



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:34 PM UTC

PDB ID : 1GFZ / pdb_00001gfz
Title : FLAVOPIRIDOL INHIBITS GLYCOGEN PHOSPHORYLASE BY BINDING AT THE INHIBITOR SITE
Authors : Oikonomakos, N.G.; Zographos, S.E.; Skamnaki, V.T.; Tsitsanou, K.E.; Johnson, L.N.
Deposited on : 2000-06-29
Resolution : 2.30 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)
Xtrriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

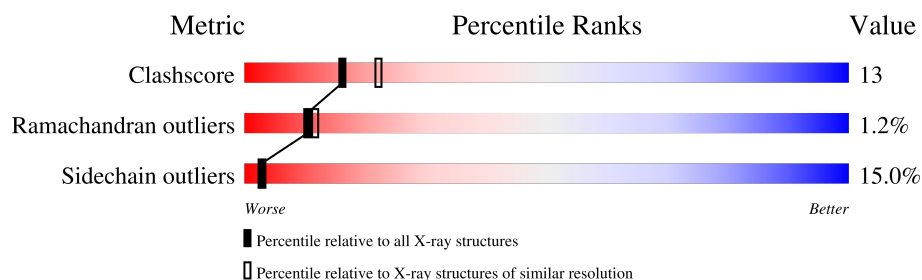
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	842	 64% 26% 6% ••

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6879 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.

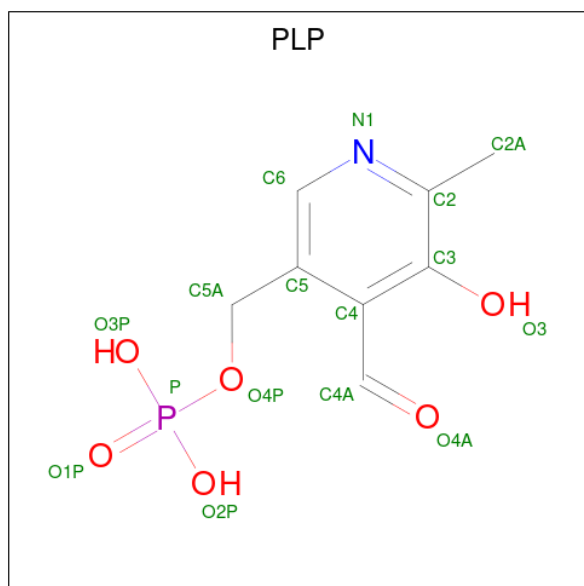
- Molecule 1 is a protein called GLYCOGEN PHOSPHORYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	814	6624	4226	1165	1204	29	0	0	0

There is a discrepancy between the modelled and reference sequences:

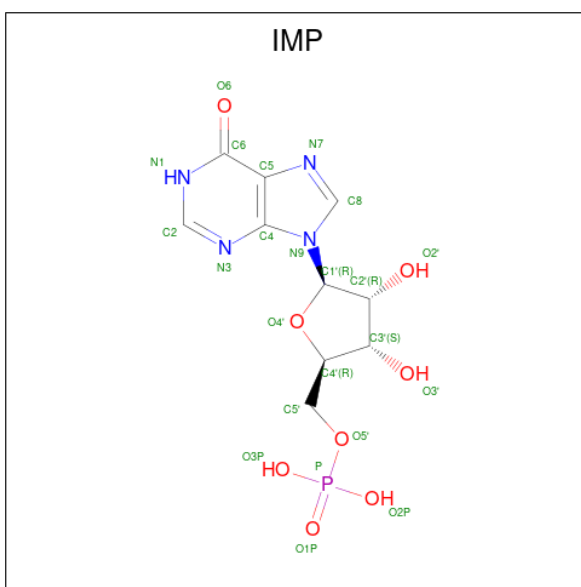
Chain	Residue	Modelled	Actual	Comment	Reference
A	380	ILE	LEU	conflict	UNP P00489

