



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:47 PM UTC

PDB ID : 1RDY / pdb_00001rdy
Title : T-STATE STRUCTURE OF THE ARG 243 TO ALA MUTANT OF PIG KIDNEY FRUCTOSE 1,6-BISPHOSPHATASE EXPRESSED IN E. COLI
Authors : Stec, B.; Abraham, R.; Giroux, E.; Kantrowitz, E.R.
Deposited on : 1996-05-17
Resolution : 2.20 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

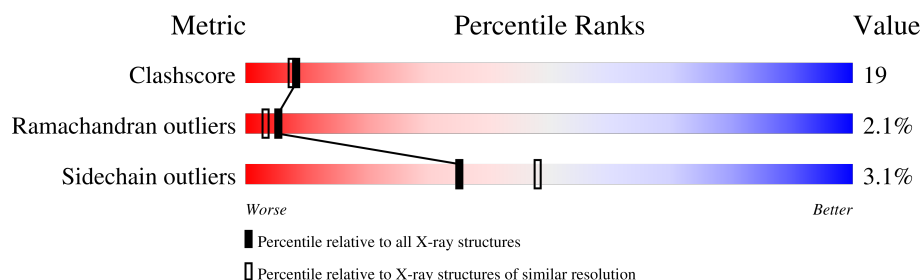
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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	337	
1	B	337	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5322 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.

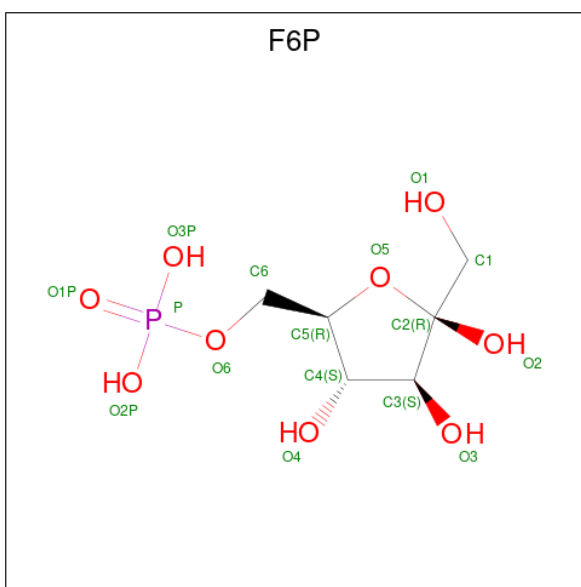
- Molecule 1 is a protein called FRUCTOSE 1,6-BISPHOSPHATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2566	1630	431	490	15			
1	B	337	Total	C	N	O	S	0	0	0
			2566	1630	431	490	15			

There are 8 discrepancies between the modelled and reference sequences:

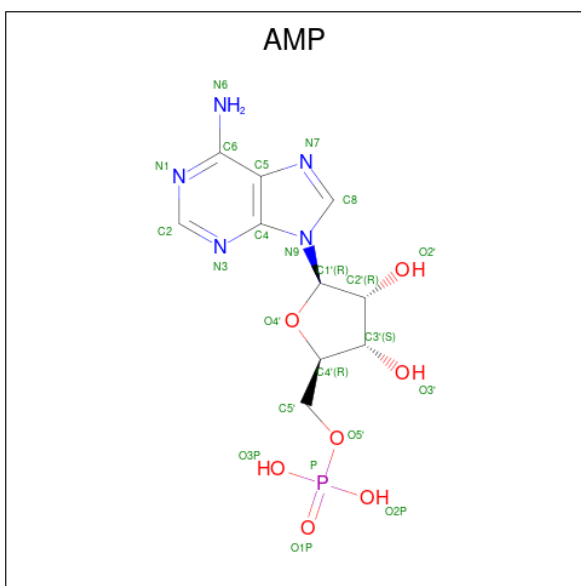
Chain	Residue	Modelled	Actual	Comment	Reference
A	20	GLN	GLU	conflict	UNP P00636
A	96	THR	SER	conflict	UNP P00636
A	199	ASN	ASP	conflict	UNP P00636
A	243	ALA	ARG	engineered mutation	UNP P00636
B	20	GLN	GLU	conflict	UNP P00636
B	96	THR	SER	conflict	UNP P00636
B	199	ASN	ASP	conflict	UNP P00636
B	243	ALA	ARG	engineered mutation	UNP P00636

- Molecule 2 is 6-O-phosphono-beta-D-fructofuranose (CCD ID: F6P) (formula: C₆H₁₃O₉P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			16	6	9	1		
2	B	1	Total	C	O	P	0	0
			16	6	9	1		

- Molecule 3 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 4 is water.

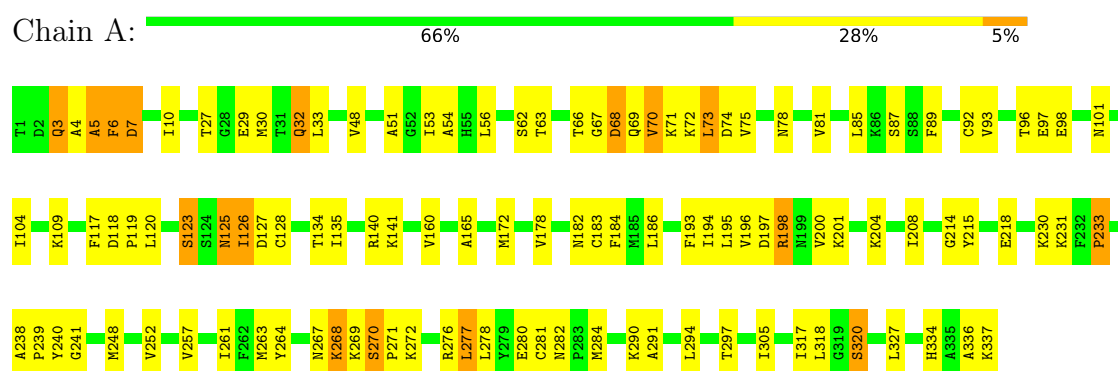
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	49	Total 49	O 49	0	0
4	B	63	Total 63	O 63	0	0

