



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:05 AM UTC

PDB ID : 7GTP / pdb_00007gtp
Title : PanDDA Analysis group deposition – Crystal structure of PTP1B in complex with FMOPL000688a
Authors : Mehlman, T.; Ginn, H.M.; Keedy, D.A.
Deposited on : 2024-01-03
Resolution : 2.47 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

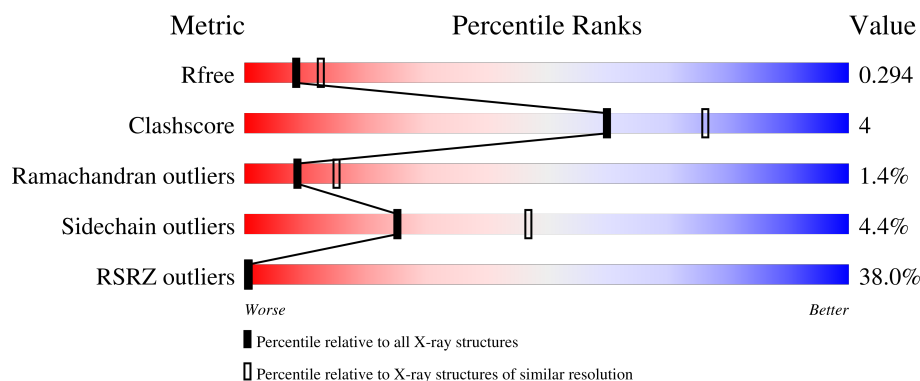
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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	321	

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	W1D	A	402	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5088 atoms, of which 2423 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-protein phosphatase non-receptor type 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	297	4829	1538	2402	418	456	15	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	32	SER	CYS	engineered mutation	UNP P18031
A	92	VAL	CYS	engineered mutation	UNP P18031

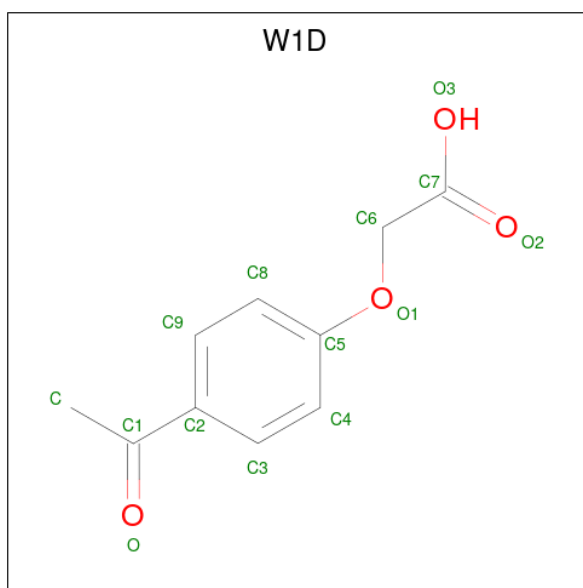
- Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O			
2	A	1	20	4	12	1	3		0	0

- Molecule 3 is (4-acetylphenoxy)acetic acid (CCD ID: W1D) (formula: $C_{10}H_{10}O_4$) (labeled