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



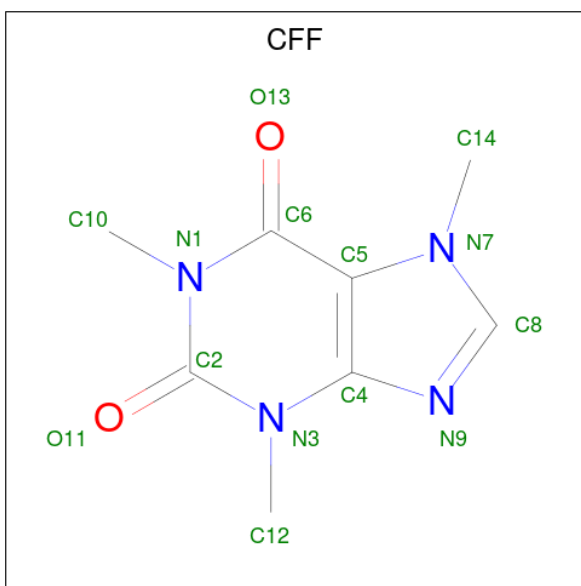
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	15	8	1	5	1	0	0

- Molecule 3 is INOSINIC ACID (CCD ID: IMP) (formula: $C_{10}H_{13}N_4O_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			23	10	4	8	1		

- Molecule 4 is CAFFEINE (CCD ID: CFF) (formula: $C_8H_{10}N_4O_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O		0	0
			14	8	4	2			

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	203	Total 203	O 203	0	0

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	128.30Å 128.30Å 116.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.75 – 2.30	Depositor
% Data completeness (in resolution range)	98.3 (28.75-2.30)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.189 , 0.219	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6879	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: IMP, CFF, PLP

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	0.49	0/6773	0.97	35/9165 (0.4%)

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	554	LYS	N-CA-C	8.69	129.30	110.80
1	A	714	ARG	N-CA-C	-7.45	99.43	110.46
1	A	576	GLN	N-CA-C	-7.17	103.64	112.38
1	A	91	MET	N-CA-C	7.08	120.39	111.69
1	A	565	VAL	N-CA-C	6.91	118.43	108.48
1	A	556	HIS	N-CA-C	6.64	124.95	110.80
1	A	105	GLU	N-CA-C	6.21	118.56	111.11
1	A	575	ARG	N-CA-C	6.01	119.68	111.39
1	A	656	VAL	CB-CA-C	-5.92	104.27	112.02
1	A	272	ALA	N-CA-C	-5.74	105.38	112.90
1	A	430	LEU	N-CA-C	-5.67	105.86	112.89
1	A	466	THR	N-CA-C	5.62	117.06	108.30
1	A	602	THR	N-CA-C	-5.59	98.23	108.02
1	A	396	LEU	CA-C-N	5.58	125.05	119.24
1	A	396	LEU	C-N-CA	5.58	125.05	119.24
1	A	16	ARG	N-CA-C	-5.55	106.71	112.93
1	A	278	VAL	N-CA-C	5.50	116.59	108.45
1	A	811	PHE	N-CA-C	5.47	119.06	112.93
1	A	723	GLN	N-CA-C	-5.44	105.25	111.07
1	A	767	HIS	N-CA-C	5.42	120.76	114.04
1	A	688	THR	N-CA-C	5.36	117.25	108.52
1	A	643	ILE	N-CA-C	5.28	115.50	108.11
1	A	277	ARG	N-CA-C	5.27	116.83	111.14
1	A	630	VAL	N-CA-C	5.27	115.47	110.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	387	TRP	N-CA-C	5.20	116.21	109.65
1	A	666	ILE	N-CA-C	5.19	120.14	109.34
1	A	555	VAL	N-CA-C	5.16	120.06	109.34
1	A	216	VAL	N-CA-C	5.15	117.03	108.99
1	A	507	ILE	N-CA-C	5.14	117.84	112.90
1	A	479	PHE	N-CA-C	5.12	117.90	109.76
1	A	115	LEU	N-CA-C	-5.10	107.03	113.20
1	A	37	PHE	N-CA-C	5.05	118.59	112.23
1	A	454	GLY	N-CA-C	-5.02	104.97	112.89
1	A	491	TRP	N-CA-C	5.01	119.11	113.15
1	A	677	GLY	N-CA-C	-5.01	106.72	112.73

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	6624	0	6571	174	0
2	A	15	0	6	0	0
3	A	23	0	11	0	0
4	A	14	0	10	0	0
5	A	203	0	0	18	0
All	All	6879	0	6598	174	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 13.