3 Residue-property plots

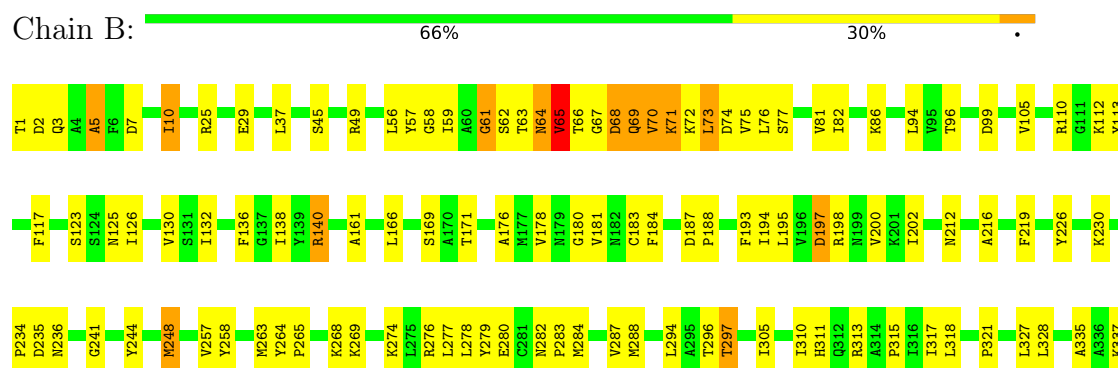
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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: FRUCTOSE 1,6-BISPHOSPHATASE



• Molecule 1: FRUCTOSE 1,6-BISPHOSPHATASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	61.20Å 166.75Å 79.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.20)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.198 , 0.234	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5322	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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: AMP, F6P

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	0.76	3/2609 (0.1%)	1.21	27/3531 (0.8%)
1	B	0.62	1/2609 (0.0%)	1.23	30/3531 (0.8%)
All	All	0.69	4/5218 (0.1%)	1.22	57/7062 (0.8%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	ILE	N-CA	19.70	1.74	1.46
1	B	234	PRO	C-N	14.08	1.56	1.33
1	A	5	ALA	C-N	8.25	1.45	1.33
1	A	7	ASP	N-CA	6.92	1.53	1.46

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	234	PRO	O-C-N	-16.08	101.48	122.22
1	B	335	ALA	N-CA-CB	13.23	130.46	109.95
1	A	334	HIS	CA-C-O	-11.69	106.20	119.41
1	B	234	PRO	CA-C-N	9.51	141.34	122.31
1	B	234	PRO	C-N-CA	9.51	141.34	122.31
1	A	68	ASP	O-C-N	9.39	132.09	123.26
1	A	125	ASN	CA-C-N	-9.31	105.32	120.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	125	ASN	C-N-CA	-9.31	105.32	120.86
1	B	235	ASP	CA-C-N	8.80	134.60	122.19
1	B	235	ASP	C-N-CA	8.80	134.60	122.19
1	A	126	ILE	N-CA-CB	-8.62	93.72	110.78
1	A	6	PHE	N-CA-C	8.51	123.71	109.76
1	B	264	TYR	CA-C-N	8.44	130.39	119.84
1	B	264	TYR	C-N-CA	8.44	130.39	119.84
1	B	278	LEU	N-CA-C	8.35	120.39	111.28
1	A	123	SER	CB-CA-C	-7.98	94.95	110.11
1	B	277	LEU	N-CA-C	7.94	119.62	110.97
1	A	126	ILE	N-CA-C	7.43	122.59	111.89
1	A	123	SER	CA-C-O	-7.41	109.41	119.12
1	A	334	HIS	CA-C-N	-6.97	109.88	121.80
1	A	334	HIS	C-N-CA	-6.97	109.88	121.80
1	B	140	ARG	CA-CB-CG	6.95	128.01	114.10
1	A	4	ALA	N-CA-C	-6.94	105.16	112.93
1	B	257	VAL	N-CA-C	6.93	116.93	110.42
1	A	233	PRO	CA-C-N	6.82	126.62	119.05
1	A	233	PRO	C-N-CA	6.82	126.62	119.05
1	B	5	ALA	O-C-N	6.70	131.50	122.59
1	A	278	LEU	N-CA-C	6.43	118.83	111.11
1	A	257	VAL	N-CA-C	6.37	117.00	110.82
1	B	335	ALA	N-CA-C	-6.33	99.69	109.76
1	B	236	ASN	O-C-N	6.20	130.72	122.41
1	B	68	ASP	N-CA-C	6.14	117.77	111.14
1	A	277	LEU	N-CA-C	6.11	117.82	111.03
1	A	264	TYR	CA-C-N	6.00	126.46	120.52
1	A	264	TYR	C-N-CA	6.00	126.46	120.52
1	A	118	ASP	CA-C-N	5.87	125.35	119.24
1	A	118	ASP	C-N-CA	5.87	125.35	119.24
1	B	110	ARG	N-CA-C	5.85	118.31	110.35
1	B	248	MET	N-CA-C	5.80	118.08	111.11
1	A	68	ASP	N-CA-C	5.80	117.07	107.73
1	B	70	VAL	O-C-N	5.80	129.57	122.72
1	B	56	LEU	N-CA-C	-5.79	105.10	111.82
1	A	320	SER	CA-C-N	5.61	125.28	119.05
1	A	320	SER	C-N-CA	5.61	125.28	119.05
1	B	70	VAL	N-CA-C	5.60	116.83	109.21
1	B	68	ASP	O-C-N	-5.47	116.19	122.09
1	B	241	GLY	N-CA-C	-5.46	104.34	113.02
1	B	279	TYR	N-CA-C	5.45	119.90	112.88
1	B	235	ASP	N-CA-CB	-5.43	102.09	110.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	187	ASP	N-CA-C	-5.28	98.15	109.81
1	A	140	ARG	N-CA-C	-5.26	100.67	109.24
1	A	27	THR	N-CA-C	-5.25	106.83	113.18
1	B	297	THR	N-CA-C	-5.13	107.07	113.38
1	B	58	GLY	N-CA-C	5.12	120.95	113.48
1	B	140	ARG	N-CA-C	-5.10	101.40	109.76
1	B	180	GLY	N-CA-C	5.09	117.74	110.46
1	A	214	GLY	N-CA-C	-5.03	107.55	113.99

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	123	SER	Mainchain

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	2566	0	2620	122	0
1	B	2566	0	2620	88	0
2	A	16	0	11	0	0
2	B	16	0	11	0	0
3	A	23	0	12	1	0
3	B	23	0	12	5	0
4	A	49	0	0	0	0
4	B	63	0	0	1	0
All	All	5322	0	5286	201	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 19.

