



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 2, 2023 – 04:46 AM EDT

PDB ID : 1B99
Title : 3'-FLUORO-URIDINE DIPHOSPHATE BINDING TO NUCLEOSIDE
DIPHOSPHATE KINASE
Authors : Janin, J.; Xu, Y.
Deposited on : 1999-02-22
Resolution : 2.70 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.34

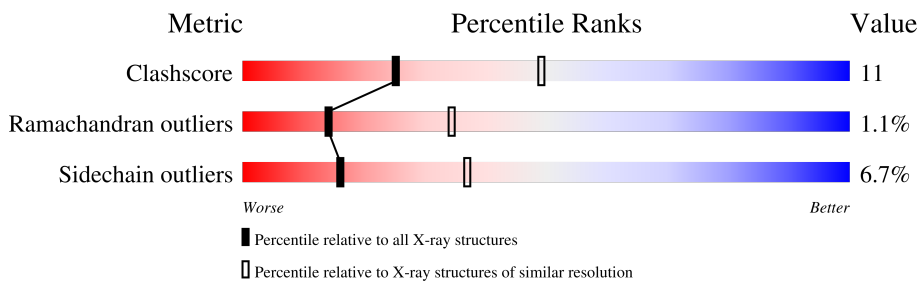
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)

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	155	69% 23% . . .
1	B	155	66% 28% . .
1	C	155	76% 18% . . .
1	D	155	74% 21% . .
1	E	155	64% 30% . .
1	F	155	75% 19% . .

2 Entry composition [i](#)

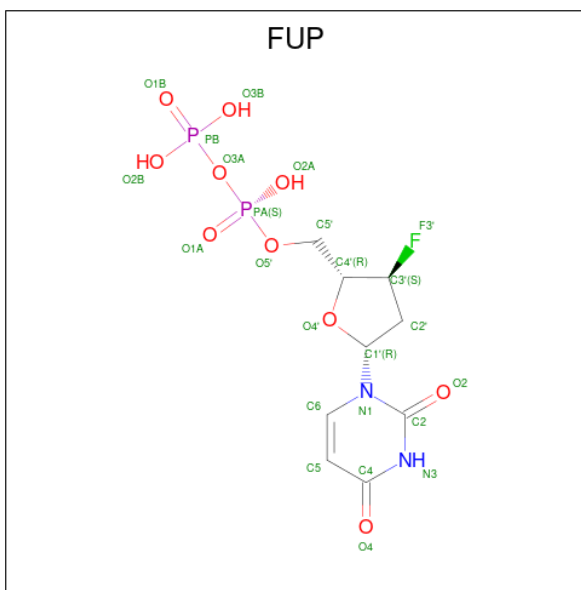
There are 4 unique types of molecules in this entry. The entry contains 6970 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 (NUCLEOSIDE DIPHOSPHATE KINASE).

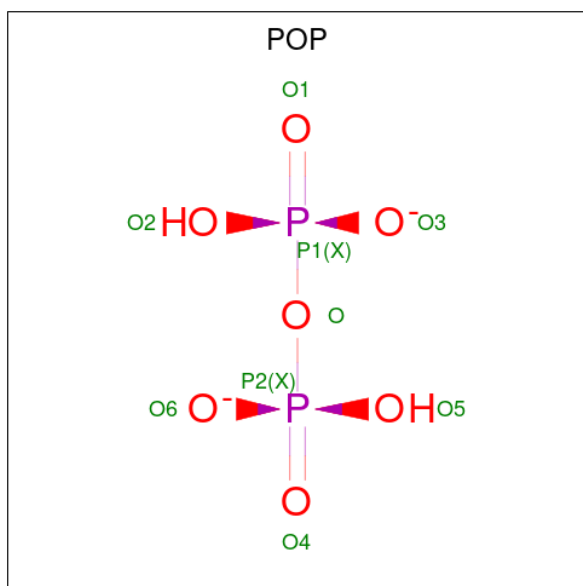
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	150	Total 1147	C 736	N 198	O 211	S 2	0	0	0
1	B	150	Total 1147	C 736	N 198	O 211	S 2	0	0	0
1	C	150	Total 1147	C 736	N 198	O 211	S 2	0	0	0
1	D	150	Total 1147	C 736	N 198	O 211	S 2	0	0	0
1	E	150	Total 1147	C 736	N 198	O 211	S 2	0	0	0
1	F	150	Total 1147	C 736	N 198	O 211	S 2	0	0	0

- Molecule 2 is 2',3'-DIDEOXY-3'-FLUORO-URIDIDINE-5'-DIPHOSPHATE (three-letter code: FUP) (formula: C₉H₁₃FN₂O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	F	N	O			P
2	A	1	24	9	1	2	10	2	0	0

- Molecule 3 is PYROPHOSPHATE 2- (three-letter code: POP) (formula: $H_2O_7P_2$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O P		
3	B	1	9	7 2	0	0
3	C	1	9	7 2	0	0
3	D	1	9	7 2	0	0
3	E	1	9	7 2	0	0
3	F	1	9	7 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	4	4	4	0	0
4	B	2	2	2	0	0
4	C	3	3	3	0	0
4	D	4	4	4	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	3	Total O 3 3	0	0
4	F	3	Total O 3 3	0	0

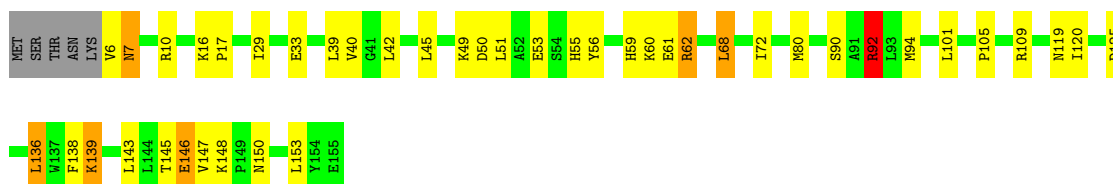
3 Residue-property plots [i](#)

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. 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.

