



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:59 PM UTC

PDB ID : 1C8K / pdb_00001c8k
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 : 2000-05-11
Resolution : 1.76 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)	:	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

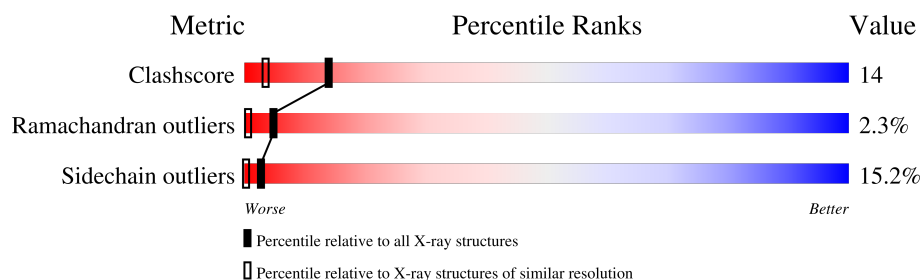
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.76 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	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)

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	67% 22% 7% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7123 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.

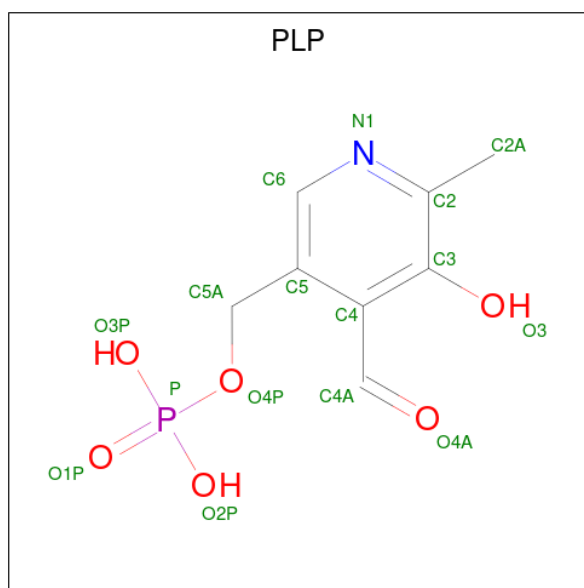
- Molecule 1 is a protein called PROTEIN (GLYCOGEN PHOSPHORYLASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	830	6749	4303	1189	1227	30	0	0	0

There is a discrepancy between the modelled and reference sequences:

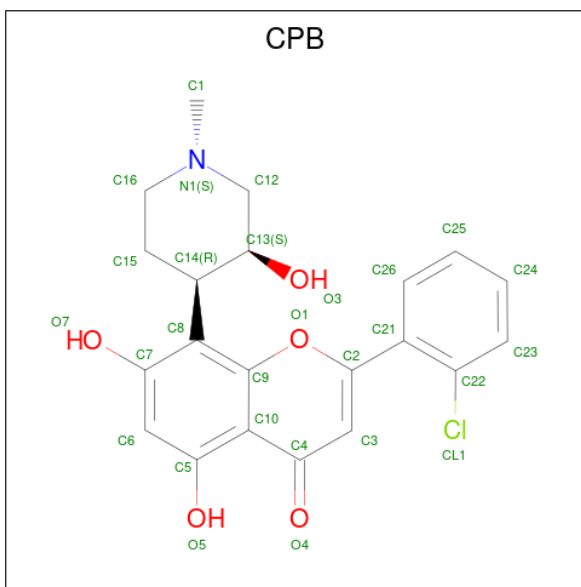
Chain	Residue	Modelled	Actual	Comment	Reference
A	380	ILE	LEU	conflict	UNP P00489

- 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 2-(2-CHLORO-PHENYL)-5,7-DIHYDROXY-8-(3-HYDROXY-1-METHYL-P IPERIDIN-4-YL)-4H-BENZOPYRAN-4-ONE (CCD ID: CPB) (formula: $C_{21}H_{20}ClNO_5$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			28	21	1	1	5		

- Molecule 4 is water.

