



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 06:33 PM EDT

PDB ID : 3IS4 / pdb_00003is4
Title : Crystal structure of Leishmania mexicana pyruvate kinase (LmPYK) in complex with 1,3,6,8-pyrenetetrasulfonic acid
Authors : Walkinshaw, M.D.; Morgan, H.P.
Deposited on : 2009-08-25
Resolution : 2.10 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

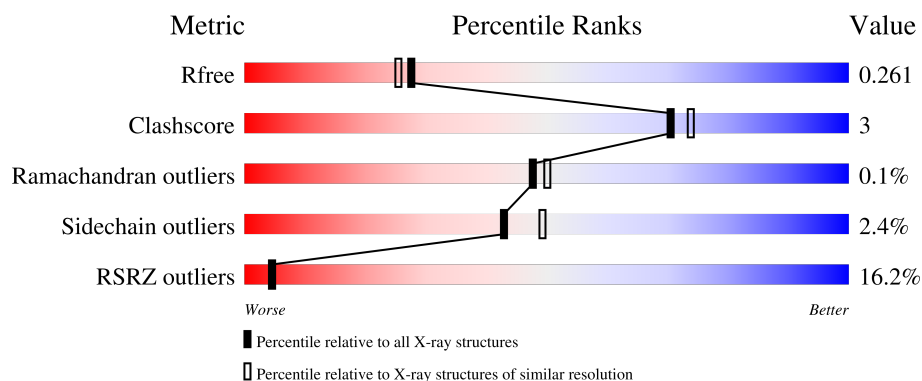
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	499	9% 90% 8% ..
1	B	499	23% 92% 6% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8211 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 Pyruvate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	492	Total	C	N	O	S	0	0	0
			3753	2339	660	728	26			
1	B	492	Total	C	N	O	S	0	0	0
			3753	2339	660	728	26			

There are 8 discrepancies between the modelled and reference sequences:

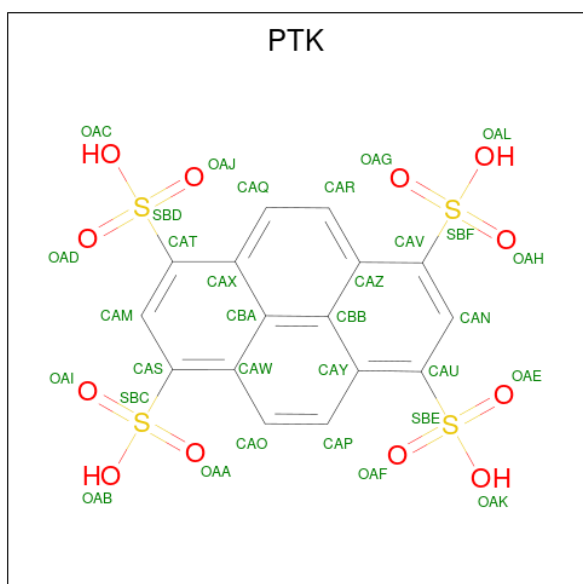
Chain	Residue	Modelled	Actual	Comment	Reference
A	382	SER	GLY	SEE REMARK 999	UNP Q27686
A	389	TYR	SER	SEE REMARK 999	UNP Q27686
A	404	ARG	ALA	SEE REMARK 999	UNP Q27686
A	405	SER	GLY	SEE REMARK 999	UNP Q27686
B	382	SER	GLY	SEE REMARK 999	UNP Q27686
B	389	TYR	SER	SEE REMARK 999	UNP Q27686
B	404	ARG	ALA	SEE REMARK 999	UNP Q27686
B	405	SER	GLY	SEE REMARK 999	UNP Q27686

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is pyrene-1,3,6,8-tetrasulfonic acid (CCD ID: PTK) (formula: $C_{16}H_{10}O_{12}S_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	S	0	0
			32	16	12	4		