All (201) 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:126:ILE:N	1:A:126:ILE:CA	1.74	1.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:ARG:HB2	1:A:198:ARG:NH1	1.61	1.15
1:B:112:LYS:NZ	3:B:339:AMP:O3P	1.86	1.08
1:A:198:ARG:HB2	1:A:198:ARG:HH11	0.94	1.05
1:A:198:ARG:NH1	1:A:198:ARG:CB	2.21	1.03
1:A:62:SER:HA	1:A:66:THR:OG1	1.60	1.01
1:B:72:LYS:O	1:B:74:ASP:N	1.94	1.00
1:A:198:ARG:HH11	1:A:198:ARG:CB	1.76	0.95
1:A:56:LEU:HA	1:A:66:THR:HG23	1.49	0.95
1:A:62:SER:HB3	1:A:66:THR:HB	1.47	0.94
1:A:215:TYR:HE1	1:A:269:LYS:NZ	1.64	0.93
1:A:337:LYS:HD3	1:A:337:LYS:H	1.33	0.93
1:A:125:ASN:C	1:A:126:ILE:CA	2.45	0.90
1:A:248:MET:HE2	1:A:284:MET:HG3	1.53	0.89
1:A:126:ILE:N	1:A:126:ILE:CB	2.39	0.84
1:A:32:GLN:HE21	1:A:32:GLN:HA	1.40	0.84
1:B:1:THR:HG22	1:B:2:ASP:H	1.43	0.82
1:B:65:VAL:HG22	1:B:66:THR:N	1.94	0.81
1:B:72:LYS:C	1:B:74:ASP:N	2.38	0.80
1:B:72:LYS:C	1:B:74:ASP:H	1.91	0.79
1:A:62:SER:HB3	1:A:66:THR:CB	2.13	0.77
1:A:69:GLN:HG2	1:A:69:GLN:O	1.84	0.77
1:B:96:THR:HG23	1:B:99:ASP:HB2	1.66	0.76
1:A:3:GLN:HA	1:A:3:GLN:HE21	1.50	0.76
1:B:64:ASN:O	1:B:67:GLY:N	2.19	0.76
1:A:32:GLN:HA	1:A:32:GLN:NE2	1.98	0.75
1:A:218:GLU:HB3	1:A:268:LYS:HB2	1.69	0.75
1:B:1:THR:HG22	1:B:2:ASP:N	2.00	0.75
1:A:267:ASN:HD21	1:A:270:SER:HB2	1.52	0.75
1:A:78:ASN:OD1	1:A:96:THR:HG21	1.87	0.73
1:A:215:TYR:HE1	1:A:269:LYS:HZ1	0.81	0.73
1:A:3:GLN:HA	1:A:3:GLN:NE2	2.03	0.71
1:B:71:LYS:O	1:B:71:LYS:HG3	1.91	0.70
1:B:112:LYS:HZ3	3:B:339:AMP:P	2.15	0.70
1:B:112:LYS:NZ	3:B:339:AMP:P	2.64	0.70
1:B:337:LYS:OXT	1:B:337:LYS:HG3	1.90	0.70
1:A:182:ASN:HD22	1:A:198:ARG:HA	1.56	0.69
1:A:267:ASN:O	1:A:268:LYS:C	2.35	0.69
1:A:72:LYS:O	1:A:75:VAL:N	2.25	0.69
1:B:63:THR:O	1:B:64:ASN:HB2	1.92	0.67
1:B:1:THR:CG2	1:B:2:ASP:H	2.09	0.66
1:A:72:LYS:HB3	1:A:74:ASP:OD1	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:LYS:O	1:A:74:ASP:N	2.29	0.66
1:A:276:ARG:HD3	1:A:276:ARG:N	2.11	0.65
1:A:198:ARG:NH1	1:A:198:ARG:HB3	2.09	0.65
1:B:297:THR:HB	1:B:315:PRO:HG2	1.77	0.65
1:A:267:ASN:C	1:A:268:LYS:O	2.39	0.64
1:A:96:THR:HG22	1:A:98:GLU:H	1.61	0.64
1:A:297:THR:HG21	1:A:305:ILE:HD11	1.79	0.64
1:B:10:ILE:O	1:B:10:ILE:HG23	1.98	0.64
1:B:230:LYS:HA	1:B:230:LYS:HE2	1.79	0.64
1:A:182:ASN:HD21	1:A:198:ARG:HG3	1.62	0.64
1:A:317:ILE:HD13	1:A:327:LEU:HD23	1.80	0.63
1:B:212:ASN:HB2	1:B:244:TYR:CE2	2.34	0.62
1:A:267:ASN:O	1:A:267:ASN:ND2	2.32	0.62
1:A:198:ARG:CB	1:A:198:ARG:CZ	2.76	0.62
1:A:56:LEU:O	1:A:66:THR:HG21	2.00	0.62
1:A:186:LEU:HB2	1:A:193:PHE:CE1	2.35	0.61
1:A:119:PRO:HA	1:A:134:THR:HG23	1.82	0.60
1:A:337:LYS:HD3	1:A:337:LYS:N	2.08	0.60
1:B:72:LYS:O	1:B:73:LEU:C	2.44	0.60
1:A:89:PHE:CD1	1:A:109:LYS:HA	2.35	0.60
1:B:113:TYR:HE1	1:B:140:ARG:HG2	1.67	0.60
1:A:62:SER:CB	1:A:66:THR:HB	2.25	0.59
1:A:267:ASN:O	1:A:268:LYS:O	2.20	0.59
1:B:72:LYS:O	1:B:75:VAL:N	2.35	0.59
1:A:10:ILE:HD12	1:B:59:ILE:HB	1.83	0.59
1:A:268:LYS:O	1:A:269:LYS:HB3	2.01	0.59