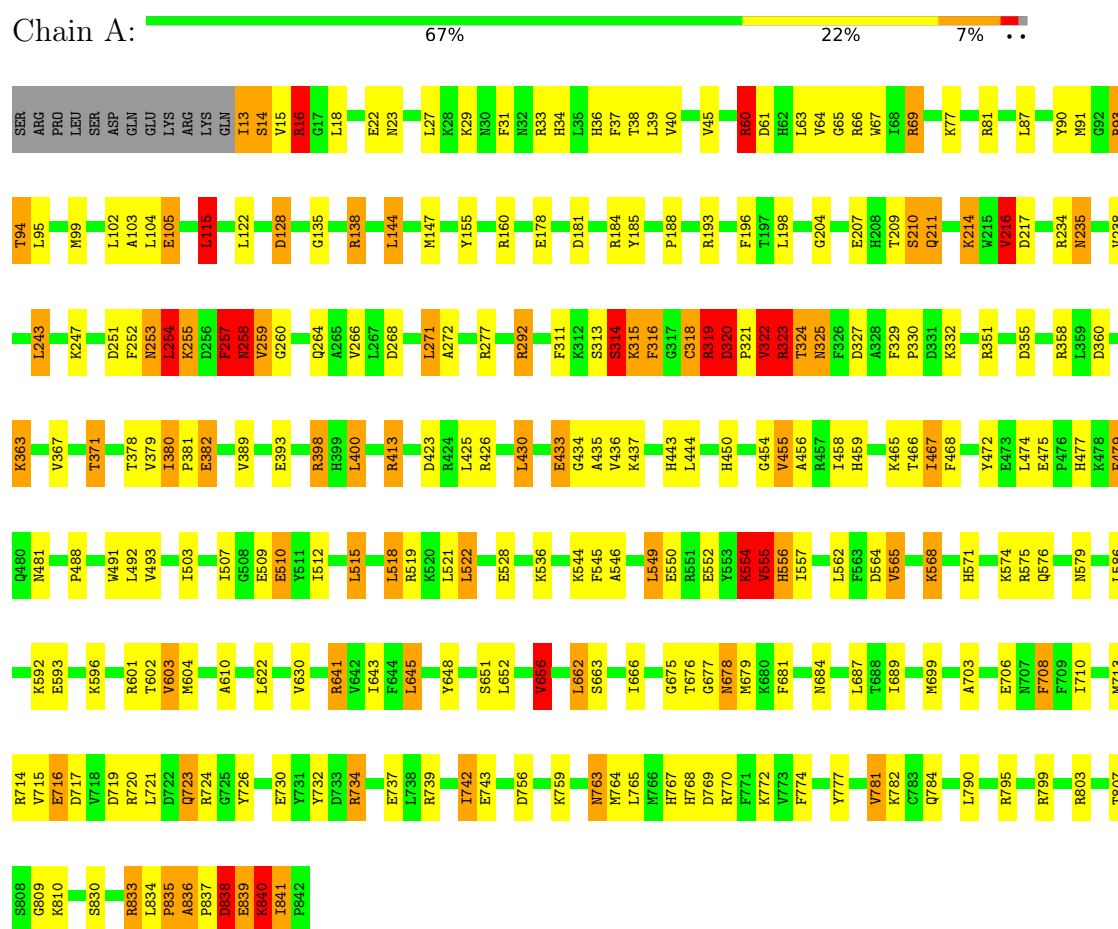
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	331	Total	O	0	0
			331	331		

3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

• Molecule 1: PROTEIN (GLYCOGEN PHOSPHORYLASE)



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	128.40 Å 128.40 Å 116.10 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.50 – 1.76	Depositor
% Data completeness (in resolution range)	99.3 (29.50-1.76)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.212 , 0.244	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7123	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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: PLP, CPB

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	0.48	1/6903 (0.0%)	0.98	41/9344 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	60	ARG	CZ-NH1	5.06	1.39	1.32