as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
3	A	1	23	10	9	4	0	0

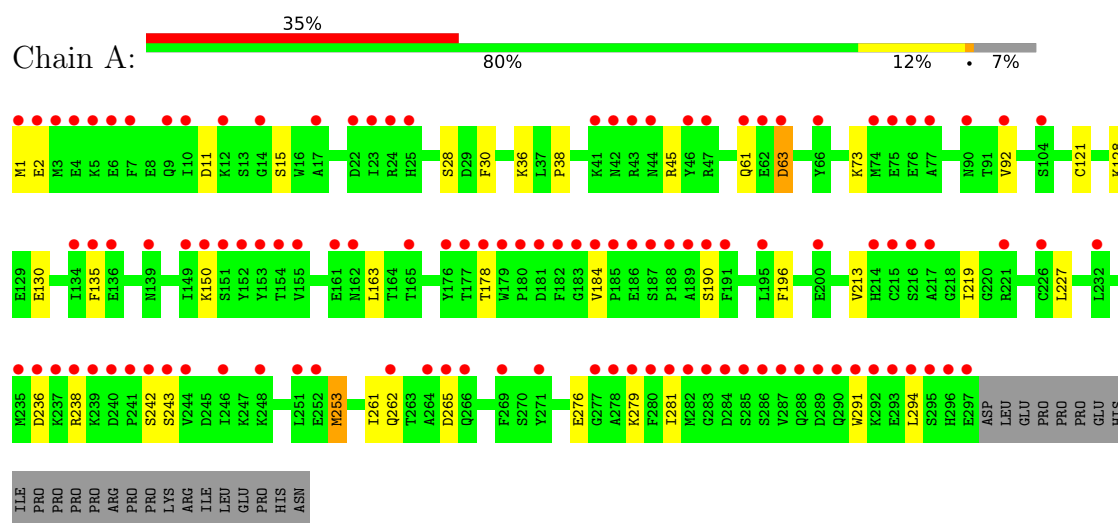
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	214	Total	O	0	4
			216	216		

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-protein phosphatase non-receptor type 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	90.10Å 90.10Å 107.17Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	78.03 – 2.47 78.03 – 2.47	Depositor EDS
% Data completeness (in resolution range)	100.0 (78.03-2.47) 99.9 (78.03-2.47)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.196 , 0.252 0.263 , 0.294	Depositor DCC
R_{free} test set	725 reflections (3.92%)	wwPDB-VP
Wilson B-factor (Å ²)	42.1	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 64.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	5088	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: W1D, TRS

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	1.03	0/2482	1.31	0/3345

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 [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	2427	2402	2393	20	2
2	A	8	12	12	1	2
3	A	14	9	0	0	0
4	A	216	0	0	8	1
All	All	2665	2423	2405	21	3

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 4.

All (21) 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:196:PHE:CE2	1:A:291:TRP:CZ3	2.61	0.89
1:A:73:LYS:HE3	4:A:563:HOH:O	1.80	0.81
1:A:73:LYS:HE2	4:A:506:HOH:O	1.95	0.66
1:A:128:LYS:NZ	4:A:509:HOH:O	2.30	0.64
1:A:92:VAL:HG12	1:A:135:PHE:CE1	2.34	0.62
1:A:38:PRO:HD2	4:A:543:HOH:O	2.00	0.60
2:A:401:TRS:O1	4:A:506:HOH:O	2.14	0.59
1:A:227:LEU:HB2	1:A:253:MET:HE1	1.87	0.57
1:A:196:PHE:HE2	1:A:291:TRP:CZ3	2.18	0.55
1:A:61:GLN:OE1	1:A:63:ASP:HB3	2.08	0.53
1:A:276:GLU:O	1:A:279:LYS:HG3	2.09	0.52
1:A:196:PHE:CE2	1:A:291:TRP:HZ3	2.21	0.51
1:A:163:LEU:O	4:A:507:HOH:O	2.18	0.48
1:A:45:ARG:NH2	1:A:121:CYS:HA	2.28	0.48
1:A:28:SER:HB3	1:A:30:PHE:CE2	2.48	0.48
1:A:219:ILE:HD12	1:A:262:GLN:NE2	2.30	0.46
1:A:227:LEU:CB	1:A:253:MET:HE1	2.46	0.46
1:A:92:VAL:CG1	1:A:135:PHE:CE1	2.99	0.45
1:A:36:LYS:HD2	4:A:628:HOH:O	2.20	0.41
1:A:1:MET:O	1:A:2:GLU:C	2.63	0.41
1:A:294:LEU:HB3	4:A:615:HOH:O	2.20	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:GLU:OE2	2:A:401:TRS:O2[4_565]	1.76	0.44
1:A:130:GLU:OE2	2:A:401:TRS:HO2[4_565]	1.45	0.15
4:A:521:HOH:O	4:A:651:HOH:O[3_454]	2.19	0.01

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	295/321 (92%)	276 (94%)	15 (5%)	4 (1%)	9	15

