



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:20 AM UTC

PDB ID : 6U3J / pdb_00006u3j
Title : Structure of the 2-oxoadipate dehydrogenase DHTKD1
Authors : Khamrui, S.; Lazarus, M.B.
Deposited on : 2019-08-21
Resolution : 2.25 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

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>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

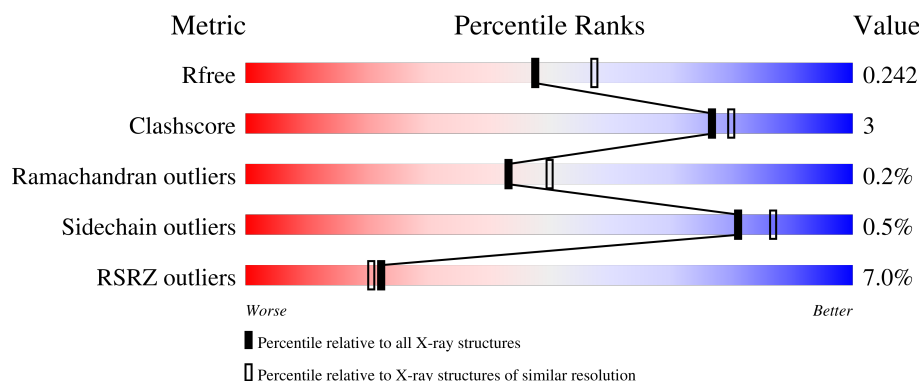
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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}	180053	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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 $\leq 5\%$. 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	904	
1	B	904	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14773 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.

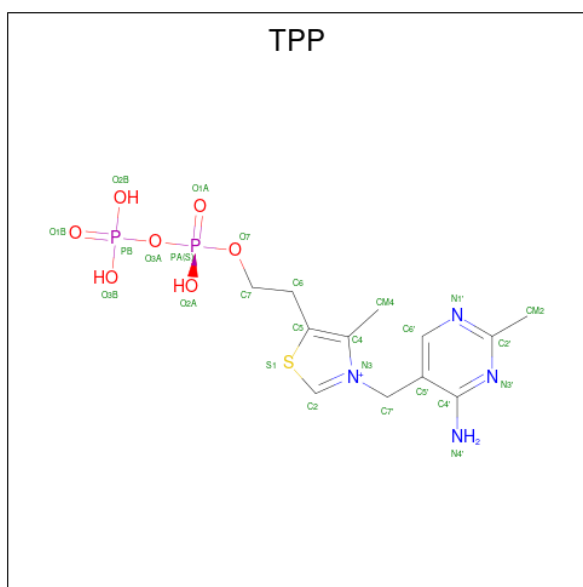
- Molecule 1 is a protein called 2-oxoglutarate dehydrogenase E1 component DHKTD1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	876	Total	C	N	O	S	0	1	0
			6878	4381	1196	1261	40			
1	B	876	Total	C	N	O	S	0	1	0
			6878	4381	1196	1261	40			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MET	-	initiating methionine	UNP Q96HY7
A	24	GLY	-	expression tag	UNP Q96HY7
A	920	HIS	-	expression tag	UNP Q96HY7
A	921	HIS	-	expression tag	UNP Q96HY7
A	922	HIS	-	expression tag	UNP Q96HY7
A	923	HIS	-	expression tag	UNP Q96HY7
A	924	HIS	-	expression tag	UNP Q96HY7
A	925	HIS	-	expression tag	UNP Q96HY7
A	926	HIS	-	expression tag	UNP Q96HY7
B	23	MET	-	initiating methionine	UNP Q96HY7
B	24	GLY	-	expression tag	UNP Q96HY7
B	920	HIS	-	expression tag	UNP Q96HY7
B	921	HIS	-	expression tag	UNP Q96HY7
B	922	HIS	-	expression tag	UNP Q96HY7
B	923	HIS	-	expression tag	UNP Q96HY7
B	924	HIS	-	expression tag	UNP Q96HY7
B	925	HIS	-	expression tag	UNP Q96HY7
B	926	HIS	-	expression tag	UNP Q96HY7

- Molecule 2 is THIAMINE DIPHOSPHATE (CCD ID: TPP) (formula: $C_{12}H_{19}N_4O_7P_2S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total 26	C 12	N 4	O 7	P 2	S 1	0	0
2	B	1	Total 26	C 12	N 4	O 7	P 2	S 1	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