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	ARG	NE-CZ-NH2	-12.86	107.63	119.20
1	A	257	PHE	N-CA-C	7.90	119.99	110.19
1	A	554	LYS	N-CA-C	7.75	127.30	110.80
1	A	576	GLN	N-CA-C	-7.30	103.48	112.38
1	A	91	MET	N-CA-C	7.26	121.38	112.23
1	A	656	VAL	CB-CA-C	-6.91	102.69	112.22
1	A	630	VAL	N-CA-C	6.59	116.74	110.42
1	A	565	VAL	N-CA-C	6.56	118.23	108.46
1	A	105	GLU	N-CA-C	6.36	118.74	111.11
1	A	316	PHE	N-CA-C	-6.32	104.62	112.90
1	A	37	PHE	N-CA-C	6.19	119.00	111.82
1	A	272	ALA	N-CA-C	-6.14	104.86	112.90
1	A	556	HIS	N-CA-C	6.02	123.61	110.80
1	A	466	THR	N-CA-C	5.92	117.53	108.30
1	A	575	ARG	N-CA-C	5.84	119.69	111.52
1	A	767	HIS	N-CA-C	5.80	121.23	114.04
1	A	507	ILE	N-CA-C	5.79	118.46	112.90
1	A	455	VAL	N-CA-C	5.75	117.64	112.17
1	A	714	ARG	N-CA-C	-5.71	101.55	110.17
1	A	115	LEU	N-CA-C	-5.68	106.33	113.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	454	GLY	N-CA-C	-5.68	103.92	112.89
1	A	491	TRP	N-CA-C	5.67	119.90	113.15
1	A	60	ARG	CD-NE-CZ	5.66	132.33	124.40
1	A	675	GLY	N-CA-C	-5.62	105.68	112.48
1	A	135	GLY	N-CA-C	5.59	119.41	112.64
1	A	216	VAL	N-CA-C	5.57	117.68	108.99
1	A	60	ARG	NE-CZ-NH1	5.56	127.06	121.50
1	A	430	LEU	N-CA-C	-5.50	106.07	112.89
1	A	277	ARG	N-CA-C	5.48	117.06	111.14
1	A	666	ILE	N-CA-C	5.46	120.70	109.34
1	A	676	THR	N-CA-C	-5.44	108.41	114.62
1	A	684	ASN	N-CA-C	5.42	118.90	112.72
1	A	16	ARG	N-CA-C	-5.34	106.95	112.93
1	A	610	ALA	N-CA-C	-5.30	102.91	109.64
1	A	31	PHE	N-CA-C	-5.29	105.41	111.07
1	A	723	GLN	N-CA-C	-5.26	105.44	111.07
1	A	320	ASP	CA-C-N	-5.24	114.42	119.87
1	A	320	ASP	C-N-CA	-5.24	114.42	119.87
1	A	479	PHE	N-CA-C	5.16	117.97	109.76
1	A	677	GLY	N-CA-C	-5.14	106.56	112.73
1	A	602	THR	N-CA-C	-5.00	99.27	108.02

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	6749	0	6695	195	0
2	A	15	0	7	0	0
3	A	28	0	20	1	0
4	A	331	0	0	22	0
All	All	7123	0	6722	195	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 14.