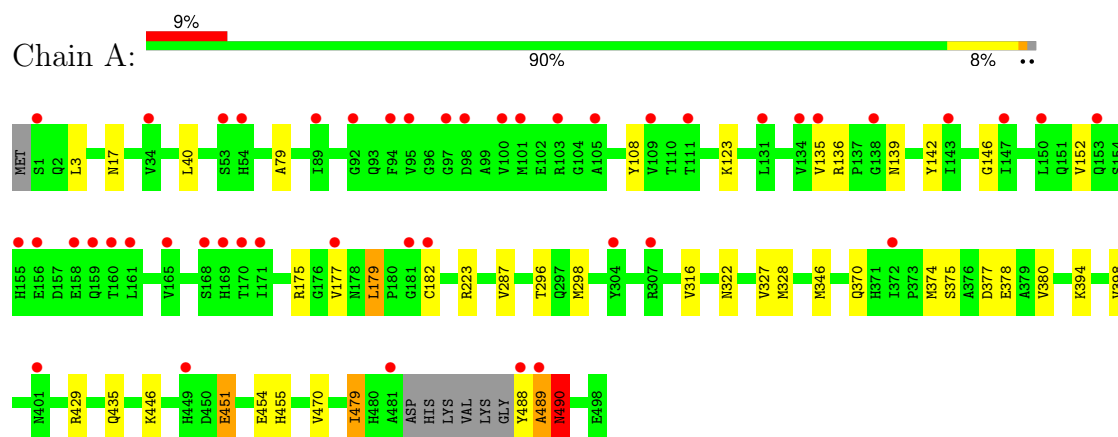
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	350	Total	O	0	0
			350	350		
4	B	293	Total	O	0	0
			293	293		

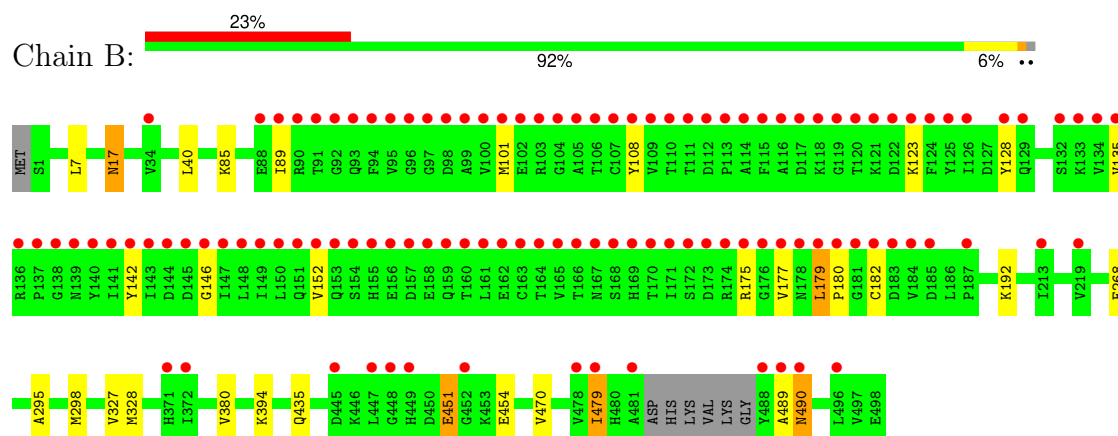
3 Residue-property plots [i](#)

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: Pyruvate kinase



• Molecule 1: Pyruvate kinase



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	122.86Å 129.86Å 165.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.69 – 2.10 20.69 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (20.69-2.10) 99.9 (20.69-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.33 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.231 , 0.263 0.230 , 0.261	Depositor DCC
R_{free} test set	3865 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	32.6	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8211	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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: PTK, GOL

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	0/3808	0.78	2/5154 (0.0%)
1	B	0.52	0/3808	0.77	0/5154
All	All	0.52	0/7616	0.77	2/10308 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	489	ALA	N-CA-C	-6.18	97.64	110.80
1	A	490	ASN	N-CA-CB	5.30	120.85	110.88