1:A:3:GLN:HE21	1:A:3:GLN:CA	2.15	0.59
1:A:269:LYS:O	1:A:270:SER:HB2	2.04	0.58
1:A:337:LYS:H	1:A:337:LYS:CD	2.14	0.58
1:A:128:CYS:SG	1:B:258:TYR:HE2	2.26	0.58
1:B:65:VAL:HG22	1:B:66:THR:H	1.66	0.58
1:B:29:GLU:HB3	3:B:339:AMP:O3P	2.04	0.58
1:B:64:ASN:O	1:B:65:VAL:C	2.47	0.57
1:A:53:ILE:HD12	1:A:56:LEU:HD22	1.87	0.57
1:A:125:ASN:O	1:A:126:ILE:CA	2.53	0.57
1:B:112:LYS:NZ	3:B:339:AMP:O1P	2.37	0.57
1:A:3:GLN:O	1:A:6:PHE:CZ	2.58	0.56
1:A:127:ASP:HB2	1:B:258:TYR:OH	2.04	0.56
1:B:248:MET:HE2	1:B:284:MET:HG3	1.87	0.56
1:B:296:THR:HG21	1:B:328:LEU:HD21	1.87	0.56
1:B:1:THR:CG2	1:B:2:ASP:N	2.67	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:ALA:HA	1:B:219:PHE:CD2	2.40	0.56
1:A:267:ASN:ND2	1:A:270:SER:HB2	2.20	0.55
1:A:198:ARG:HB3	1:A:198:ARG:CZ	2.37	0.55
1:A:62:SER:CB	1:A:66:THR:CB	2.81	0.55
1:B:113:TYR:CE1	1:B:140:ARG:HG2	2.42	0.55
1:B:123:SER:HA	1:B:126:ILE:HG23	1.88	0.55
1:B:297:THR:HG21	1:B:305:ILE:HD11	1.89	0.54
1:B:138:ILE:HD12	1:B:161:ALA:HB3	1.89	0.54
1:A:54:ALA:HA	1:B:194:ILE:CD1	2.37	0.54
1:B:226:TYR:HB2	1:B:327:LEU:HD13	1.90	0.54
1:A:317:ILE:HD13	1:A:327:LEU:CD2	2.37	0.54
1:B:65:VAL:CG2	1:B:66:THR:N	2.65	0.54
1:B:68:ASP:OD1	1:B:68:ASP:C	2.51	0.54
1:B:294:LEU:HD12	1:B:321:PRO:HA	1.87	0.54
1:B:183:CYS:SG	1:B:200:VAL:HG21	2.49	0.53
1:A:72:LYS:O	1:A:73:LEU:C	2.51	0.53
1:A:182:ASN:ND2	1:A:198:ARG:HG3	2.23	0.52
1:A:269:LYS:O	1:A:269:LYS:HG2	2.08	0.52
1:B:96:THR:CG2	1:B:99:ASP:HB2	2.38	0.52
1:A:252:VAL:HG11	1:A:284:MET:SD	2.50	0.52
1:A:184:PHE:HB3	1:A:193:PHE:HB3	1.90	0.52
1:A:29:GLU:HB3	3:A:339:AMP:O1P	2.10	0.52
1:B:337:LYS:OXT	1:B:337:LYS:CG	2.57	0.52
1:B:5:ALA:CB	1:B:7:ASP:OD1	2.58	0.51
1:B:123:SER:O	1:B:126:ILE:HG12	2.10	0.51
1:A:182:ASN:ND2	1:A:198:ARG:HA	2.24	0.51
1:A:62:SER:HA	1:A:66:THR:CB	2.40	0.51
1:A:337:LYS:N	1:A:337:LYS:CD	2.73	0.51
1:B:45:SER:O	1:B:49:ARG:HD3	2.11	0.51
1:B:61:GLY:HA2	4:B:352:HOH:O	2.11	0.51
1:B:68:ASP:OD1	1:B:68:ASP:O	2.29	0.51
1:B:166:LEU:O	1:B:171:THR:HA	2.11	0.51
1:B:5:ALA:HB3	1:B:7:ASP:OD1	2.11	0.51
1:B:202:ILE:HD12	1:B:288:MET:HE1	1.92	0.50
1:A:263:MET:HG2	1:A:317:ILE:HG23	1.93	0.50
1:A:269:LYS:O	1:A:270:SER:CB	2.58	0.50
1:B:183:CYS:HB2	1:B:197:ASP:HB2	1.93	0.50
1:A:62:SER:CA	1:A:66:THR:OG1	2.48	0.50
1:A:276:ARG:HD3	1:A:276:ARG:H	1.74	0.50
1:A:135:ILE:HD13	1:A:284:MET:HE2	1.94	0.50
1:A:196:VAL:HG12	1:B:57:TYR:CZ	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:LEU:HD21	1:B:198:ARG:HG3	1.93	0.50
1:A:204:LYS:O	1:A:320:SER:HB3	2.12	0.49
1:A:56:LEU:O	1:A:66:THR:CG2	2.60	0.49
1:A:89:PHE:HA	1:A:109:LYS:O	2.13	0.49
1:B:184:PHE:HB3	1:B:193:PHE:HB3	1.95	0.49
1:B:283:PRO:O	1:B:287:VAL:HG23	2.13	0.49
1:B:96:THR:HB	1:B:117:PHE:CZ	2.48	0.49
1:A:201:LYS:HA	1:A:291:ALA:O	2.13	0.48
1:A:277:LEU:HA	1:A:281:CYS:HB2	1.96	0.48
1:B:310:ILE:HG13	1:B:311:HIS:CD2	2.48	0.48
1:B:65:VAL:HG13	1:B:66:THR:H	1.79	0.48
1:B:318:LEU:C	1:B:318:LEU:HD12	2.38	0.47