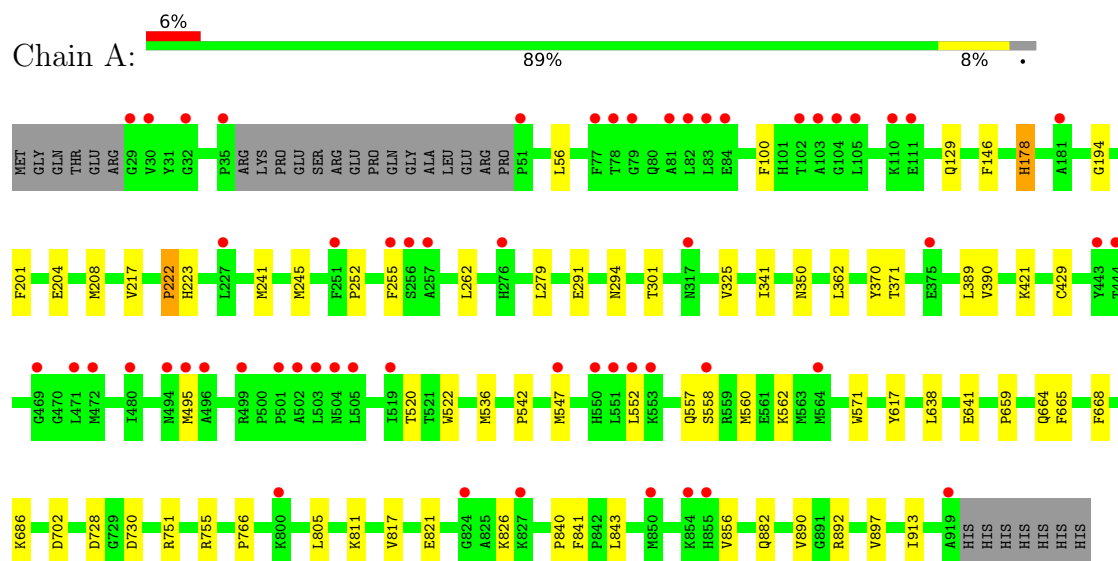
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	472	Total O 472 472	0	0
4	B	490	Total O 490 490	0	0

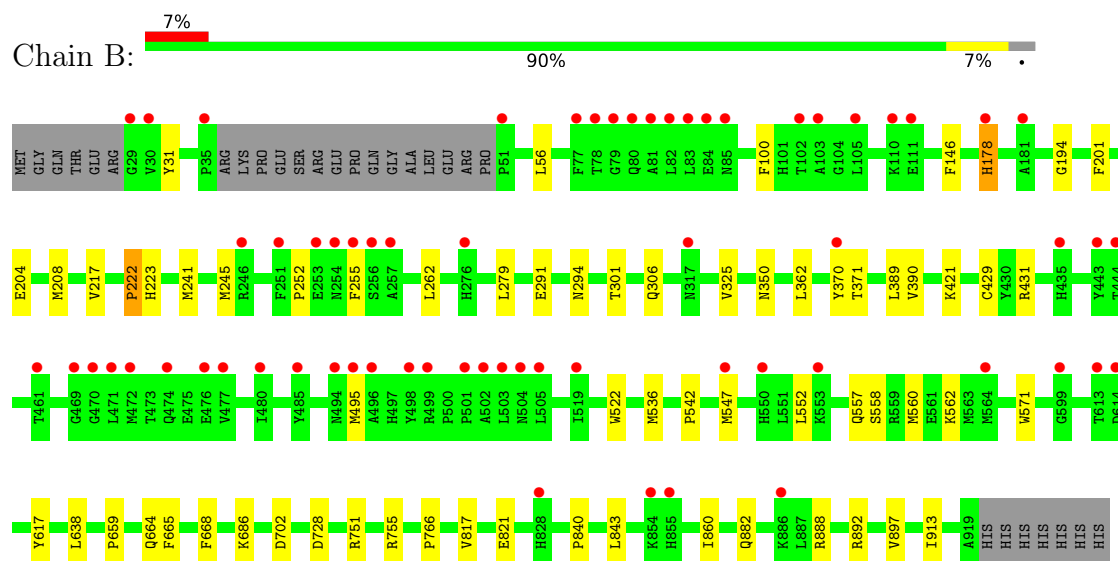
3 Residue-property plots [i](#)

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: 2-oxoglutarate dehydrogenase E1 component DHKTD1, mitochondrial



- Molecule 1: 2-oxoglutarate dehydrogenase E1 component DHKTD1, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	332.31Å 72.61Å 79.67Å 90.00° 91.87° 90.00°	Depositor
Resolution (Å)	49.01 – 2.25 49.01 – 2.25	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.01-2.25) 99.9 (49.01-2.25)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.209 , 0.241 0.211 , 0.242	Depositor DCC
R_{free} test set	4591 reflections (4.12%)	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.069 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14773	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MLZ, MG, TPP

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	0.10	0/7029	0.29	0/9525
1	B	0.10	0/7029	0.29	0/9525
All	All	0.10	0/14058	0.29	0/19050

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6878	0	6756	40	0
1	B	6878	0	6756	37	0
2	A	26	0	16	2	0
2	B	26	0	16	2	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
4	A	472	0	0	2	0
4	B	490	0	0	2	0
All	All	14773	0	13544	79	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 3.