All (174) 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:251:ASP:CG	1:A:252:PHE:H	1.70	0.98
1:A:413:ARG:HH11	1:A:413:ARG:HG2	1.33	0.92
1:A:571:HIS:HB2	1:A:574:LYS:HG3	1.56	0.86
1:A:554:LYS:O	1:A:555:VAL:HG22	1.77	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:678:ASN:HD22	1:A:679:MET:H	1.25	0.84
1:A:455:VAL:H	1:A:459:HIS:HD2	1.28	0.81
1:A:764:MET:HE2	1:A:769:ASP:HA	1.61	0.81
1:A:316:PHE:HE1	1:A:332:LYS:HZ2	1.32	0.78
1:A:66:ARG:HG3	1:A:837:PRO:HB3	1.66	0.77
1:A:413:ARG:HG2	1:A:413:ARG:NH1	2.00	0.76
1:A:251:ASP:CG	1:A:252:PHE:N	2.43	0.76
1:A:15:VAL:HG21	1:A:509:GLU:CD	2.12	0.74
1:A:719:ASP:O	1:A:723:GLN:HG2	1.88	0.73
1:A:65:GLY:O	1:A:69:ARG:HG2	1.89	0.71
1:A:764:MET:HE1	5:A:1059:HOH:O	1.91	0.71
1:A:678:ASN:ND2	1:A:679:MET:H	1.88	0.70
1:A:515:LEU:HD22	1:A:518:LEU:HD22	1.72	0.69
1:A:90:TYR:HE1	5:A:1039:HOH:O	1.76	0.68
1:A:85:LEU:HD13	1:A:335:ILE:HG23	1.75	0.68
1:A:160:ARG:HB2	1:A:243:LEU:HB3	1.75	0.68
1:A:545:PHE:CZ	1:A:656:VAL:HG13	2.30	0.67
1:A:235:ASN:H	1:A:235:ASN:HD22	1.40	0.66
1:A:716:GLU:H	1:A:716:GLU:CD	2.02	0.66
1:A:546:ALA:HA	1:A:557:ILE:HD11	1.78	0.66
1:A:795:ARG:O	1:A:799:ARG:HG3	1.96	0.66
1:A:329:PHE:HB3	1:A:330:PRO:HD3	1.79	0.65
1:A:554:LYS:C	1:A:555:VAL:HG22	2.21	0.64
1:A:708:PHE:HB3	1:A:710:ILE:CD1	2.27	0.64
1:A:764:MET:CE	1:A:769:ASP:HA	2.26	0.64
1:A:13:ILE:O	1:A:14:SER:HB2	1.96	0.64
1:A:678:ASN:HD22	1:A:679:MET:N	1.96	0.64
1:A:15:VAL:HG13	1:A:15:VAL:O	1.98	0.62
1:A:764:MET:HE3	5:A:1055:HOH:O	1.99	0.61
1:A:703:ALA:HA	1:A:807:THR:HG21	1.82	0.61
1:A:138:ARG:O	1:A:138:ARG:HD3	2.01	0.61
1:A:61:ASP:O	1:A:64:VAL:HG22	2.01	0.60
1:A:708:PHE:HB3	1:A:710:ILE:HD12	1.83	0.60
1:A:360:ASP:OD1	1:A:363:LYS:HB2	2.02	0.60
1:A:16:ARG:HB3	1:A:105:GLU:HB3	1.83	0.59
1:A:648:TYR:HA	1:A:652:LEU:HD23	1.84	0.59
1:A:181:ASP:HB3	1:A:184:ARG:NH1	2.18	0.59
1:A:367:VAL:O	1:A:371:THR:HG23	2.03	0.58
1:A:601:ARG:HD2	5:A:1030:HOH:O	2.03	0.58