All (195) 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:320:ASP:HB3	1:A:321:PRO:HD2	1.29	1.10
1:A:251:ASP:CG	1:A:252:PHE:H	1.67	1.01
1:A:320:ASP:HB3	1:A:321:PRO:CD	1.97	0.93
1:A:571:HIS:HB2	1:A:574:LYS:HG3	1.54	0.89
1:A:764:MET:HE2	1:A:769:ASP:HA	1.52	0.89
1:A:515:LEU:HD22	1:A:518:LEU:HD22	1.58	0.83
1:A:455:VAL:H	1:A:459:HIS:HD2	1.27	0.81
1:A:839:GLU:O	1:A:840:LYS:HG3	1.81	0.81
1:A:322:VAL:O	1:A:323:ARG:HB2	1.78	0.80
1:A:554:LYS:O	1:A:555:VAL:HG22	1.82	0.80
1:A:251:ASP:CG	1:A:252:PHE:N	2.40	0.79
1:A:568:LYS:HG3	1:A:574:LYS:HD2	1.64	0.78
1:A:413:ARG:HH11	1:A:413:ARG:HG2	1.48	0.78
1:A:319:ARG:HG2	1:A:320:ASP:H	1.49	0.77
1:A:319:ARG:CG	1:A:320:ASP:H	1.98	0.77
1:A:138:ARG:O	1:A:138:ARG:HD3	1.84	0.76
1:A:316:PHE:HE1	1:A:332:LYS:HZ2	1.35	0.75
1:A:678:ASN:HD22	1:A:679:MET:H	1.33	0.74
1:A:545:PHE:CZ	1:A:656:VAL:HG13	2.23	0.74
1:A:15:VAL:HB	4:A:1301:HOH:O	1.88	0.74
1:A:703:ALA:HA	1:A:807:THR:HG21	1.69	0.73
1:A:184:ARG:HE	1:A:185:TYR:HE1	1.37	0.72
1:A:66:ARG:HG3	1:A:837:PRO:HB3	1.70	0.72
1:A:601:ARG:HD2	4:A:1033:HOH:O	1.88	0.72
1:A:65:GLY:O	1:A:69:ARG:HG2	1.90	0.71
1:A:13:ILE:O	1:A:14:SER:HB2	1.92	0.69
1:A:235:ASN:H	1:A:235:ASN:HD22	1.39	0.69
1:A:795:ARG:O	1:A:799:ARG:HG3	1.92	0.68
1:A:764:MET:SD	4:A:1309:HOH:O	2.51	0.68
1:A:841:ILE:HG23	1:A:841:ILE:O	1.93	0.67
1:A:15:VAL:HG21	1:A:509:GLU:CD	2.20	0.67
1:A:380:ILE:HD12	3:A:940:CPB:H251	1.76	0.67
1:A:716:GLU:HG2	4:A:1310:HOH:O	1.93	0.67
1:A:323:ARG:O	1:A:325:ASN:N	2.30	0.65
1:A:257:PHE:O	1:A:258:ASN:HB2	1.95	0.65
1:A:703:ALA:CA	1:A:807:THR:HG21	2.26	0.64
1:A:360:ASP:OD1	1:A:363:LYS:HB2	1.97	0.64
1:A:319:ARG:N	1:A:319:ARG:HD3	2.11	0.64
1:A:713:MET:HB3	1:A:717:ASP:HB2	1.79	0.63
1:A:69:ARG:NH1	1:A:840:LYS:HG2	2.13	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:VAL:HG13	1:A:15:VAL:O	1.98	0.63
1:A:367:VAL:O	1:A:371:THR:HG23	1.99	0.62
1:A:160:ARG:HB2	1:A:243:LEU:HB3	1.80	0.62
1:A:764:MET:CE	1:A:769:ASP:HA	2.27	0.62
1:A:554:LYS:C	1:A:555:VAL:HG22	2.23	0.62
1:A:841:ILE:O	1:A:841:ILE:CG2	2.49	0.61
1:A:378:THR:OG1	1:A:380:ILE:HG12	2.01	0.61
1:A:316:PHE:HE1	1:A:332:LYS:NZ	1.98	0.61
1:A:739:ARG:O	1:A:743:GLU:HG3	2.01	0.60
1:A:550:GLU:O	1:A:554:LYS:HG2	2.01	0.60
1:A:678:ASN:ND2	1:A:679:MET:H	1.98	0.60
1:A:325:ASN:HA	4:A:1173:HOH:O	2.02	0.59
1:A:764:MET:HE1	4:A:1210:HOH:O	2.02	0.59
1:A:60:ARG:HD2	1:A:188:PRO:O	2.01	0.59
1:A:329:PHE:HB3	1:A:330:PRO:HD3	1.85	0.59
1:A:641:ARG:HG3	4:A:1215:HOH:O	2.02	0.59
1:A:493:VAL:CG2	1:A:512:ILE:HD12	2.33	0.58
1:A:719:ASP:O	1:A:723:GLN:HG2	2.04	0.58
1:A:16:ARG:HB3	1:A:105:GLU:HB3	1.86	0.57
1:A:455:VAL:H	1:A:459:HIS:CD2	2.17	0.57
1:A:34:HIS:HD2	1:A:38:THR:OG1	1.88	0.57
1:A:678:ASN:HD22	1:A:679:MET:N	2.02	0.57
1:A:413:ARG:HG2	1:A:413:ARG:NH1	2.15	0.56
1:A:15:VAL:HG12	4:A:1190:HOH:O	2.04	0.56
1:A:90:TYR:HE1	4:A:1041:HOH:O	1.88	0.56