All (79) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:GLU:OE1	1:B:294:ASN:ND2	2.22	0.72
1:A:291:GLU:OE1	1:A:294:ASN:ND2	2.22	0.71
1:A:389:LEU:HD12	1:A:390:VAL:HG13	1.75	0.68
1:B:389:LEU:HD12	1:B:390:VAL:HG13	1.75	0.66
1:B:241:MET:HG2	1:B:245:MET:HE3	1.85	0.58
1:B:178:HIS:C	1:B:178:HIS:HD1	2.12	0.58
1:A:129:GLN:NE2	4:A:1111:HOH:O	2.37	0.57
1:A:241:MET:HG2	1:A:245:MET:HE3	1.85	0.56
1:A:350:ASN:HD21	1:A:421:LYS:HE3	1.72	0.54
1:B:350:ASN:HD21	1:B:421:LYS:HE3	1.72	0.54
1:B:431:ARG:NH1	4:B:1117:HOH:O	2.39	0.51
1:A:204:GLU:HG3	1:A:208:MET:HE2	1.93	0.51
1:B:552:LEU:O	1:B:557:GLN:HG2	2.11	0.51
1:B:204:GLU:HG3	1:B:208:MET:HE2	1.93	0.51
1:A:552:LEU:O	1:A:557:GLN:HG2	2.11	0.50
1:A:558:SER:O	1:A:562:LYS:HG2	2.11	0.50
1:B:558:SER:O	1:B:562:LYS:HG2	2.11	0.50
1:B:547:MET:HG2	1:B:552:LEU:HG	1.94	0.49
1:A:817:VAL:O	1:A:821:GLU:HG2	2.12	0.49
1:B:571:TRP:CE2	1:B:766:PRO:HG3	2.48	0.49
1:A:547:MET:HG2	1:A:552:LEU:HG	1.94	0.49
1:B:817:VAL:O	1:B:821:GLU:HG2	2.12	0.49
1:A:571:TRP:CE2	1:A:766:PRO:HG3	2.48	0.49
1:B:536:MET:HE1	1:B:560:MET:HA	1.95	0.48
1:A:536:MET:HE1	1:A:560:MET:HA	1.95	0.48
1:B:702:ASP:OD2	4:B:1101:HOH:O	2.19	0.48
1:B:178:HIS:C	1:B:178:HIS:ND1	2.71	0.47
1:A:686:LYS:HG3	1:B:897:VAL:HG11	1.97	0.46
1:A:751:ARG:O	1:A:755:ARG:HG3	2.16	0.46
1:B:751:ARG:O	1:B:755:ARG:HG3	2.16	0.46
1:B:892:ARG:HG3	1:B:913:ILE:HD11	1.98	0.45
1:B:843:LEU:HD13	1:B:882:GLN:HB3	1.98	0.45
2:A:1001:TPP:HN42	2:A:1001:TPP:H2	1.81	0.45
1:A:194:GLY:O	1:A:429:CYS:HB2	2.17	0.45
1:A:178:HIS:HD1	1:A:178:HIS:C	2.25	0.45
1:A:665:PHE:HB2	1:A:668:PHE:CE2	2.52	0.45
1:B:665:PHE:HB2	1:B:668:PHE:CE2	2.52	0.45
1:A:843:LEU:HD13	1:A:882:GLN:HB3	1.98	0.44
1:B:194:GLY:O	1:B:429:CYS:HB2	2.17	0.44
1:A:897:VAL:HG11	1:B:686:LYS:HG3	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1001:TPP:HN42	2:B:1001:TPP:H2	1.81	0.44
2:B:1001:TPP:HN42	2:B:1001:TPP:C2	2.31	0.44
1:A:702:ASP:OD1	1:A:811:LYS:NZ	2.50	0.44
1:A:892:ARG:HG3	1:A:913:ILE:HD11	1.99	0.44
2:A:1001:TPP:HN42	2:A:1001:TPP:C2	2.31	0.44
1:A:728:ASP:HB3	1:A:840:PRO:HB3	2.00	0.43
1:A:522:TRP:CE2	1:A:755:ARG:HG2	2.54	0.43
1:B:522:TRP:CE2	1:B:755:ARG:HG2	2.54	0.43
1:B:370:TYR:HB3	1:B:371:THR:H	1.68	0.43
1:B:728:ASP:HB3	1:B:840:PRO:HB3	2.00	0.43
1:B:56:LEU:HD22	1:B:100:PHE:HB3	2.01	0.42
1:B:217:VAL:HG23	1:B:279:LEU:HD11	2.02	0.42
1:B:245:MET:HE2	1:B:262:LEU:HD21	2.01	0.42
1:A:495:MET:HE2	1:A:495:MET:HB3	1.93	0.42
1:B:495:MET:HE2	1:B:495:MET:HB3	1.93	0.42
1:A:542:PRO:HD3	1:A:617:TYR:CE2	2.54	0.42
1:B:542:PRO:HD3	1:B:617:TYR:CE2	2.55	0.42
1:B:252:PRO:HD2	1:B:255:PHE:CD2	2.56	0.41
1:B:201:PHE:CD2	1:B:362:LEU:HD22	2.55	0.41
1:A:217:VAL:HG23	1:A:279:LEU:HD11	2.02	0.41
1:A:341:ILE:HD12	1:A:641:GLU:HG2	2.02	0.41
1:A:841:PHE:O	4:A:1101:HOH:O	2.21	0.41
1:B:222:PRO:HB2	1:B:223:HIS:H	1.69	0.41
1:A:201:PHE:CD2	1:A:362:LEU:HD22	2.55	0.41
1:A:245:MET:HE2	1:A:262:LEU:HD21	2.02	0.41
1:A:520:THR:OG1	1:A:730:ASP:OD2	2.38	0.41
1:A:826:LYS:HB3	1:A:826:LYS:HE2	1.87	0.41
1:A:890:VAL:HG12	1:A:913:ILE:HG23	2.03	0.41
1:B:301:THR:HG23	1:B:325:VAL:HG12	2.03	0.41
1:A:222:PRO:HB2	1:A:223:HIS:H	1.69	0.40
1:A:301:THR:HG23	1:A:325:VAL:HG12	2.03	0.40
1:A:56:LEU:HD22	1:A:100:PHE:HB3	2.01	0.40
1:A:252:PRO:HD2	1:A:255:PHE:CD2	2.56	0.40
1:A:370:TYR:HB3	1:A:371:THR:H	1.69	0.40
1:A:178:HIS:C	1:A:178:HIS:ND1	2.76	0.40
1:A:805:LEU:HG	1:A:856:VAL:HG11	2.02	0.40
1:B:522:TRP:CD1	1:B:755:ARG:HG2	2.57	0.40
1:B:31:TYR:OH	1:B:306:GLN:NE2	2.49	0.40
1:B:860:ILE:HD11	1:B:888:ARG:NH2	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	871/904 (96%)	847 (97%)	22 (2%)	2 (0%)	43	50
1	B	871/904 (96%)	845 (97%)	24 (3%)	2 (0%)	43	50
All	All	1742/1808 (96%)	1692 (97%)	46 (3%)	4 (0%)	43	50