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	3753	0	3752	31	0
1	B	3753	0	3752	21	0
2	A	24	0	32	3	0
2	B	6	0	8	0	0
3	A	32	0	10	0	0
4	A	350	0	0	4	0
4	B	293	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8211	0	7554	52	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 (52) 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:298:MET:HE3	1:A:316:VAL:HG22	1.60	0.81
1:A:479:ILE:HD11	1:A:489:ALA:HB1	1.66	0.77
1:B:135:VAL:HG11	1:B:152:VAL:HG21	1.75	0.69
1:A:135:VAL:HG11	1:A:152:VAL:HG21	1.76	0.67
1:A:374:MET:HE2	1:A:378:GLU:HG3	1.84	0.60
1:A:398:VAL:HG13	1:A:479:ILE:HD13	1.84	0.59
1:B:479:ILE:HD11	1:B:489:ALA:CB	2.33	0.59
1:B:179:LEU:HB3	1:B:182:CYS:HB2	1.85	0.58
1:B:101:MET:HE1	4:B:600:HOH:O	2.03	0.57
1:B:394:LYS:HB2	1:B:470:VAL:HG12	1.86	0.57
1:A:298:MET:HE2	1:A:346:MET:HE2	1.88	0.56
1:B:298:MET:HE2	1:B:327:VAL:HB	1.87	0.55
1:A:179:LEU:HB3	1:A:182:CYS:HB2	1.89	0.54
1:A:223:ARG:HD2	4:A:553:HOH:O	2.06	0.54
1:B:479:ILE:HD11	1:B:489:ALA:HB1	1.89	0.54
1:A:455:HIS:HD2	4:A:559:HOH:O	1.91	0.53
1:B:17:ASN:H	1:B:17:ASN:HD22	1.56	0.53
1:A:377:ASP:HB3	1:A:488:TYR:CD2	2.44	0.53
1:A:451:GLU:H	1:A:451:GLU:CD	2.17	0.53
1:B:298:MET:CE	1:B:327:VAL:HB	2.39	0.53
1:A:380:VAL:HG21	1:A:490:ASN:HA	1.92	0.52
1:A:374:MET:HE2	1:A:378:GLU:CG	2.40	0.50
1:B:295:ALA:HB1	1:B:328:MET:HE2	1.94	0.50
1:B:380:VAL:HG21	1:B:490:ASN:HA	1.94	0.49
1:A:370:GLN:HG3	1:A:374:MET:HE3	1.94	0.49
1:A:375:SER:OG	1:A:377:ASP:OD2	2.23	0.47
1:B:180:PRO:HB3	1:B:268:GLU:HB2	1.97	0.47
1:A:479:ILE:HD11	1:A:489:ALA:CB	2.41	0.46
1:A:322:ASN:HA	2:A:502:GOL:H11	1.97	0.46
1:A:322:ASN:HA	2:A:502:GOL:C1	2.46	0.45
1:B:451:GLU:H	1:B:451:GLU:CD	2.25	0.45
1:A:287:VAL:HG21	1:B:7:LEU:HD11	1.99	0.45
1:B:85:LYS:HD3	1:B:192:LYS:HD3	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:TYR:O	1:B:123:LYS:HA	2.17	0.44
1:B:89:ILE:HG23	1:B:128:TYR:HB2	2.00	0.44
1:A:108:TYR:O	1:A:123:LYS:HA	2.18	0.44
1:B:17:ASN:HD22	1:B:17:ASN:N	2.13	0.43
1:A:394:LYS:HB2	1:A:470:VAL:HG12	1.99	0.43
1:A:488:TYR:CD2	1:A:489:ALA:O	2.71	0.43
1:B:142:TYR:HB3	1:B:146:GLY:HA2	2.00	0.43
1:B:490:ASN:HD22	1:B:490:ASN:H	1.66	0.43
1:A:298:MET:HG3	1:A:328:MET:O	2.19	0.42
1:A:3:LEU:C	1:A:3:LEU:HD23	2.45	0.41
1:A:298:MET:HE2	1:A:327:VAL:HB	2.01	0.41
1:B:101:MET:CE	4:B:600:HOH:O	2.64	0.41
1:A:136:ARG:H	1:A:139:ASN:ND2	2.19	0.41
1:A:142:TYR:HB3	1:A:146:GLY:HA2	2.02	0.41
1:A:398:VAL:HG22	1:A:479:ILE:HG23	2.03	0.41
1:A:298:MET:CE	1:A:316:VAL:HG22	2.40	0.41
1:A:446:LYS:HD3	4:A:815:HOH:O	2.19	0.41
2:A:502:GOL:H12	4:A:547:HOH:O	2.21	0.41
1:A:79:ALA:HB2	1:A:429:ARG:O	2.20	0.41

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	488/499 (98%)	476 (98%)	11 (2%)	1 (0%)	43	44
1	B	488/499 (98%)	479 (98%)	9 (2%)	0	100	100
All	All	976/998 (98%)	955 (98%)	20 (2%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	296	THR