1:A:730:GLU:O	1:A:734:ARG:HG2	2.04	0.58
1:A:94:THR:HG22	5:A:1191:HOH:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:703:ALA:CA	1:A:807:THR:HG21	2.34	0.57
1:A:307:ILE:HD13	1:A:335:ILE:HD11	1.85	0.57
1:A:554:LYS:O	1:A:555:VAL:HG13	2.04	0.57
1:A:592:LYS:HG3	1:A:593:GLU:HG2	1.86	0.57
1:A:678:ASN:HD22	1:A:678:ASN:N	2.02	0.57
1:A:15:VAL:HG21	1:A:509:GLU:OE2	2.05	0.57
1:A:434:GLY:O	1:A:435:ALA:C	2.48	0.56
1:A:703:ALA:HB2	1:A:807:THR:HG21	1.87	0.56
1:A:493:VAL:CG2	1:A:512:ILE:HD12	2.35	0.56
1:A:34:HIS:HD2	1:A:38:THR:OG1	1.89	0.56
1:A:568:LYS:HE2	5:A:1106:HOH:O	2.06	0.56
1:A:144:LEU:HD12	1:A:147:MET:CE	2.36	0.56
1:A:379:VAL:HG22	5:A:1094:HOH:O	2.06	0.56
1:A:128:ASP:OD1	1:A:651:SER:HB3	2.06	0.55
1:A:90:TYR:CE1	5:A:1039:HOH:O	2.52	0.55
1:A:488:PRO:O	1:A:492:LEU:HB3	2.07	0.55
1:A:703:ALA:CB	1:A:807:THR:HG21	2.37	0.55
1:A:253:ASN:O	1:A:254:LEU:HB2	2.05	0.55
1:A:378:THR:OG1	1:A:380:ILE:HG12	2.08	0.54
1:A:144:LEU:HA	1:A:147:MET:HE3	1.89	0.54
1:A:458:ILE:HD11	1:A:715:VAL:HG21	1.89	0.54
1:A:571:HIS:HB2	1:A:574:LYS:CG	2.32	0.53
1:A:66:ARG:CG	1:A:837:PRO:HB3	2.38	0.53
1:A:184:ARG:HE	1:A:185:TYR:HE1	1.56	0.52
1:A:235:ASN:HA	1:A:833:ARG:HG3	1.91	0.52
1:A:325:ASN:HA	5:A:1164:HOH:O	2.10	0.52
1:A:355:ASP:OD2	1:A:398:ARG:HD3	2.08	0.52
1:A:311:PHE:O	1:A:314:SER:HB2	2.09	0.52
1:A:235:ASN:HD22	1:A:235:ASN:N	2.05	0.52
1:A:424:ARG:NH2	1:A:473:GLU:OE1	2.36	0.52
1:A:836:ALA:HB1	1:A:837:PRO:CD	2.40	0.52
1:A:467:ILE:HG23	1:A:468:PHE:CG	2.45	0.51
1:A:504:ALA:HA	1:A:508:GLY:O	2.09	0.51
1:A:604:MET:HB3	1:A:645:LEU:HD22	1.91	0.51
1:A:313:SER:O	1:A:314:SER:C	2.53	0.51
1:A:343:SER:OG	1:A:441:MET:HE3	2.11	0.51
1:A:351:ARG:O	1:A:355:ASP:HB2	2.11	0.51
1:A:423:ASP:O	1:A:426:ARG:HG3	2.10	0.51
1:A:550:GLU:O	1:A:554:LYS:HG2	2.12	0.50
1:A:94:THR:CG2	5:A:1191:HOH:O	2.58	0.50
1:A:457:ARG:HH22	1:A:701:GLU:CD	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:764:MET:HE2	1:A:769:ASP:CA	2.39	0.49