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	222	PRO
1	B	222	PRO
1	A	659	PRO
1	B	659	PRO

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	742/779 (95%)	738 (100%)	4 (0%)	81	87
1	B	742/779 (95%)	738 (100%)	4 (0%)	81	87
All	All	1484/1558 (95%)	1476 (100%)	8 (0%)	81	87

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	146	PHE
1	A	178	HIS

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Mol	Chain	Res	Type
1	A	638	LEU
1	A	664	GLN
1	B	146	PHE
1	B	178	HIS
1	B	638	LEU
1	B	664	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	HIS
1	A	129	GLN
1	A	289	HIS
1	A	350	ASN
1	A	445	ASN
1	A	465	HIS
1	A	489	ASN
1	A	555	HIS
1	A	605	HIS
1	A	752	GLN
1	B	66	HIS
1	B	129	GLN
1	B	140	GLN
1	B	289	HIS
1	B	306	GLN
1	B	350	ASN
1	B	365	ASN
1	B	445	ASN
1	B	465	HIS
1	B	489	ASN
1	B	555	HIS
1	B	605	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	MLZ	A	72	1	8,9,10	0.95	0	4,9,11	0.46	0
1	MLZ	A	537	1	8,9,10	0.84	0	4,9,11	0.38	0
1	MLZ	B	72	1	8,9,10	0.96	0	4,9,11	0.47	0
1	MLZ	B	537	1	8,9,10	0.83	0	4,9,11	0.37	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	MLZ	A	72	1	-	1/7/8/10	-
1	MLZ	A	537	1	-	1/7/8/10	-
1	MLZ	B	72	1	-	1/7/8/10	-
1	MLZ	B	537	1	-	1/7/8/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	72	MLZ	CG-CD-CE-NZ
1	B	72	MLZ	CG-CD-CE-NZ
1	A	537	MLZ	C-CA-CB-CG
1	B	537	MLZ	C-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	TPP	B	1001	3	26,27,27	1.82	6 (23%)	38,40,40	1.63	10 (26%)
2	TPP	A	1001	3	26,27,27	1.83	7 (26%)	38,40,40	1.60	9 (23%)

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	TPP	B	1001	3	-	4/17/17/17	0/2/2/2
2	TPP	A	1001	3	-	3/17/17/17	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	TPP	C5-S1	-5.10	1.59	1.72
2	A	1001	TPP	C5-S1	-5.10	1.59	1.72
2	A	1001	TPP	C2-S1	-2.99	1.60	1.69
2	B	1001	TPP	C2-S1	-2.98	1.60	1.69
2	A	1001	TPP	C2'-N1'	2.71	1.38	1.34
2	B	1001	TPP	C2'-N1'	2.71	1.38	1.34
2	A	1001	TPP	C4'-N3'	2.53	1.38	1.35
2	B	1001	TPP	C4'-N3'	2.48	1.38	1.35
2	A	1001	TPP	C2'-N3'	2.45	1.38	1.34
2	B	1001	TPP	C2'-N3'	2.45	1.38	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	TPP	C2-N3	2.32	1.38	1.32
2	A	1001	TPP	C2-N3	2.29	1.37	1.32
2	A	1001	TPP	C6'-N1'	2.02	1.38	1.34