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	63	ASP
1	A	261	ILE
1	A	243	SER
1	A	281	ILE

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	270/294 (92%)	258 (96%)	12 (4%)	25	47

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASP
1	A	15	SER
1	A	150	LYS
1	A	178	THR
1	A	184	VAL
1	A	190	SER
1	A	213	VAL
1	A	236	ASP
1	A	238	ARG
1	A	242	SER
1	A	253	MET
1	A	265	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	193	ASN
1	A	262	GLN
1	A	296	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 oligosaccharides 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	$\# Z > 2$
3	W1D	A	402	-	14,14,14	1.47	2 (14%)	18,18,18	1.61	3 (16%)
2	TRS	A	401	-	7,7,7	0.35	0	9,9,9	0.57	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
3	W1D	A	402	-	-	4/9/9/9	0/1/1/1
2	TRS	A	401	-	-	4/9/9/9	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	W1D	C6-C7	4.10	1.62	1.50
3	A	402	W1D	C8-C5	2.14	1.42	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	W1D	C8-C5-C4	-3.21	115.47	120.16
3	A	402	W1D	C9-C2-C3	-2.29	115.66	118.57
3	A	402	W1D	C6-O1-C5	2.06	121.32	117.63

There are no chirality outliers.

All (8) torsion outliers are listed below:

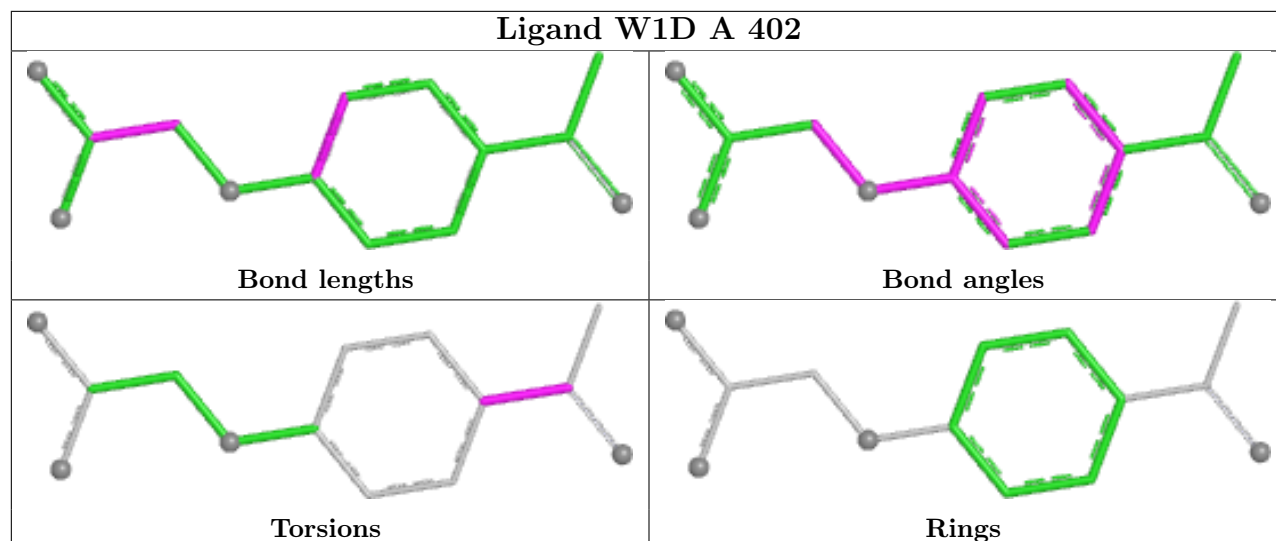
Mol	Chain	Res	Type	Atoms
2	A	401	TRS	C1-C-C2-O2
2	A	401	TRS	C3-C-C2-O2
2	A	401	TRS	N-C-C2-O2
3	A	402	W1D	C-C1-C2-C3
3	A	402	W1D	C-C1-C2-C9
3	A	402	W1D	O-C1-C2-C9
3	A	402	W1D	O-C1-C2-C3
2	A	401	TRS	C3-C-C1-O1

