry.