1:A:678:ASN:ND2	1:A:679:MET:N	2.56	0.49
1:A:720:ARG:CD	5:A:1193:HOH:O	2.61	0.49
1:A:34:HIS:HE1	1:A:61:ASP:OD2	1.96	0.48
1:A:252:PHE:C	1:A:254:LEU:H	2.20	0.48
1:A:720:ARG:HD3	5:A:1193:HOH:O	2.13	0.48
1:A:316:PHE:HE1	1:A:332:LYS:NZ	2.08	0.48
1:A:678:ASN:ND2	1:A:678:ASN:N	2.60	0.48
1:A:810:LYS:O	1:A:810:LYS:HG2	2.14	0.48
1:A:835:PRO:O	1:A:836:ALA:O	2.32	0.47
1:A:216:VAL:HG13	1:A:217:ASP:CG	2.39	0.47
1:A:639:ARG:HG2	5:A:1160:HOH:O	2.14	0.47
1:A:604:MET:HB3	1:A:645:LEU:CD2	2.45	0.47
1:A:493:VAL:HG21	1:A:512:ILE:HD12	1.96	0.47
1:A:91:MET:HB2	1:A:129:ALA:HB3	1.97	0.46
1:A:170:ILE:HG12	1:A:646:GLU:HG3	1.96	0.46
1:A:554:LYS:O	1:A:555:VAL:CG2	2.55	0.46
1:A:716:GLU:CD	1:A:716:GLU:N	2.71	0.46
1:A:181:ASP:HB3	1:A:184:ARG:HH11	1.80	0.46
1:A:758:PHE:HD1	1:A:761:ILE:HD12	1.81	0.46
1:A:95:LEU:HD11	1:A:99:MET:HE3	1.98	0.46
1:A:689:ILE:HG23	1:A:689:ILE:O	2.15	0.46
1:A:268:ASP:O	1:A:271:LEU:HB2	2.16	0.46
1:A:463:LEU:HA	1:A:467:ILE:HG22	1.97	0.46
1:A:340:THR:O	1:A:343:SER:HB3	2.16	0.45
1:A:693:ASP:O	1:A:696:ASN:HB2	2.16	0.45
1:A:562:LEU:HD21	1:A:662:LEU:HB2	1.98	0.45
1:A:834:LEU:O	1:A:836:ALA:N	2.50	0.45
1:A:499:LEU:HD12	1:A:502:ILE:HD11	1.99	0.45
1:A:93:ARG:HH11	1:A:93:ARG:CG	2.30	0.45
1:A:251:ASP:C	1:A:255:LYS:HB3	2.42	0.45
1:A:834:LEU:HG	1:A:835:PRO:HD2	1.99	0.45
1:A:413:ARG:HH11	1:A:413:ARG:CG	2.13	0.45
1:A:836:ALA:HB1	1:A:837:PRO:HD2	1.97	0.45
1:A:209:THR:HG21	5:A:1092:HOH:O	2.16	0.44
1:A:575:ARG:HD3	1:A:666:ILE:O	2.17	0.44
1:A:726:TYR:OH	1:A:774:PHE:HB2	2.17	0.44
1:A:518:LEU:HD12	1:A:518:LEU:HA	1.90	0.44
1:A:703:ALA:HB2	1:A:807:THR:CG2	2.47	0.44
1:A:709:PHE:CD1	1:A:787:VAL:HG12	2.52	0.44
1:A:14:SER:HB3	1:A:16:ARG:HG2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:ASN:HA	1:A:269:ARG:HH12	1.83	0.44
1:A:254:LEU:HD21	1:A:266:VAL:HG23	2.00	0.44
1:A:81:ARG:CG	1:A:155:TYR:HE1	2.31	0.44
1:A:455:VAL:H	1:A:459:HIS:CD2	2.20	0.44
1:A:689:ILE:HA	1:A:709:PHE:HB2	2.01	0.43