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	TRS	1	2

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	297/321 (92%)	3.99	113 (38%) 1 0	10, 34, 71, 115	68 (22%)

All (113) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	286	SER	18.9
1	A	190	SER	18.1
1	A	292	LYS	17.8
1	A	280	PHE	17.1
1	A	284	ASP	16.9
1	A	178	THR	16.5
1	A	187	SER	16.5
1	A	177	THR	16.4
1	A	271	TYR	16.0
1	A	290	GLN	16.0
1	A	153	TYR	15.6
1	A	289	ASP	15.6
1	A	240	ASP	15.3
1	A	179	TRP	15.3
1	A	287	VAL	15.1
1	A	241	PRO	15.1
1	A	184	VAL	15.0
1	A	296	HIS	15.0
1	A	188	PRO	15.0
1	A	152	TYR	14.9
1	A	180	PRO	14.8
1	A	269	PHE	14.7
1	A	151	SER	14.6
1	A	243	SER	14.6
1	A	183	GLY	14.5
1	A	189	ALA	14.4
1	A	149	ILE	14.3

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Mol	Chain	Res	Type	RSRZ
1	A	279	LYS	14.2
1	A	176	TYR	14.1
1	A	293	GLU	14.0
1	A	242	SER	13.9
1	A	244	VAL	13.9
1	A	288	GLN	13.9
1	A	294	LEU	13.7
1	A	75	GLU	13.7
1	A	185	PRO	13.7
1	A	154	THR	13.6
1	A	150	LYS	13.6
1	A	191	PHE	13.5
1	A	237	LYS	13.5
1	A	283	GLY	13.4
1	A	265	ASP	13.3
1	A	238	ARG	13.2
1	A	76	GLU	13.0
1	A	195	LEU	13.0
1	A	182	PHE	13.0
1	A	285	SER	12.9
1	A	215	CYS	12.8
1	A	236	ASP	12.7
1	A	181	ASP	12.5
1	A	186	GLU	12.3
1	A	262	GLN	12.2
1	A	155	VAL	12.2
1	A	295	SER	12.1
1	A	74	MET	12.0
1	A	252	GLU	12.0
1	A	281	ILE	12.0
1	A	282	MET	12.0
1	A	46	TYR	11.9
1	A	278	ALA	11.8
1	A	221	ARG	11.8
1	A	214	HIS	11.6
1	A	200	GLU	11.5
1	A	235	MET	11.3
1	A	216	SER	10.9
1	A	239	LYS	10.8
1	A	297	GLU	10.7
1	A	291	TRP	10.7
1	A	1	MET	7.1

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Mol	Chain	Res	Type	RSRZ
1	A	4	GLU	6.2
1	A	61	GLN	4.8
1	A	5	LYS	4.6
1	A	63	ASP	4.0
1	A	17	ALA	3.6
1	A	77	ALA	3.6
1	A	23	ILE	3.5
1	A	161	GLU	3.4
1	A	136	GLU	3.4
1	A	90	ASN	3.2
1	A	277	GLY	3.1
1	A	226	CYS	3.0
1	A	66	TYR	3.0
1	A	25	HIS	3.0
1	A	246	ILE	2.9
1	A	217	ALA	2.9
1	A	24	ARG	2.9
1	A	3	MET	2.9
1	A	62	GLU	2.8
1	A	2	GLU	2.8
1	A	12	LYS	2.8
1	A	47	ARG	2.7
1	A	92	VAL	2.7
1	A	251	LEU	2.7
1	A	7	PHE	2.6
1	A	43	ARG	2.5
1	A	264	ALA	2.5
1	A	165	THR	2.4
1	A	10	ILE	2.4
1	A	14	GLY	2.3
1	A	42	ASN	2.2
1	A	9	GLN	2.2
1	A	44	ASN	2.2
1	A	139	ASN	2.2
1	A	134	ILE	2.2
1	A	104	SER	2.2
1	A	22	ASP	2.2
1	A	41	LYS	2.1
1	A	266	GLN	2.1
1	A	6	GLU	2.1
1	A	162	ASN	2.0
1	A	248	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	135	PHE	2.0
1	A	232	LEU	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](#)