1:A:60:ARG:HG2	1:A:60:ARG:HH11	1.70	0.56
1:A:15:VAL:CG1	4:A:1190:HOH:O	2.54	0.55
1:A:94:THR:CG2	4:A:1083:HOH:O	2.55	0.55
1:A:320:ASP:CB	1:A:321:PRO:CD	2.81	0.55
1:A:708:PHE:HB3	1:A:710:ILE:CD1	2.37	0.55
1:A:355:ASP:OD2	1:A:398:ARG:HD3	2.06	0.55
1:A:554:LYS:O	1:A:555:VAL:HG13	2.06	0.55
1:A:703:ALA:CB	1:A:807:THR:HG21	2.36	0.55
1:A:61:ASP:O	1:A:64:VAL:HG22	2.07	0.54
1:A:252:PHE:C	1:A:254:LEU:H	2.14	0.54
1:A:678:ASN:HD22	1:A:678:ASN:N	2.05	0.54
1:A:93:ARG:HH11	1:A:93:ARG:CG	2.20	0.54
1:A:467:ILE:HG23	1:A:468:PHE:CG	2.43	0.54
1:A:15:VAL:HG21	1:A:509:GLU:OE2	2.08	0.53
1:A:235:ASN:HA	1:A:833:ARG:HG3	1.90	0.53
1:A:546:ALA:HA	1:A:557:ILE:HD11	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:VAL:O	1:A:322:VAL:HG12	2.08	0.52
1:A:810:LYS:O	1:A:810:LYS:HG2	2.09	0.52
1:A:253:ASN:O	1:A:254:LEU:HB2	2.07	0.52
1:A:777:TYR:O	1:A:781:VAL:HG13	2.10	0.52
1:A:93:ARG:HH11	1:A:93:ARG:HG3	1.75	0.51
1:A:138:ARG:HD3	1:A:138:ARG:C	2.36	0.51
1:A:699:MET:HE2	1:A:708:PHE:HZ	1.76	0.51
1:A:592:LYS:HD2	1:A:592:LYS:O	2.12	0.50
1:A:313:SER:O	1:A:314:SER:C	2.54	0.50
1:A:763:ASN:HD22	1:A:763:ASN:C	2.19	0.50
1:A:319:ARG:HD3	1:A:319:ARG:H	1.77	0.50
1:A:708:PHE:HB3	1:A:710:ILE:HD12	1.93	0.50
1:A:475:GLU:HB3	1:A:477:HIS:CE1	2.47	0.50
1:A:423:ASP:O	1:A:426:ARG:HG3	2.12	0.50
1:A:434:GLY:O	1:A:435:ALA:C	2.55	0.50
1:A:268:ASP:O	1:A:271:LEU:HB2	2.11	0.50
1:A:367:VAL:O	1:A:371:THR:CG2	2.60	0.50
1:A:144:LEU:HD12	1:A:147:MET:CE	2.42	0.49
1:A:251:ASP:C	1:A:255:LYS:HB3	2.37	0.49
1:A:379:VAL:HG22	4:A:1104:HOH:O	2.13	0.49
1:A:292:ARG:HD2	1:A:292:ARG:O	2.13	0.49
1:A:703:ALA:HB2	1:A:807:THR:HG21	1.94	0.49
1:A:36:HIS:O	1:A:40:VAL:HA	2.13	0.49
1:A:314:SER:HB3	1:A:315:LYS:H	1.41	0.49
1:A:836:ALA:HB1	1:A:837:PRO:CD	2.43	0.49
1:A:648:TYR:HA	1:A:652:LEU:HD23	1.94	0.48
1:A:382:GLU:H	1:A:382:GLU:CD	2.17	0.48
1:A:181:ASP:HB3	1:A:184:ARG:NH1	2.29	0.48
1:A:316:PHE:C	1:A:318:CYS:H	2.21	0.48
1:A:65:GLY:O	1:A:69:ARG:CG	2.61	0.48
1:A:216:VAL:HG13	1:A:217:ASP:N	2.29	0.48
1:A:94:THR:HG22	4:A:1083:HOH:O	2.13	0.47
1:A:95:LEU:HD11	1:A:99:MET:HE3	1.94	0.47
1:A:450:HIS:HD2	4:A:1264:HOH:O	1.97	0.47
1:A:389:VAL:O	1:A:393:GLU:HB2	2.15	0.47
1:A:235:ASN:HD22	1:A:235:ASN:N	2.02	0.47
1:A:592:LYS:HG3	1:A:593:GLU:HG2	1.96	0.47
1:A:510:GLU:HB2	4:A:1256:HOH:O	2.13	0.47
1:A:13:ILE:HG22	4:A:1296:HOH:O	2.14	0.47
1:A:329:PHE:HD2	1:A:371:THR:HG21	1.80	0.47
1:A:834:LEU:O	1:A:836:ALA:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:519:ARG:HH11	1:A:519:ARG:HG2	1.80	0.46
1:A:836:ALA:HB1	1:A:837:PRO:HD2	1.97	0.46
1:A:67:TRP:HA	1:A:238:VAL:HB	1.98	0.46
1:A:678:ASN:ND2	1:A:679:MET:N	2.62	0.46
1:A:742:ILE:HD11	1:A:774:PHE:CZ	2.51	0.45
1:A:319:ARG:CG	1:A:320:ASP:N	2.69	0.45
1:A:389:VAL:HG13	1:A:400:LEU:HD11	1.97	0.45
1:A:458:ILE:HD11	1:A:715:VAL:HG21	1.98	0.45
1:A:519:ARG:HG2	1:A:519:ARG:NH1	2.31	0.45