1:A:699:MET:HE2	1:A:708:PHE:HZ	1.83	0.43
1:A:579:ASN:C	1:A:579:ASN:HD22	2.25	0.43
1:A:503:ILE:HG12	1:A:521:LEU:HD11	2.00	0.43
1:A:687:LEU:HD12	1:A:687:LEU:HA	1.81	0.43
1:A:737:GLU:HG3	5:A:1150:HOH:O	2.18	0.43
1:A:790:LEU:HD22	1:A:790:LEU:O	2.19	0.43
1:A:564:ASP:O	1:A:603:VAL:HA	2.19	0.43
1:A:67:TRP:HA	1:A:238:VAL:HB	2.00	0.43
1:A:254:LEU:HD21	1:A:266:VAL:CG2	2.49	0.43
1:A:417:ALA:C	1:A:419:PRO:HD3	2.43	0.43
1:A:450:HIS:HE1	5:A:1135:HOH:O	2.02	0.43
1:A:732:TYR:CZ	1:A:739:ARG:HG3	2.54	0.42
1:A:735:ILE:HG22	1:A:738:LEU:H	1.84	0.42
1:A:777:TYR:O	1:A:781:VAL:HG13	2.19	0.42
1:A:36:HIS:O	1:A:40:VAL:HA	2.19	0.42
1:A:115:LEU:HD12	1:A:115:LEU:HA	1.90	0.42
1:A:398:ARG:O	1:A:398:ARG:HG3	2.19	0.42
1:A:466:THR:HG22	1:A:467:ILE:N	2.34	0.42
1:A:720:ARG:NH1	5:A:1193:HOH:O	2.52	0.42
1:A:144:LEU:HD12	1:A:147:MET:HE1	2.00	0.42
1:A:690:GLY:O	1:A:710:ILE:HA	2.20	0.42
1:A:484:ASN:ND2	5:A:1014:HOH:O	2.53	0.42
1:A:142:CYS:SG	1:A:487:THR:HG22	2.59	0.42
1:A:63:LEU:HB3	1:A:102:LEU:HD21	2.02	0.41
1:A:389:VAL:HG13	1:A:400:LEU:HD11	2.01	0.41
1:A:663:SER:HB2	1:A:681:PHE:CG	2.56	0.41
1:A:387:TRP:HA	1:A:388:PRO:HD3	1.97	0.41
1:A:467:ILE:HG23	1:A:468:PHE:N	2.35	0.41
1:A:568:LYS:HE3	1:A:574:LYS:HD2	2.01	0.41
1:A:430:LEU:HD22	1:A:443:HIS:HB3	2.03	0.41
1:A:627:GLY:HA2	1:A:642:VAL:HB	2.03	0.41
1:A:467:ILE:CG2	1:A:468:PHE:N	2.84	0.41
1:A:558:ASN:HA	1:A:559:PRO:HD3	1.89	0.41
1:A:204:GLY:HA2	1:A:217:ASP:O	2.21	0.41
1:A:386:ARG:HA	1:A:439:ILE:O	2.20	0.41
1:A:758:PHE:CD1	1:A:761:ILE:HD12	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:713:MET:HE3	1:A:713:MET:HB2	1.90	0.40
1:A:424:ARG:HH22	1:A:473:GLU:CD	2.25	0.40
1:A:605:ILE:O	1:A:644:PHE:HA	2.21	0.40
1:A:110:GLU:O	1:A:113:TYR:HB3	2.21	0.40
1:A:209:THR:O	1:A:211:GLN:N	2.53	0.40
1:A:252:PHE:C	1:A:254:LEU:N	2.78	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	808/842 (96%)	760 (94%)	38 (5%)	10 (1%)	10 12