1:B:282:ASN:HB2	1:B:283:PRO:HD3	1.96	0.47
1:B:71:LYS:HE3	1:B:76:LEU:HD22	1.97	0.47
1:A:48:VAL:HG12	1:B:169:SER:O	2.14	0.47
1:A:67:GLY:O	1:A:68:ASP:HB2	2.14	0.47
1:A:72:LYS:O	1:A:74:ASP:OD1	2.32	0.47
1:A:3:GLN:C	1:A:5:ALA:H	2.23	0.47
1:A:7:ASP:CG	1:A:198:ARG:NH2	2.73	0.47
1:B:280:GLU:O	1:B:284:MET:HG2	2.15	0.47
1:A:183:CYS:SG	1:A:200:VAL:HG21	2.56	0.46
1:B:71:LYS:O	1:B:71:LYS:CG	2.62	0.46
1:A:70:VAL:CG1	1:A:71:LYS:HG3	2.46	0.46
1:A:165:ALA:HA	1:A:172:MET:O	2.16	0.46
1:A:3:GLN:O	1:A:6:PHE:CE1	2.69	0.45
1:A:10:ILE:HG23	1:A:10:ILE:O	2.17	0.45
1:A:261:ILE:HD11	1:A:317:ILE:CG2	2.46	0.45
1:A:183:CYS:HB2	1:A:197:ASP:HB2	1.99	0.45
1:A:56:LEU:CA	1:A:66:THR:HG23	2.36	0.45
1:A:93:VAL:HG22	1:A:104:ILE:HG13	1.99	0.45
1:B:176:ALA:HB2	1:B:181:VAL:HG22	1.99	0.45
1:B:276:ARG:HA	1:B:313:ARG:HA	1.98	0.45
1:A:10:ILE:CD1	1:B:59:ILE:HB	2.47	0.45
1:A:231:LYS:C	1:A:233:PRO:HD3	2.42	0.45
1:A:126:ILE:N	1:A:126:ILE:HB	2.26	0.44
1:B:125:ASN:O	1:B:130:VAL:HG12	2.17	0.44
1:B:226:TYR:CB	1:B:327:LEU:HD13	2.47	0.44
1:A:62:SER:CA	1:A:66:THR:CB	2.95	0.44
1:A:117:PHE:HA	1:A:135:ILE:O	2.17	0.44
1:B:73:LEU:O	1:B:77:SER:OG	2.35	0.44
1:B:86:LYS:HG2	1:B:105:VAL:HG22	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:SER:HB3	1:A:66:THR:CG2	2.48	0.44
1:A:73:LEU:HD23	1:A:73:LEU:HA	1.77	0.43
1:A:337:LYS:OXT	1:A:337:LYS:HG2	2.19	0.43
1:A:73:LEU:HB3	1:A:120:LEU:CD2	2.48	0.43
1:A:51:ALA:HA	1:B:188:PRO:HD2	2.01	0.43
1:A:238:ALA:HA	1:A:239:PRO:HD3	1.89	0.43
1:B:66:THR:O	1:B:66:THR:HG22	2.19	0.43
1:A:30:MET:O	1:A:33:LEU:HB3	2.19	0.43
1:A:184:PHE:HA	1:A:194:ILE:O	2.18	0.42
1:B:274:LYS:O	1:B:313:ARG:NH1	2.52	0.42
1:A:92:CYS:SG	1:A:141:LYS:HD2	2.58	0.42
1:A:272:LYS:HD2	1:A:272:LYS:N	2.34	0.42
1:B:248:MET:CE	1:B:284:MET:HG3	2.48	0.42
1:B:25:ARG:HH11	1:B:25:ARG:HG3	1.83	0.42
1:A:70:VAL:CG1	1:A:71:LYS:N	2.83	0.42
1:B:61:GLY:O	1:B:63:THR:N	2.52	0.42
1:A:70:VAL:HG13	1:A:71:LYS:HG3	2.02	0.42
1:A:69:GLN:O	1:A:69:GLN:CG	2.49	0.42
1:A:270:SER:H	1:A:271:PRO:HD3	1.84	0.42
1:B:69:GLN:HE21	1:B:69:GLN:HB3	1.74	0.41
1:A:267:ASN:ND2	1:A:269:LYS:O	2.53	0.41
1:A:184:PHE:CE1	1:A:195:LEU:HD13	2.56	0.41
1:B:81:VAL:HG11	1:B:117:PHE:CE2	2.56	0.41
1:A:10:ILE:HB	1:B:57:TYR:O	2.21	0.41
1:A:81:VAL:O	1:A:85:LEU:HG	2.20	0.41
1:A:208:ILE:HA	1:A:241:GLY:O	2.21	0.41
1:B:82:ILE:HG23	1:B:94:LEU:HD12	2.02	0.41
1:B:288:MET:HG3	1:B:318:LEU:HB2	2.03	0.41
1:A:282:ASN:HD22	1:A:282:ASN:HA	1.68	0.41
1:B:66:THR:O	1:B:66:THR:CG2	2.68	0.40
1:B:263:MET:HG2	1:B:317:ILE:HG23	2.03	0.40
1:A:230:LYS:HB3	1:A:240:TYR:HB2	2.02	0.40
1:A:276:ARG:O	1:A:280:GLU:HB2	2.21	0.40
1:A:290:LYS:HA	1:A:290:LYS:HD3	1.83	0.40
1:A:294:LEU:O	1:A:318:LEU:HA	2.21	0.40
1:B:37:LEU:HD21	1:B:136:PHE:CD1	2.56	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	335/337 (99%)	306 (91%)	22 (7%)	7 (2%)	5	3
1	B	335/337 (99%)	312 (93%)	16 (5%)	7 (2%)	5	3
All	All	670/674 (99%)	618 (92%)	38 (6%)	14 (2%)	5	3