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	TPP	N1'-C2'-N3'	-3.29	120.06	125.53
2	A	1001	TPP	N1'-C2'-N3'	-3.22	120.17	125.53
2	A	1001	TPP	C6'-C5'-C4'	3.17	119.44	115.55
2	B	1001	TPP	C6'-C5'-C4'	3.12	119.39	115.55
2	A	1001	TPP	C6-C5-S1	-3.02	113.84	120.90
2	B	1001	TPP	C6-C5-S1	-2.98	113.91	120.90
2	B	1001	TPP	CM2-C2'-N1'	2.90	120.29	117.20
2	A	1001	TPP	CM2-C2'-N1'	2.80	120.19	117.20
2	B	1001	TPP	C6'-N1'-C2'	2.44	120.08	116.07
2	A	1001	TPP	C2-N3-C4	-2.40	110.82	114.06
2	B	1001	TPP	C2-N3-C4	-2.38	110.84	114.06
2	A	1001	TPP	C6'-N1'-C2'	2.38	119.98	116.07
2	A	1001	TPP	CM4-C4-N3	2.31	125.90	120.57
2	B	1001	TPP	CM4-C4-N3	2.29	125.84	120.57
2	B	1001	TPP	C5'-C6'-N1'	-2.25	120.18	123.83
2	A	1001	TPP	C4-C5-S1	2.24	114.05	110.56
2	A	1001	TPP	C5'-C6'-N1'	-2.23	120.20	123.83
2	B	1001	TPP	C4-C5-S1	2.22	114.02	110.56
2	B	1001	TPP	O2A-PA-O3A	2.05	112.80	107.27

There are no chirality outliers.

All (7) torsion outliers are listed below:

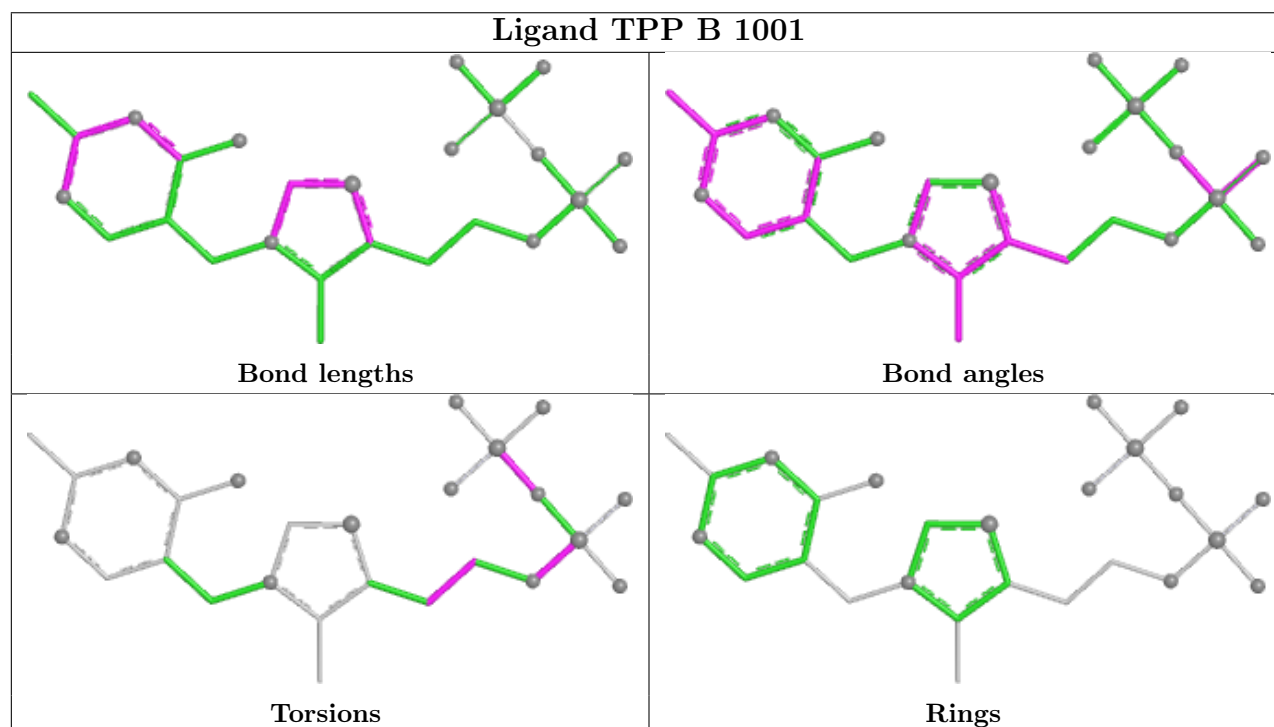
Mol	Chain	Res	Type	Atoms
2	A	1001	TPP	PA-O3A-PB-O2B
2	A	1001	TPP	PA-O3A-PB-O3B
2	B	1001	TPP	PA-O3A-PB-O3B
2	B	1001	TPP	C7-O7-PA-O1A
2	B	1001	TPP	PA-O3A-PB-O1B
2	A	1001	TPP	C5-C6-C7-O7
2	B	1001	TPP	C5-C6-C7-O7