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	14	SER
1	A	210	SER
1	A	254	LEU
1	A	555	VAL
1	A	556	HIS
1	A	836	ALA
1	A	314	SER
1	A	835	PRO
1	A	43	ARG
1	A	554	LYS

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	705/731 (96%)	599 (85%)	106 (15%)	3 3

All (106) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ILE
1	A	16	ARG
1	A	18	LEU
1	A	22	GLU
1	A	23	ASN
1	A	27	LEU
1	A	29	LYS
1	A	33	ARG
1	A	39	LEU
1	A	45	VAL
1	A	69	ARG
1	A	87	LEU
1	A	93	ARG
1	A	94	THR
1	A	102	LEU
1	A	104	LEU
1	A	115	LEU
1	A	122	LEU
1	A	128	ASP
1	A	138	ARG
1	A	144	LEU
1	A	178	GLU
1	A	198	LEU
1	A	210	SER
1	A	211	GLN
1	A	214	LYS
1	A	216	VAL
1	A	234	ARG
1	A	235	ASN
1	A	243	LEU

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Mol	Chain	Res	Type
1	A	247	LYS
1	A	250	ASN
1	A	253	ASN
1	A	254	LEU
1	A	255	LYS
1	A	264	GLN
1	A	271	LEU
1	A	279	LEU
1	A	287	GLU
1	A	292	ARG
1	A	306	ASP
1	A	314	SER
1	A	315	LYS
1	A	324	THR
1	A	337	LEU
1	A	358	ARG
1	A	363	LYS
1	A	371	THR
1	A	398	ARG
1	A	400	LEU
1	A	413	ARG
1	A	425	LEU
1	A	430	LEU
1	A	433	GLU
1	A	436	VAL
1	A	444	LEU
1	A	465	LYS
1	A	474	LEU
1	A	510	GLU
1	A	515	LEU
1	A	518	LEU
1	A	522	LEU
1	A	528	GLU
1	A	536	LYS
1	A	544	LYS
1	A	549	LEU
1	A	552	GLU
1	A	555	VAL
1	A	565	VAL
1	A	568	LYS
1	A	586	LEU
1	A	593	GLU

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Mol	Chain	Res	Type
1	A	596	LYS
1	A	603	VAL
1	A	622	LEU
1	A	641	ARG
1	A	643	ILE
1	A	645	LEU
1	A	656	VAL
1	A	662	LEU
1	A	666	ILE
1	A	678	ASN
1	A	687	LEU
1	A	706	GLU
1	A	708	PHE
1	A	716	GLU
1	A	720	ARG
1	A	721	LEU
1	A	724	ARG
1	A	734	ARG
1	A	742	ILE
1	A	759	LYS
1	A	763	ASN
1	A	764	MET
1	A	765	LEU
1	A	768	HIS
1	A	772	LYS
1	A	781	VAL
1	A	782	LYS
1	A	787	VAL
1	A	790	LEU
1	A	807	THR
1	A	812	SER
1	A	813	SER
1	A	830	SER
1	A	833	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	208	HIS
1	A	211	GLN
1	A	219	GLN

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Mol	Chain	Res	Type
1	A	235	ASN
1	A	325	ASN
1	A	390	HIS
1	A	412	ASN
1	A	450	HIS
1	A	459	HIS
1	A	477	HIS
1	A	481	ASN
1	A	484	ASN
1	A	517	GLN
1	A	560	ASN
1	A	566	GLN
1	A	579	ASN
1	A	678	ASN
1	A	763	ASN
1	A	832	GLN

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](#)

3 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	IMP	A	930	-	25,25,25	1.63	5 (20%)	37,38,38	1.51	7 (18%)
2	PLP	A	999	1	15,15,16	1.45	3 (20%)	21,22,23	1.12	3 (14%)
4	CFF	A	940	-	15,15,15	2.59	5 (33%)	23,23,23	2.06	7 (30%)

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	IMP	A	930	-	-	5/10/26/26	0/3/3/3
2	PLP	A	999	1	-	2/6/6/8	0/1/1/1
4	CFF	A	940	-	-	-	0/2/2/2

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	940	CFF	C2-N1	7.78	1.54	1.39
3	A	930	IMP	C8-N7	3.85	1.43	1.32
3	A	930	IMP	C5'-C4'	3.30	1.61	1.51
3	A	930	IMP	C4-N3	-3.10	1.29	1.35
2	A	999	PLP	C5A-C5	2.92	1.58	1.50
4	A	940	CFF	O13-C6	2.87	1.28	1.23
2	A	999	PLP	C3-C2	-2.72	1.38	1.41
4	A	940	CFF	C10-N1	2.64	1.51	1.47
4	A	940	CFF	C4-N3	2.64	1.42	1.38
3	A	930	IMP	P-O3P	-2.47	1.45	1.54
3	A	930	IMP	C1'-N9	2.24	1.53	1.47
4	A	940	CFF	C2-N3	-2.23	1.34	1.39
2	A	999	PLP	P-O2P	2.10	1.62	1.54