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	63	THR
1	A	73	LEU
1	A	270	SER
1	B	62	SER
1	B	64	ASN
1	B	65	VAL
1	B	73	LEU
1	B	269	LYS
1	A	268	LYS
1	A	178	VAL
1	B	61	GLY
1	B	268	LYS
1	A	87	SER
1	A	336	ALA

5.3.2 Protein sidechains [i](#)

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	278/278 (100%)	271 (98%)	7 (2%)	42	56
1	B	278/278 (100%)	268 (96%)	10 (4%)	31	42
All	All	556/556 (100%)	539 (97%)	17 (3%)	35	48

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	32	GLN
1	A	70	VAL
1	A	97	GLU
1	A	101	ASN
1	A	160	VAL
1	A	198	ARG
1	B	3	GLN
1	B	10	ILE
1	B	65	VAL
1	B	69	GLN
1	B	70	VAL
1	B	71	LYS
1	B	132	ILE
1	B	178	VAL
1	B	197	ASP
1	B	265	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	20	GLN
1	A	32	GLN
1	A	35	ASN
1	A	182	ASN
1	A	199	ASN
1	A	267	ASN
1	A	282	ASN
1	B	20	GLN
1	B	35	ASN
1	B	69	GLN
1	B	78	ASN
1	B	83	ASN

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Mol	Chain	Res	Type
1	B	125	ASN
1	B	142	ASN
1	B	228	GLN
1	B	236	ASN
1	B	282	ASN
1	B	332	GLN

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 ⓘ

4 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	AMP	A	339	-	25,25,25	1.32	5 (20%)	37,38,38	1.83	8 (21%)
2	F6P	B	338	-	15,16,16	1.06	0	16,25,25	1.22	2 (12%)
2	F6P	A	338	-	15,16,16	0.98	1 (6%)	16,25,25	1.30	2 (12%)
3	AMP	B	339	-	25,25,25	1.70	5 (20%)	37,38,38	1.96	14 (37%)

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	AMP	A	339	-	-	4/10/26/26	0/3/3/3
2	F6P	B	338	-	-	8/9/28/28	0/1/1/1
2	F6P	A	338	-	-	3/9/28/28	0/1/1/1
3	AMP	B	339	-	-	4/10/26/26	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	339	AMP	C8-N7	4.92	1.41	1.31
3	A	339	AMP	C8-N7	3.70	1.38	1.31
3	B	339	AMP	C2-N3	3.69	1.40	1.33
3	B	339	AMP	C5'-C4'	-2.76	1.43	1.51
3	B	339	AMP	C1'-N9	2.43	1.53	1.46
3	B	339	AMP	C4-N9	-2.37	1.32	1.37
3	A	339	AMP	C1'-N9	2.26	1.52	1.46
3	A	339	AMP	C5'-C4'	-2.25	1.44	1.51
3	A	339	AMP	C2-N1	2.18	1.37	1.33
2	A	338	F6P	O5-C2	-2.02	1.40	1.43
3	A	339	AMP	C4-N9	-2.02	1.33	1.37

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	339	AMP	O4'-C1'-N9	5.12	117.92	108.09
3	A	339	AMP	N3-C2-N1	-4.83	121.28	128.58
3	A	339	AMP	O4'-C1'-N9	3.68	115.16	108.09
3	B	339	AMP	O5'-P-O1P	3.34	115.47	106.44
3	A	339	AMP	C2-N1-C6	3.16	123.93	118.73
3	A	339	AMP	O3P-P-O5'	3.14	114.84	106.67
3	A	339	AMP	O5'-P-O1P	3.10	114.83	106.44
3	B	339	AMP	O2P-P-O5'	3.08	114.71	106.67
2	B	338	F6P	O1-C1-C2	2.94	118.18	111.67
3	B	339	AMP	O4'-C4'-C3'	-2.93	99.33	105.15
3	B	339	AMP	N3-C2-N1	-2.91	124.17	128.58
3	B	339	AMP	C6-C5-C4	2.90	121.14	117.18
3	B	339	AMP	C6-C5-N7	-2.71	126.87	132.09
3	B	339	AMP	C2'-C1'-N9	-2.62	106.80	113.30
2	B	338	F6P	O4-C4-C5	2.58	118.50	111.08
3	B	339	AMP	O5'-C5'-C4'	2.58	117.79	108.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	339	AMP	O4'-C4'-C3'	-2.44	100.31	105.15
2	A	338	F6P	O6-C6-C5	-2.42	100.74	108.99
3	A	339	AMP	O2P-P-O5'	2.40	112.93	106.67
3	B	339	AMP	N6-C6-N1	2.38	123.68	118.38
3	B	339	AMP	C2'-C3'-C4'	-2.35	98.07	102.61
3	B	339	AMP	C5-C4-N3	-2.27	123.59	126.72
3	B	339	AMP	O3P-P-O1P	-2.05	102.86	110.83
3	B	339	AMP	C2-N1-C6	2.04	122.08	118.73
3	A	339	AMP	C5-C4-N3	-2.00	123.96	126.72
2	A	338	F6P	O6-P-O1P	2.00	111.85	106.44

There are no chirality outliers.