1:A:128:ASP:OD1	1:A:651:SER:HB3	2.17	0.45
1:A:329:PHE:CD2	1:A:371:THR:HG21	2.52	0.45
1:A:835:PRO:O	1:A:836:ALA:O	2.35	0.45
1:A:604:MET:HB3	1:A:645:LEU:HD22	1.98	0.45
1:A:550:GLU:HA	1:A:554:LYS:HA	1.98	0.45
1:A:730:GLU:O	1:A:734:ARG:HG2	2.16	0.45
1:A:15:VAL:HG22	4:A:1026:HOH:O	2.16	0.45
1:A:493:VAL:HG21	1:A:512:ILE:HD12	1.99	0.45
1:A:252:PHE:C	1:A:254:LEU:N	2.74	0.44
1:A:579:ASN:C	1:A:579:ASN:HD22	2.24	0.44
1:A:433:GLU:CD	1:A:437:LYS:HD2	2.43	0.44
1:A:60:ARG:HG2	1:A:60:ARG:NH1	2.29	0.44
1:A:311:PHE:O	1:A:314:SER:HB2	2.17	0.44
1:A:663:SER:HB2	1:A:681:PHE:CG	2.52	0.44
1:A:732:TYR:CZ	1:A:739:ARG:HG3	2.53	0.44
1:A:726:TYR:OH	1:A:774:PHE:HB2	2.17	0.44
1:A:259:VAL:CG1	1:A:260:GLY:N	2.80	0.43
1:A:488:PRO:O	1:A:492:LEU:HB3	2.18	0.43
1:A:678:ASN:ND2	1:A:678:ASN:N	2.64	0.43
1:A:764:MET:HE3	4:A:1056:HOH:O	2.18	0.43
1:A:382:GLU:OE2	1:A:770:ARG:NH1	2.50	0.43
1:A:515:LEU:HB3	1:A:809:GLY:HA2	1.99	0.43
1:A:737:GLU:HG3	4:A:1160:HOH:O	2.17	0.43
1:A:450:HIS:HE1	4:A:1142:HOH:O	2.01	0.43
1:A:319:ARG:NE	1:A:320:ASP:N	2.66	0.43
1:A:549:LEU:HA	1:A:549:LEU:HD12	1.73	0.43
1:A:193:ARG:HB3	1:A:196:PHE:CD2	2.54	0.43
1:A:456:ALA:C	1:A:481:ASN:HD21	2.27	0.43
1:A:77:LYS:HA	1:A:77:LYS:HD2	1.87	0.43
1:A:209:THR:O	1:A:211:GLN:N	2.51	0.43
1:A:254:LEU:HD21	1:A:266:VAL:HG23	2.00	0.42
1:A:322:VAL:O	1:A:322:VAL:CG1	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:601:ARG:NH2	1:A:784:GLN:OE1	2.44	0.42
1:A:60:ARG:O	1:A:63:LEU:HD23	2.19	0.42
1:A:216:VAL:HG13	1:A:217:ASP:CG	2.44	0.42
1:A:319:ARG:O	1:A:320:ASP:HB2	2.19	0.42
1:A:325:ASN:C	1:A:327:ASP:H	2.28	0.42
1:A:204:GLY:HA2	1:A:217:ASP:O	2.20	0.42
1:A:381:PRO:HG2	1:A:382:GLU:OE1	2.19	0.42
1:A:562:LEU:HD21	1:A:662:LEU:HB2	2.00	0.42
1:A:467:ILE:CG2	1:A:468:PHE:N	2.83	0.42
1:A:689:ILE:HG23	1:A:689:ILE:O	2.19	0.42
1:A:207:GLU:OE1	1:A:214:LYS:NZ	2.51	0.41
1:A:144:LEU:HA	1:A:147:MET:HE3	2.02	0.41
1:A:255:LYS:O	1:A:257:PHE:N	2.54	0.41
1:A:472:TYR:HA	1:A:479:PHE:CE1	2.56	0.41
1:A:351:ARG:O	1:A:355:ASP:HB2	2.21	0.41
1:A:34:HIS:HE1	1:A:61:ASP:OD2	2.03	0.41
1:A:519:ARG:O	1:A:522:LEU:HB2	2.19	0.41
1:A:764:MET:CE	4:A:1056:HOH:O	2.68	0.41
1:A:433:GLU:OE1	1:A:433:GLU:CA	2.68	0.41
1:A:251:ASP:O	1:A:255:LYS:N	2.45	0.41
1:A:430:LEU:HD22	1:A:443:HIS:HB3	2.03	0.41
1:A:564:ASP:O	1:A:603:VAL:HA	2.21	0.41
1:A:69:ARG:NH1	1:A:838:ASP:O	2.54	0.41
1:A:81:ARG:HG2	1:A:155:TYR:HE1	1.86	0.41
1:A:115:LEU:HD12	1:A:115:LEU:HA	1.92	0.41
1:A:193:ARG:HB3	1:A:196:PHE:HD2	1.85	0.41
1:A:103:ALA:HA	4:A:1005:HOH:O	2.21	0.41
1:A:235:ASN:H	1:A:235:ASN:ND2	2.11	0.40
1:A:503:ILE:HG12	1:A:521:LEU:HD11	2.03	0.40
1:A:292:ARG:HD2	1:A:292:ARG:C	2.46	0.40
1:A:803:ARG:NH1	1:A:803:ARG:HG3	2.37	0.40
1:A:803:ARG:HG3	1:A:803:ARG:HH11	1.86	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	828/842 (98%)	768 (93%)	41 (5%)	19 (2%)	5 0