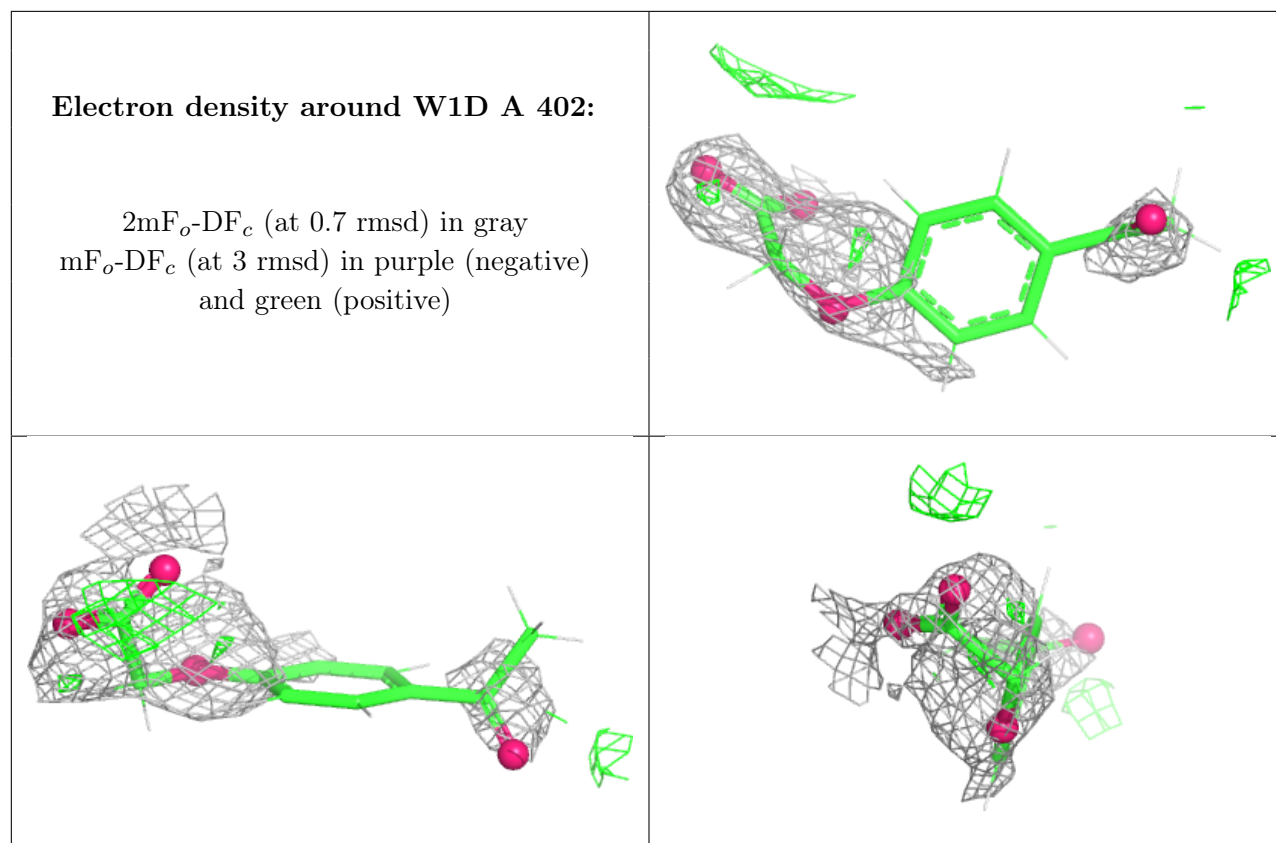
There are no oligosaccharides in this entry.

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	W1D	A	402	14/14	0.76	0.47	39,41,44,47	23
2	TRS	A	401	8/8	0.79	0.23	38,75,88,99	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 ⓘ

There are no such residues in this entry.