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	411/417 (99%)	401 (98%)	10 (2%)	43	49
1	B	411/417 (99%)	401 (98%)	10 (2%)	43	49
All	All	822/834 (99%)	802 (98%)	20 (2%)	43	49

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	40	LEU
1	A	175	ARG
1	A	177	VAL
1	A	179	LEU
1	A	435	GLN
1	A	451	GLU
1	A	454	GLU
1	A	479	ILE
1	A	490	ASN
1	B	17	ASN
1	B	40	LEU
1	B	175	ARG
1	B	177	VAL
1	B	179	LEU
1	B	435	GLN
1	B	451	GLU
1	B	454	GLU
1	B	479	ILE
1	B	490	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	HIS
1	A	17	ASN
1	A	153	GLN
1	A	297	GLN
1	A	305	ASN
1	A	344	GLN
1	A	364	ASN
1	A	435	GLN
1	A	471	GLN
1	B	17	ASN
1	B	69	GLN
1	B	153	GLN
1	B	242	HIS
1	B	305	ASN
1	B	435	GLN
1	B	490	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

6 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
2	GOL	A	503	-	5,5,5	0.37	0	5,5,5	0.29	0
2	GOL	B	499	-	5,5,5	0.37	0	5,5,5	0.38	0
3	PTK	A	501	-	31,35,35	1.47	4 (12%)	52,60,60	1.06	4 (7%)
2	GOL	A	500	-	5,5,5	0.45	0	5,5,5	0.27	0
2	GOL	A	502	-	5,5,5	0.46	0	5,5,5	0.73	0
2	GOL	A	499	-	5,5,5	0.45	0	5,5,5	0.32	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
2	GOL	A	503	-	-	0/4/4/4	-
2	GOL	B	499	-	-	2/4/4/4	-
3	PTK	A	501	-	-	23/24/24/24	0/4/4/4
2	GOL	A	500	-	-	0/4/4/4	-
2	GOL	A	502	-	-	4/4/4/4	-
2	GOL	A	499	-	-	2/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	PTK	CAV-CAZ	-2.29	1.39	1.43
3	A	501	PTK	CAS-CAW	-2.23	1.39	1.43
3	A	501	PTK	CAT-CAX	-2.23	1.39	1.43
3	A	501	PTK	CAU-CAY	-2.23	1.39	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	PTK	CAR-CAZ-CAV	-2.55	119.74	123.58
3	A	501	PTK	CAQ-CAX-CAT	-2.48	119.86	123.58
3	A	501	PTK	CAP-CAY-CAU	-2.45	119.89	123.58
3	A	501	PTK	CAO-CAW-CAS	-2.43	119.93	123.58

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	502	GOL	O1-C1-C2-O2
2	A	502	GOL	O1-C1-C2-C3
2	A	502	GOL	C1-C2-C3-O3
2	A	502	GOL	O2-C2-C3-O3
3	A	501	PTK	CAM-CAS-SBC-OAB
3	A	501	PTK	CAM-CAS-SBC-OAI
3	A	501	PTK	CAW-CAS-SBC-OAA
3	A	501	PTK	CAW-CAS-SBC-OAB
3	A	501	PTK	CAW-CAS-SBC-OAI
3	A	501	PTK	CAM-CAT-SBD-OAC
3	A	501	PTK	CAM-CAT-SBD-OAJ
3	A	501	PTK	CAX-CAT-SBD-OAC
3	A	501	PTK	CAX-CAT-SBD-OAD
3	A	501	PTK	CAX-CAT-SBD-OAJ
3	A	501	PTK	CAY-CAU-SBE-OAE
3	A	501	PTK	CAY-CAU-SBE-OAF
3	A	501	PTK	CAN-CAV-SBF-OAG
3	A	501	PTK	CAN-CAV-SBF-OAL
3	A	501	PTK	CAZ-CAV-SBF-OAG
3	A	501	PTK	CAZ-CAV-SBF-OAH
3	A	501	PTK	CAZ-CAV-SBF-OAL
2	A	499	GOL	O1-C1-C2-C3
2	B	499	GOL	O1-C1-C2-C3
2	A	499	GOL	O1-C1-C2-O2
3	A	501	PTK	CAM-CAT-SBD-OAD
2	B	499	GOL	O1-C1-C2-O2
3	A	501	PTK	CAM-CAS-SBC-OAA
3	A	501	PTK	CAN-CAU-SBE-OAE
3	A	501	PTK	CAN-CAU-SBE-OAF
3	A	501	PTK	CAN-CAV-SBF-OAH
3	A	501	PTK	CAN-CAU-SBE-OAK