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	14	SER
1	A	210	SER
1	A	254	LEU
1	A	258	ASN
1	A	259	VAL
1	A	320	ASP
1	A	323	ARG
1	A	324	THR
1	A	555	VAL
1	A	556	HIS
1	A	839	GLU
1	A	840	LYS
1	A	554	LYS
1	A	835	PRO
1	A	836	ALA
1	A	838	ASP
1	A	314	SER
1	A	319	ARG
1	A	322	VAL

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	719/731 (98%)	610 (85%)	109 (15%)	3 0

All (109) 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	60	ARG
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
1	A	247	LYS
1	A	253	ASN
1	A	254	LEU
1	A	255	LYS
1	A	257	PHE
1	A	258	ASN
1	A	264	GLN
1	A	271	LEU
1	A	292	ARG
1	A	314	SER
1	A	315	LYS
1	A	318	CYS

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Mol	Chain	Res	Type
1	A	319	ARG
1	A	320	ASP
1	A	322	VAL
1	A	323	ARG
1	A	324	THR
1	A	325	ASN
1	A	358	ARG
1	A	363	LYS
1	A	371	THR
1	A	380	ILE
1	A	382	GLU
1	A	398	ARG
1	A	400	LEU
1	A	413	ARG
1	A	425	LEU
1	A	433	GLU
1	A	436	VAL
1	A	444	LEU
1	A	465	LYS
1	A	467	ILE
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	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

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Mol	Chain	Res	Type
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	756	ASP
1	A	759	LYS
1	A	763	ASN
1	A	765	LEU
1	A	768	HIS
1	A	772	LYS
1	A	781	VAL
1	A	782	LYS
1	A	790	LEU
1	A	830	SER
1	A	833	ARG
1	A	838	ASP
1	A	840	LYS
1	A	841	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	72	GLN
1	A	208	HIS
1	A	235	ASN
1	A	250	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	560	ASN
1	A	566	GLN

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Mol	Chain	Res	Type
1	A	579	ASN
1	A	678	ASN
1	A	763	ASN
1	A	767	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	$\# Z > 2$
2	PLP	A	999	1	15,15,16	1.50	3 (20%)	21,22,23	1.07	1 (4%)
3	CPB	A	940	-	31,31,31	2.26	15 (48%)	42,46,46	1.91	10 (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	PLP	A	999	1	-	1/6/6/8	0/1/1/1
3	CPB	A	940	-	-	0/8/21/21	0/4/4/4

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	940	CPB	C14-C13	5.09	1.57	1.53
3	A	940	CPB	C21-C22	4.41	1.45	1.39
2	A	999	PLP	C5A-C5	3.42	1.59	1.50
3	A	940	CPB	C21-C2	3.29	1.54	1.47
3	A	940	CPB	C6-C7	2.99	1.43	1.38
3	A	940	CPB	C7-C8	2.83	1.44	1.40
3	A	940	CPB	C26-C21	2.80	1.44	1.39
3	A	940	CPB	C25-C24	2.78	1.44	1.38
2	A	999	PLP	C4A-C4	2.66	1.56	1.51
3	A	940	CPB	C10-C4	2.63	1.53	1.46
3	A	940	CPB	O1-C9	2.60	1.42	1.38
3	A	940	CPB	C12-C13	2.49	1.55	1.52
3	A	940	CPB	O5-C5	2.34	1.41	1.36
3	A	940	CPB	C25-C26	2.33	1.42	1.38
3	A	940	CPB	C24-C23	2.23	1.42	1.38
3	A	940	CPB	C8-C14	2.19	1.55	1.52
2	A	999	PLP	C3-C2	-2.12	1.38	1.41
3	A	940	CPB	C12-N1	2.08	1.49	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	940	CPB	O1-C2-C21	6.67	118.50	110.65
3	A	940	CPB	C9-O1-C2	4.33	124.88	120.04
3	A	940	CPB	O1-C9-C8	3.54	119.57	116.10
3	A	940	CPB	O3-C13-C14	3.09	115.97	110.81
3	A	940	CPB	C15-C14-C13	2.97	112.05	109.35
3	A	940	CPB	O3-C13-C12	-2.76	104.38	109.66
2	A	999	PLP	O3P-P-O1P	2.46	120.40	110.83
3	A	940	CPB	C16-N1-C12	2.20	112.27	110.06
3	A	940	CPB	C7-C8-C9	2.17	118.61	116.47
3	A	940	CPB	C15-C14-C8	-2.17	107.59	113.40
3	A	940	CPB	C5-C10-C9	2.17	120.51	117.04

There are no chirality outliers.