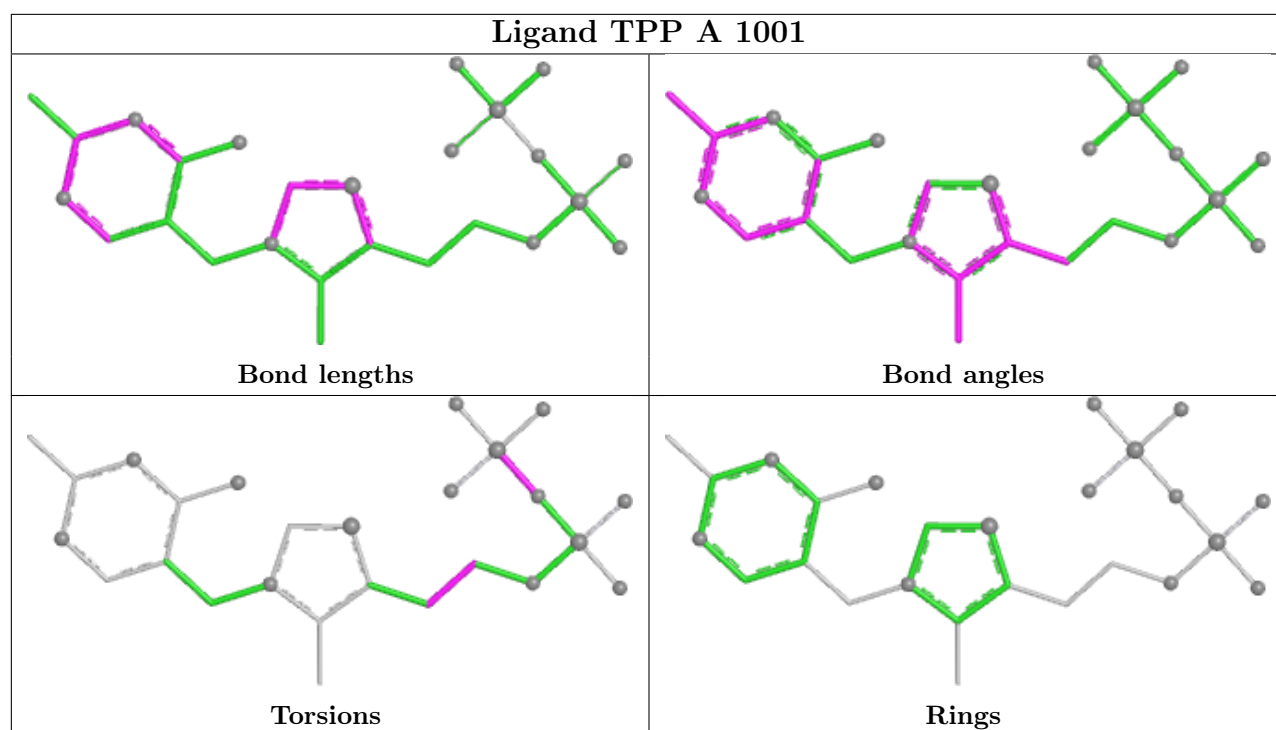
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1001	TPP	2	0
2	A	1001	TPP	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

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, 95th 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 ²)	Q<0.9
1	A	874/904 (96%)	0.53	57 (6%)	25 23	15, 37, 77, 151	1 (0%)
1	B	874/904 (96%)	0.56	65 (7%)	20 19	16, 37, 77, 137	1 (0%)
All	All	1748/1808 (96%)	0.55	122 (6%)	22 21	15, 37, 77, 151	2 (0%)