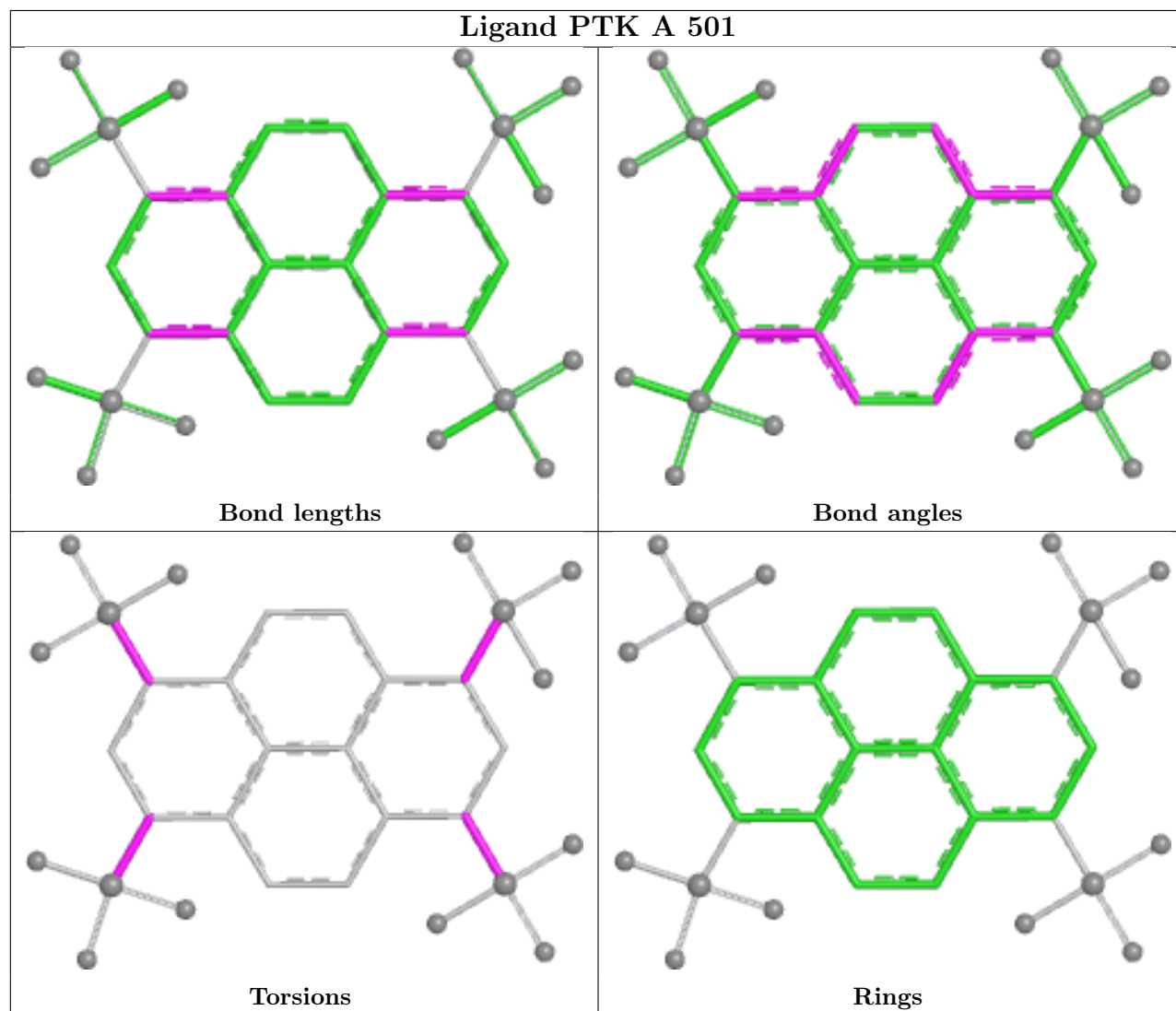
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	502	GOL	3	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	492/499 (98%)	0.55	46 (9%) 14 15	20, 33, 55, 60	0
1	B	492/499 (98%)	1.51	113 (22%) 2 2	21, 34, 55, 60	0
All	All	984/998 (98%)	1.03	159 (16%) 4 4	20, 34, 55, 60	0