All (1) torsion outliers are listed below:

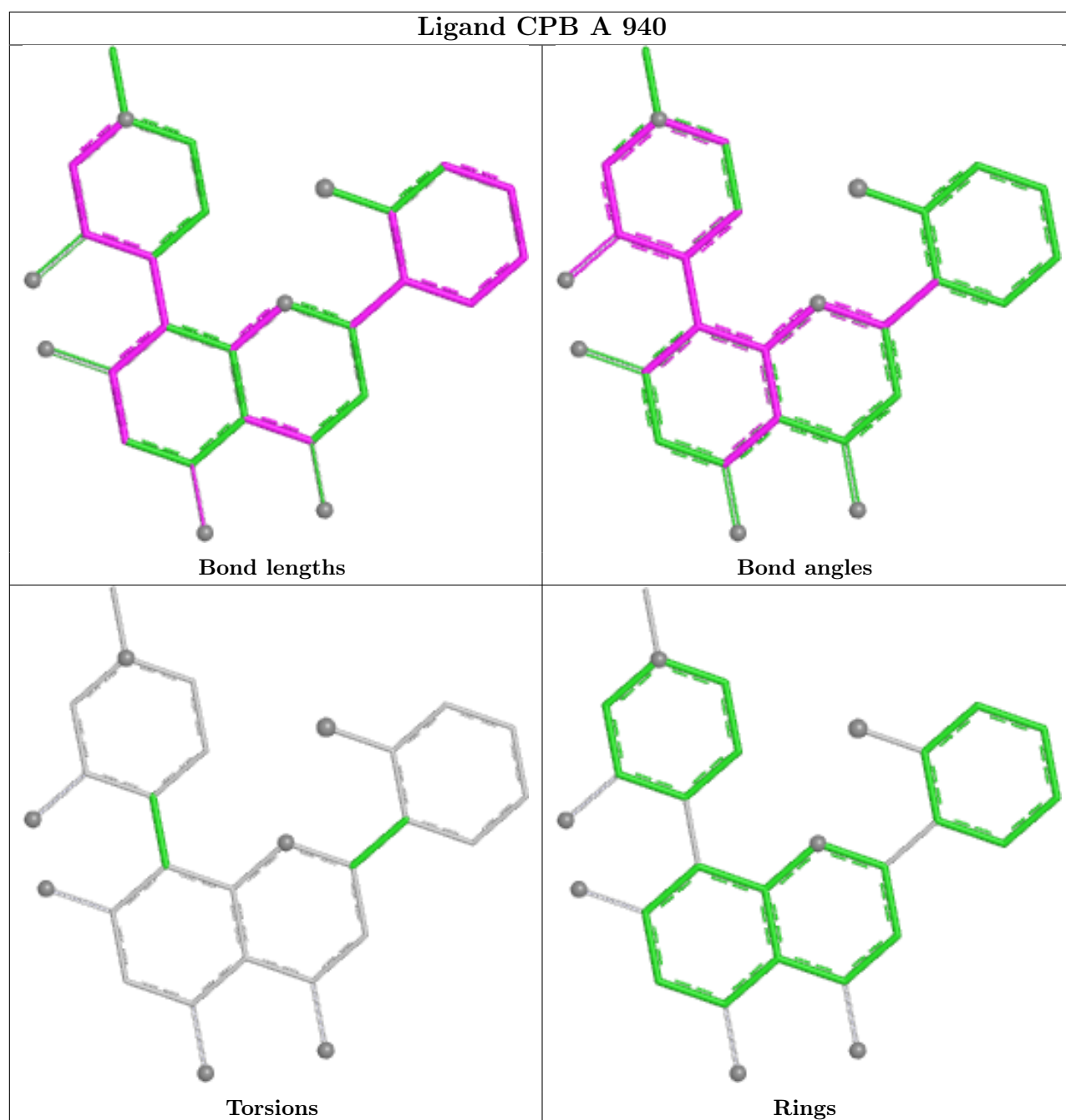
Mol	Chain	Res	Type	Atoms
2	A	999	PLP	C6-C5-C5A-O4P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	940	CPB	1	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

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.