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	29	GLY	6.6
1	B	29	GLY	5.9
1	B	78	THR	5.5
1	A	78	THR	5.3
1	A	503	LEU	5.2
1	B	854	LYS	5.1
1	B	83	LEU	4.9
1	B	82	LEU	4.8
1	B	503	LEU	4.7
1	B	103	ALA	4.6
1	A	83	LEU	4.6
1	A	103	ALA	4.5
1	A	82	LEU	4.4
1	B	77	PHE	4.3
1	B	496	ALA	4.3
1	A	504	ASN	4.2
1	B	255	PHE	4.1
1	A	854	LYS	4.0
1	A	77	PHE	4.0
1	A	547	MET	3.9
1	B	504	ASN	3.9
1	B	84	GLU	3.8
1	A	496	ALA	3.8
1	B	553	LYS	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	105	LEU	3.6
1	A	255	PHE	3.6
1	A	550	HIS	3.5
1	A	501	PRO	3.4
1	A	110	LYS	3.4
1	A	494	ASN	3.4
1	A	84	GLU	3.4
1	B	80	GLN	3.3
1	B	370	TYR	3.3
1	A	564	MET	3.2
1	B	599	GLY	3.2
1	B	102	THR	3.2
1	A	30	VAL	3.2
1	A	105	LEU	3.2
1	B	613	THR	3.2
1	B	110	LYS	3.1
1	A	552	LEU	3.1
1	A	499	ARG	3.0
1	B	501	PRO	3.0
1	B	547	MET	3.0
1	B	519	ILE	3.0
1	B	81	ALA	2.9
1	B	276	HIS	2.9
1	A	495	MET	2.9
1	B	494	ASN	2.9
1	A	257	ALA	2.9
1	A	443	TYR	2.8
1	A	553	LYS	2.8
1	B	51	PRO	2.8
1	B	181	ALA	2.8
1	A	111	GLU	2.8
1	B	79	GLY	2.7
1	A	855	HIS	2.7
1	A	51	PRO	2.7
1	B	85	ASN	2.7
1	B	443	TYR	2.7
1	A	317	ASN	2.7
1	B	253	GLU	2.7
1	B	495	MET	2.6
1	B	499	ARG	2.6
1	B	564	MET	2.6
1	A	919	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	502	ALA	2.5
1	A	102	THR	2.5
1	A	505	LEU	2.5
1	A	502	ALA	2.5
1	A	824	GLY	2.5
1	A	850	MET	2.5
1	B	461	THR	2.5
1	A	251	PHE	2.5
1	B	256	SER	2.5
1	B	477	VAL	2.5
1	A	551	LEU	2.5
1	B	471	LEU	2.5
1	B	505	LEU	2.5
1	B	614	ASP	2.5
1	B	257	ALA	2.5
1	B	251	PHE	2.4
1	B	472	MET	2.4
1	A	181	ALA	2.4
1	B	246	ARG	2.4
1	B	469	GLY	2.4
1	A	827	LYS	2.4
1	A	81	ALA	2.4
1	B	111	GLU	2.4
1	B	485	TYR	2.3
1	A	79	GLY	2.3
1	B	30	VAL	2.3
1	B	317	ASN	2.3
1	A	519	ILE	2.3
1	A	375	GLU	2.3
1	A	444	THR	2.3
1	A	32	GLY	2.3
1	B	178	HIS	2.3
1	B	474	GLN	2.2
1	B	480	ILE	2.2
1	B	886	LYS	2.2
1	A	256	SER	2.2
1	A	276	HIS	2.2
1	B	254	ASN	2.2
1	B	828	HIS	2.2
1	A	227	LEU	2.1
1	A	471	LEU	2.1
1	A	472	MET	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	435	HIS	2.1
1	B	855	HIS	2.1
1	A	35	PRO	2.1
1	B	444	THR	2.1
1	A	800	LYS	2.1
1	B	550	HIS	2.1
1	A	480	ILE	2.1
1	B	35	PRO	2.1
1	A	104	GLY	2.1
1	A	558	SER	2.0
1	B	498	TYR	2.0
1	A	469	GLY	2.0
1	B	470	GLY	2.0
1	B	476	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	MLZ	B	72	10/11	0.81	0.17	24,42,57,65	0
1	MLZ	A	537	10/11	0.82	0.14	33,39,47,48	0
1	MLZ	B	537	10/11	0.87	0.13	32,38,46,47	0
1	MLZ	A	72	10/11	0.88	0.13	22,43,56,65	0

6.3 Carbohydrates [i](#)

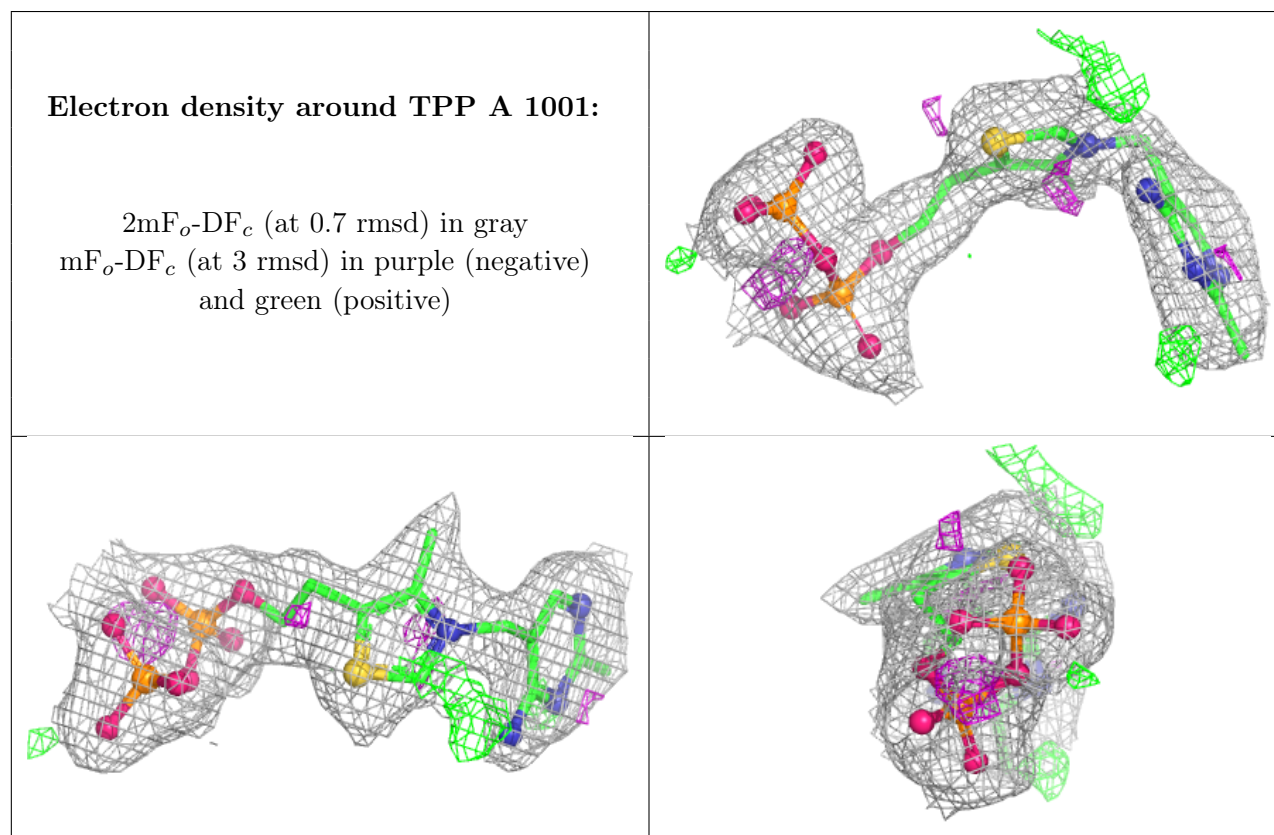
There are no oligosaccharides in this entry.