All (159) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	171	ILE	12.0
1	B	94	PHE	11.1
1	B	105	ALA	10.3
1	B	166	THR	9.5
1	B	172	SER	9.2
1	B	149	ILE	9.1
1	B	124	PHE	8.9
1	B	142	TYR	8.8
1	B	109	VAL	8.6
1	B	106	THR	8.3
1	B	167	ASN	8.2
1	B	104	GLY	8.2
1	B	95	VAL	8.1
1	B	116	ALA	8.1
1	B	165	VAL	8.0
1	B	115	PHE	7.8
1	B	125	TYR	7.6
1	B	170	THR	7.5
1	B	169	HIS	7.3
1	B	97	GLY	7.1
1	B	163	CYS	7.1
1	B	119	GLY	7.1
1	B	146	GLY	7.0
1	B	113	PRO	7.0

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Mol	Chain	Res	Type	RSRZ
1	B	108	TYR	7.0
1	B	150	LEU	6.9
1	B	122	ASP	6.9
1	B	145	ASP	6.8
1	A	488	TYR	6.8
1	B	101	MET	6.7
1	B	164	THR	6.7
1	B	92	GLY	6.7
1	B	138	GLY	6.5
1	B	155	HIS	6.5
1	B	147	ILE	6.3
1	B	99	ALA	6.3
1	B	179	LEU	6.2
1	B	137	PRO	6.0
1	B	152	VAL	6.0
1	B	103	ARG	6.0
1	B	100	VAL	5.9
1	A	489	ALA	5.9
1	B	168	SER	5.8
1	B	126	ILE	5.7
1	B	140	TYR	5.7
1	B	96	GLY	5.6
1	B	182	CYS	5.6
1	B	489	ALA	5.5
1	B	120	THR	5.5
1	B	173	ASP	5.4
1	B	159	GLN	5.4
1	B	148	LEU	5.2
1	B	136	ARG	5.2
1	B	154	SER	5.2
1	B	93	GLN	5.2
1	B	107	CYS	5.2
1	B	139	ASN	5.0
1	B	117	ASP	4.9
1	B	151	GLN	4.9
1	B	114	ALA	4.8
1	B	110	THR	4.8
1	B	481	ALA	4.7
1	B	121	LYS	4.7
1	B	175	ARG	4.7
1	B	180	PRO	4.7
1	B	123	LYS	4.6