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	940	CFF	O11-C2-N3	5.47	128.90	121.33
4	A	940	CFF	C2-N3-C4	4.46	125.31	119.58
4	A	940	CFF	N3-C2-N1	-3.57	112.67	117.14
3	A	930	IMP	N1-C2-N3	-3.42	121.56	125.50
3	A	930	IMP	O4'-C4'-C3'	-3.02	99.16	105.15
4	A	940	CFF	C12-N3-C4	-2.98	116.36	120.75
3	A	930	IMP	O4'-C4'-C5'	-2.89	100.06	109.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	930	IMP	N9-C8-N7	-2.67	108.45	113.40
2	A	999	PLP	O3P-P-O1P	2.64	121.12	110.83
3	A	930	IMP	C6-C5-C4	2.44	122.49	118.83
3	A	930	IMP	C5-C6-N1	-2.36	107.39	110.78
3	A	930	IMP	C8-N9-C4	2.28	110.29	106.03
2	A	999	PLP	C6-C5-C4	2.19	119.89	118.10
2	A	999	PLP	C5-C6-N1	-2.15	120.33	123.83
4	A	940	CFF	O11-C2-N1	-2.10	118.43	121.33
4	A	940	CFF	C6-C5-C4	-2.04	120.34	122.92
4	A	940	CFF	C5-C6-N1	2.04	115.79	112.06

There are no chirality outliers.

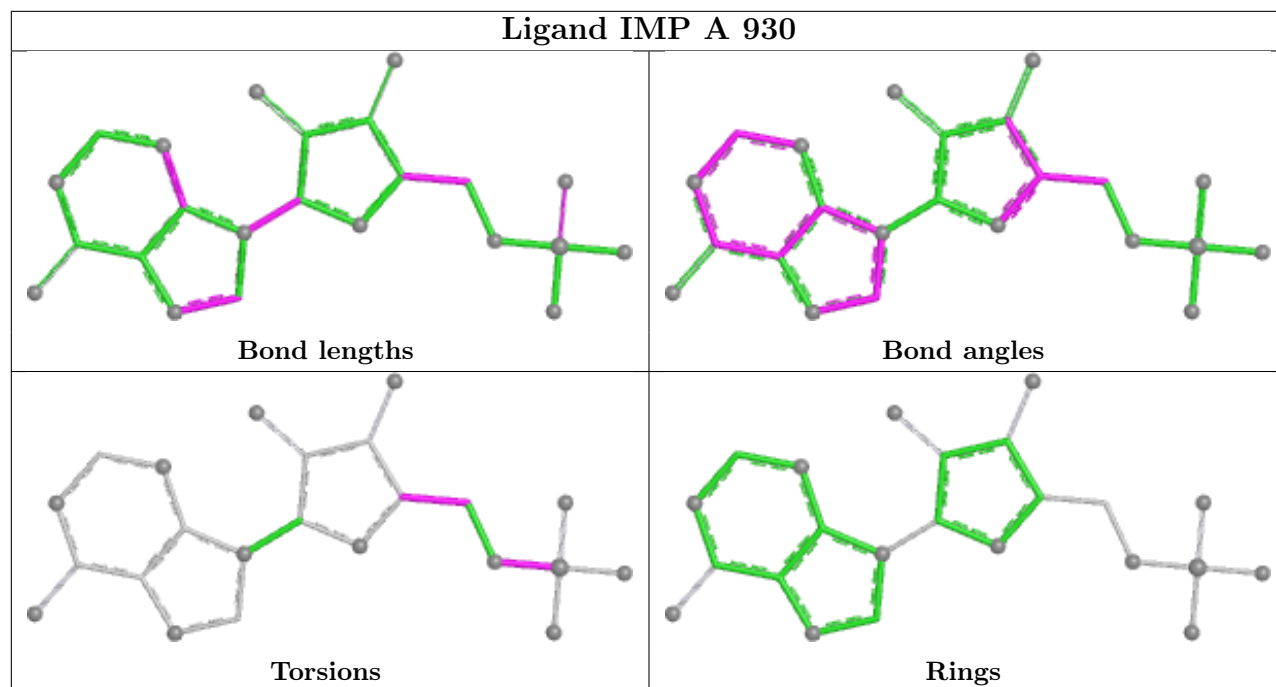
All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	930	IMP	C5'-O5'-P-O1P
3	A	930	IMP	C5'-O5'-P-O2P
3	A	930	IMP	C5'-O5'-P-O3P
3	A	930	IMP	C3'-C4'-C5'-O5'
3	A	930	IMP	O4'-C4'-C5'-O5'
2	A	999	PLP	C6-C5-C5A-O4P
2	A	999	PLP	C4-C5-C5A-O4P

There are no ring outliers.

No monomer is involved in short contacts.

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.