All (19) torsion outliers are listed below:

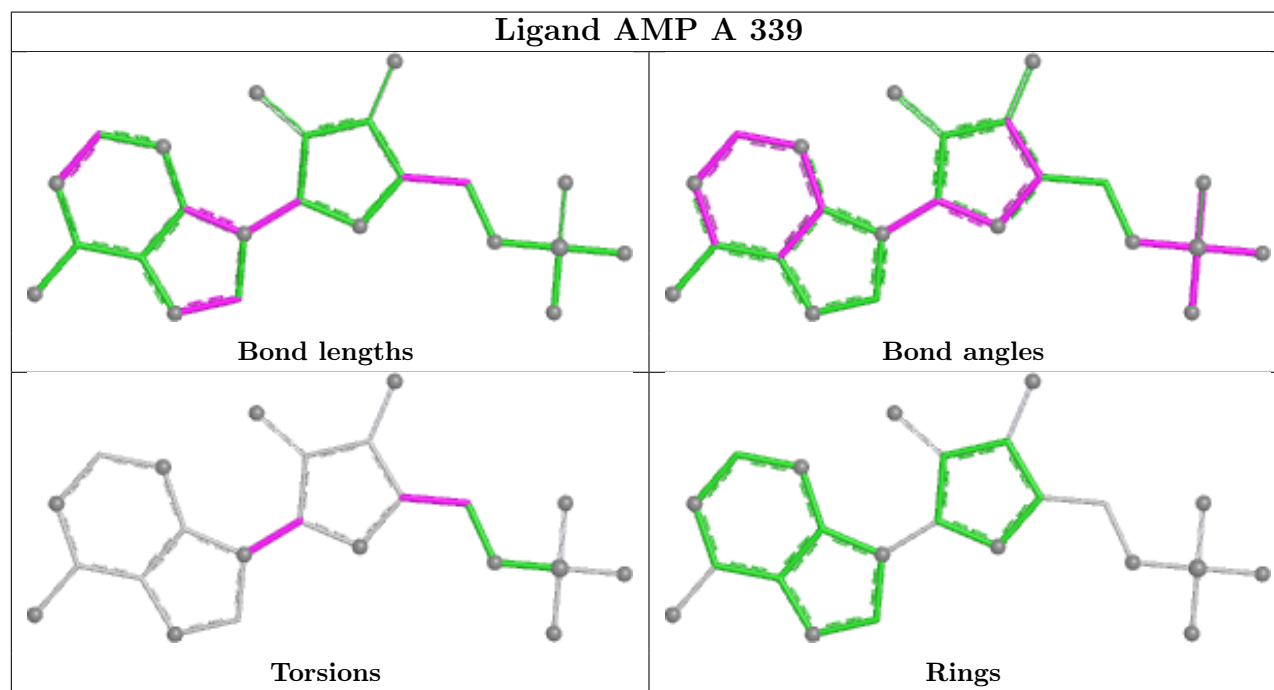
Mol	Chain	Res	Type	Atoms
2	B	338	F6P	O1-C1-C2-O2
2	B	338	F6P	O1-C1-C2-C3
2	B	338	F6P	O1-C1-C2-O5
2	B	338	F6P	C6-O6-P-O1P
2	B	338	F6P	C6-O6-P-O2P
2	B	338	F6P	C6-O6-P-O3P
3	B	339	AMP	O4'-C1'-N9-C8
2	B	338	F6P	C4-C5-C6-O6
3	B	339	AMP	C3'-C4'-C5'-O5'
3	A	339	AMP	O4'-C4'-C5'-O5'
3	B	339	AMP	O4'-C4'-C5'-O5'
3	A	339	AMP	O4'-C1'-N9-C4
2	A	338	F6P	O1-C1-C2-C3
2	B	338	F6P	O5-C5-C6-O6
2	A	338	F6P	O1-C1-C2-O5
3	A	339	AMP	C3'-C4'-C5'-O5'
3	B	339	AMP	O4'-C1'-N9-C4
2	A	338	F6P	O1-C1-C2-O2
3	A	339	AMP	O4'-C1'-N9-C8

There are no ring outliers.

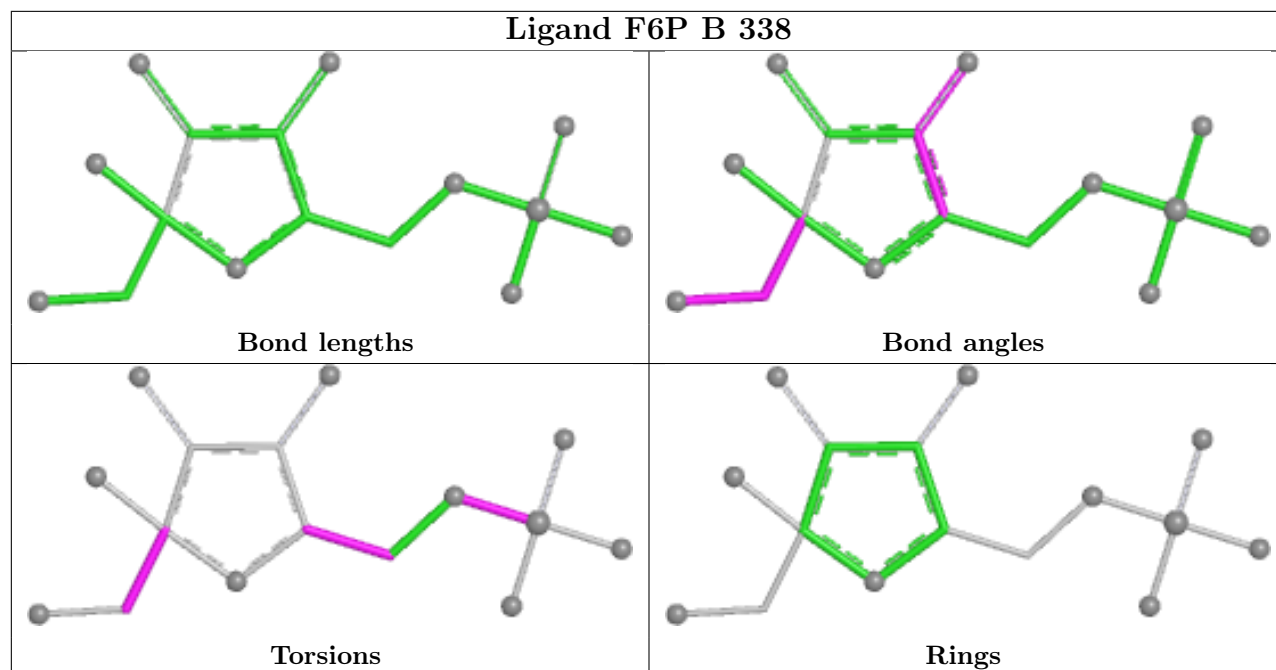
2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	339	AMP	1	0
3	B	339	AMP	5	0