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Mol	Chain	Res	Type	RSRZ
1	B	153	GLN	4.6
1	B	102	GLU	4.5
1	B	143	ILE	4.5
1	B	162	GLU	4.5
1	B	91	THR	4.5
1	B	141	ILE	4.4
1	A	138	GLY	4.3
1	B	177	VAL	4.3
1	B	112	ASP	4.3
1	B	135	VAL	4.3
1	B	183	ASP	4.2
1	B	176	GLY	4.2
1	A	481	ALA	4.2
1	B	129	GLN	4.1
1	B	161	LEU	4.1
1	B	156	GLU	4.1
1	B	134	VAL	4.1
1	B	98	ASP	4.0
1	B	111	THR	3.9
1	B	90	ARG	3.9
1	B	488	TYR	3.9
1	A	170	THR	3.9
1	B	181	GLY	3.8
1	A	94	PHE	3.8
1	B	158	GLU	3.8
1	B	160	THR	3.7
1	B	118	LYS	3.7
1	B	144	ASP	3.7
1	B	157	ASP	3.7
1	B	89	ILE	3.4
1	B	178	ASN	3.3
1	A	53	SER	3.2
1	A	169	HIS	3.2
1	A	135	VAL	3.1
1	B	479	ILE	3.1
1	A	97	GLY	3.1
1	B	371	HIS	3.1
1	A	98	ASP	3.0
1	A	168	SER	3.0
1	A	109	VAL	2.9
1	B	132	SER	2.8
1	B	490	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	184	VAL	2.8
1	B	478	VAL	2.8
1	B	174	ARG	2.7
1	A	158	GLU	2.7
1	A	165	VAL	2.7
1	B	447	LEU	2.7
1	B	219	VAL	2.7
1	A	1	SER	2.6
1	A	177	VAL	2.6
1	B	133	LYS	2.6
1	A	304	TYR	2.6
1	A	156	GLU	2.6
1	A	134	VAL	2.5
1	B	452	GLY	2.5
1	A	182	CYS	2.5
1	A	372	ILE	2.5
1	A	150	LEU	2.4
1	A	449	HIS	2.4
1	A	153	GLN	2.4
1	B	88	GLU	2.4
1	A	401	ASN	2.4
1	A	100	VAL	2.4
1	B	34	VAL	2.4
1	A	89	ILE	2.3
1	B	128	TYR	2.3
1	A	101	MET	2.3
1	B	445	ASP	2.3
1	A	54	HIS	2.3
1	A	143	ILE	2.3
1	A	147	ILE	2.3
1	B	372	ILE	2.3
1	A	95	VAL	2.2
1	A	105	ALA	2.2
1	A	159	GLN	2.2
1	B	449	HIS	2.2
1	B	187	PRO	2.2
1	B	496	LEU	2.2
1	A	111	THR	2.2
1	A	171	ILE	2.1
1	A	155	HIS	2.1
1	B	213	ILE	2.1
1	A	34	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	103	ARG	2.1
1	A	160	THR	2.1
1	A	131	LEU	2.1
1	A	161	LEU	2.1
1	B	185	ASP	2.1
1	A	92	GLY	2.1
1	A	307	ARG	2.0
1	A	181	GLY	2.0
1	B	448	GLY	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](#)

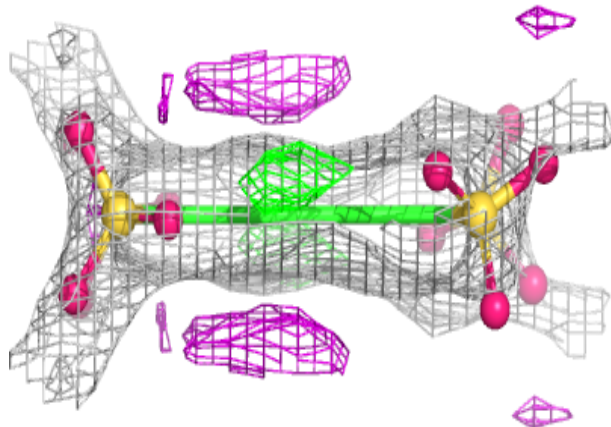
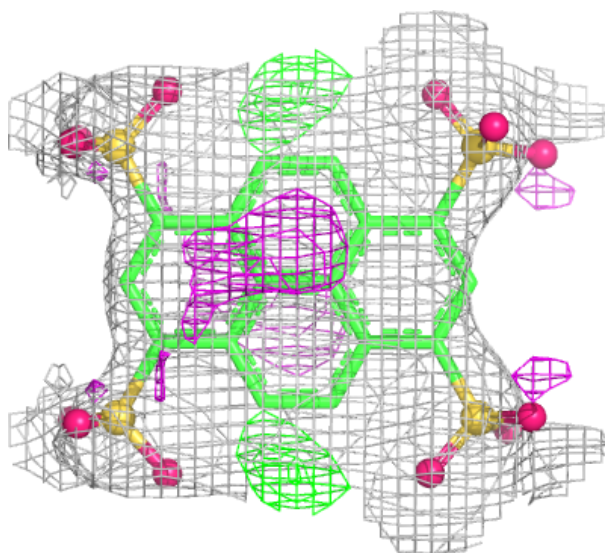
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PTK	A	501	32/32	0.54	0.20	100,100,101,101	32
2	GOL	A	499	6/6	0.69	0.24	41,44,45,46	0
2	GOL	B	499	6/6	0.75	0.18	67,68,68,68	0
2	GOL	A	500	6/6	0.79	0.17	63,63,63,63	0
2	GOL	A	503	6/6	0.82	0.17	59,61,62,62	0
2	GOL	A	502	6/6	0.83	0.17	36,41,42,43	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 PTK 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 [i](#)

There are no such residues in this entry.