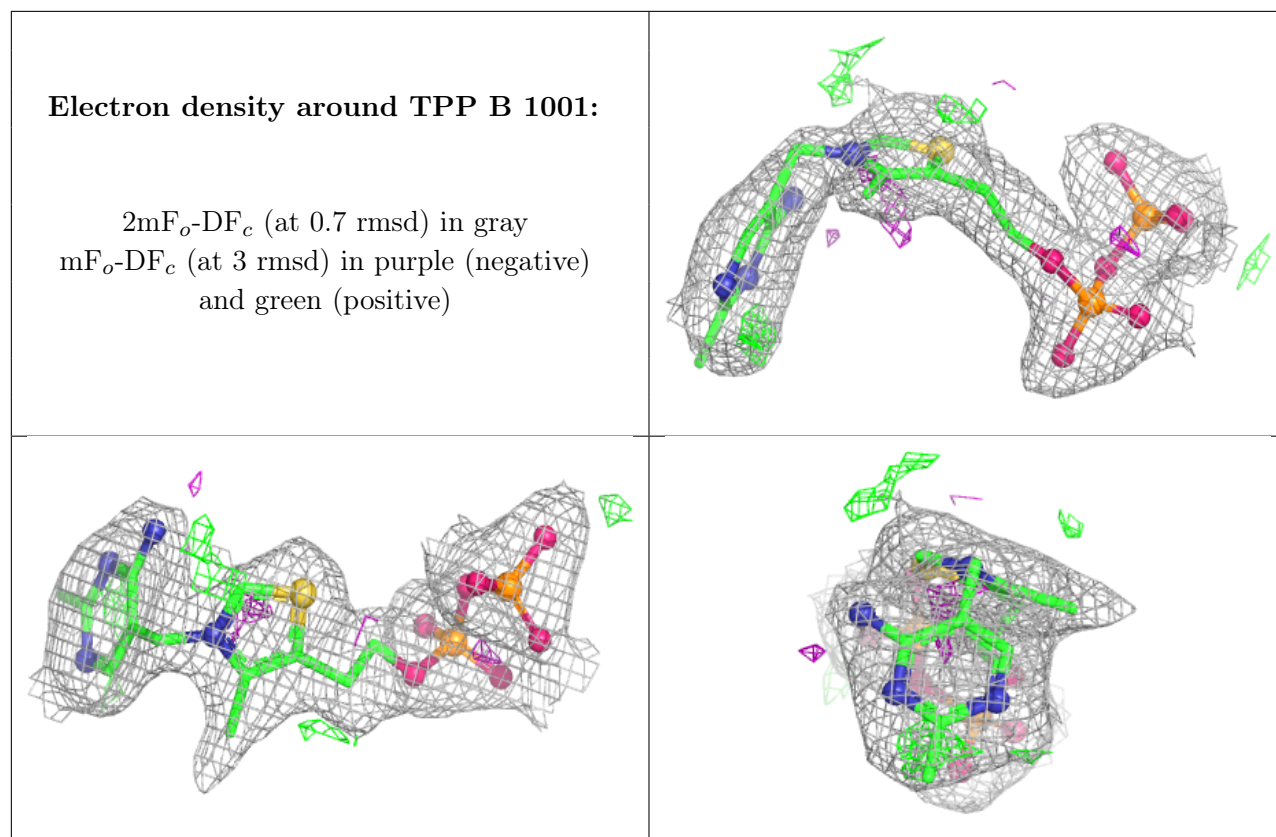
6.4 Ligands [i](#)

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, 95th 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
3	MG	B	1002	1/1	0.89	0.08	27,27,27,27	0
3	MG	A	1002	1/1	0.91	0.12	36,36,36,36	0
2	TPP	A	1001	26/26	0.91	0.10	18,35,51,59	0
2	TPP	B	1001	26/26	0.92	0.11	18,36,50,60	0
3	MG	A	1003	1/1	0.96	0.08	28,28,28,28	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.