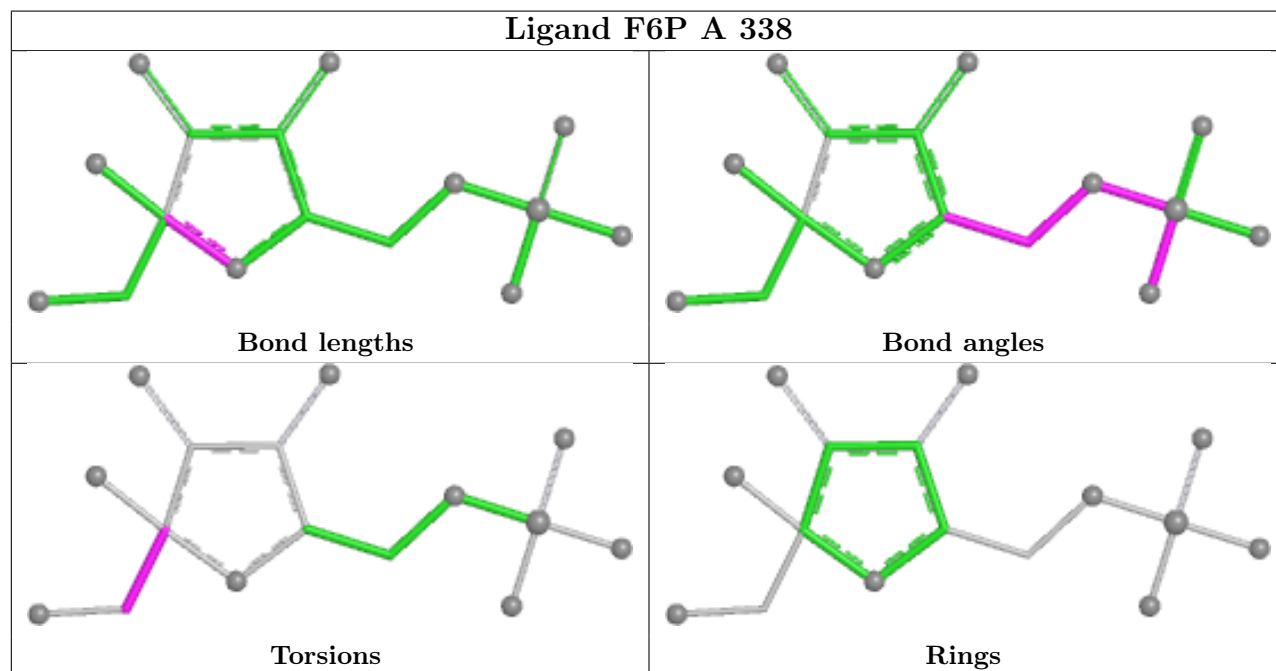
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

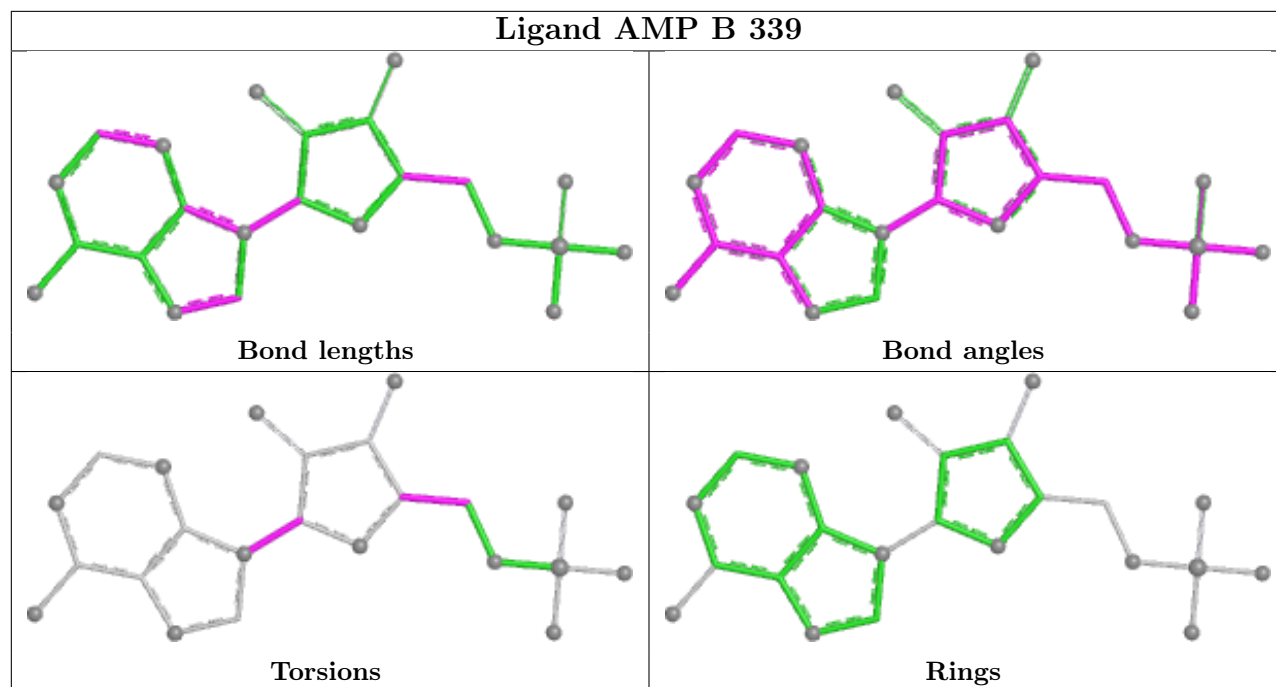


Ligand F6P B 338



Ligand F6P A 338





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.