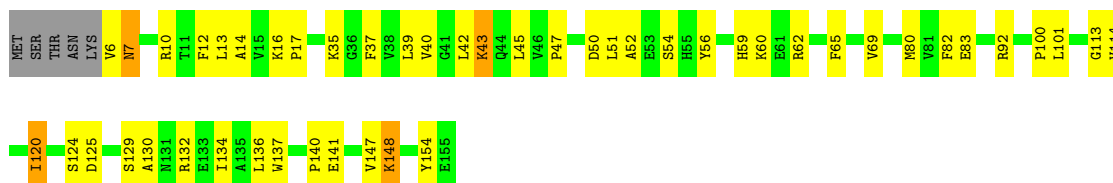
- Molecule 1: PROTEIN (NUCLEOSIDE DIPHOSPHATE KINASE)

Chain A: 




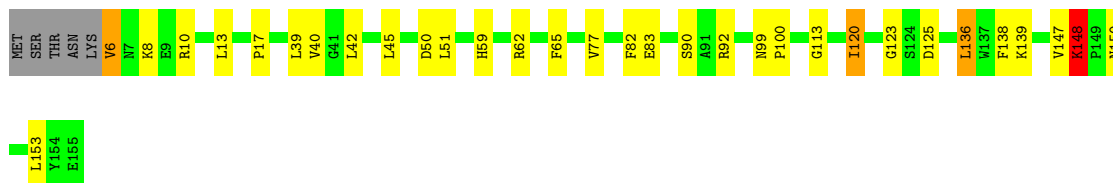
- Molecule 1: PROTEIN (NUCLEOSIDE DIPHOSPHATE KINASE)

Chain B: 



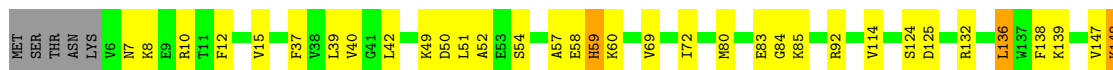
- Molecule 1: PROTEIN (NUCLEOSIDE DIPHOSPHATE KINASE)

Chain C: 



- Molecule 1: PROTEIN (NUCLEOSIDE DIPHOSPHATE KINASE)

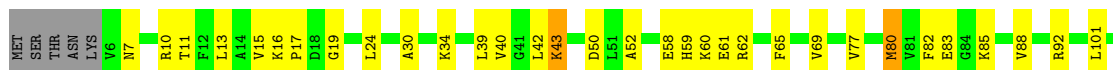
Chain D: 





- Molecule 1: PROTEIN (NUCLEOSIDE DIPHOSPHATE KINASE)

Chain E: 64% 30%



- Molecule 1: PROTEIN (NUCLEOSIDE DIPHOSPHATE KINASE)

Chain F: 75% 19%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.02Å 105.31Å 70.35Å 90.00° 117.43° 90.00°	Depositor
Resolution (Å)	6.00 – 2.70	Depositor
% Data completeness (in resolution range)	81.0 (6.00-2.70)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.207 , 0.304	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	6970	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

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: POP, FUP

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	0/1171	0.75	1/1583 (0.1%)
1	B	0.51	0/1171	0.73	1/1583 (0.1%)
1	C	0.50	0/1171	0.71	1/1583 (0.1%)
1	D	0.51	0/1171	0.71	0/1583
1	E	0.53	0/1171	0.72	1/1583 (0.1%)
1	F	0.50	0/1171	0.72	1/1583 (0.1%)
All	All	0.51	0/7026	0.72	5/9498 (0.1%)

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
1	F	0	1
All	All	0	2

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	132	ARG	NE-CZ-NH2	-5.61	117.49	120.30
1	C	120	ILE	N-CA-C	5.57	126.03	111.00
1	F	120	ILE	N-CA-C	5.50	125.86	111.00
1	A	39	LEU	CA-CB-CG	5.41	127.74	115.30
1	B	120	ILE	N-CA-C	5.37	125.50	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	92	ARG	Sidechain
1	F	32	TYR	Sidechain

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	1147	0	1167	33	0
1	B	1147	0	1167	29	0
1	C	1147	0	1167	21	0
1	D	1147	0	1167	23	0
1	E	1147	0	1167	30	0
1	F	1147	0	1167	20	0
2	A	24	0	10	0	0
3	B	9	0	0	0	0
3	C	9	0	0	0	0
3	D	9	0	0	0	0
3	E	9	0	0	0	0
3	F	9	0	0	0	0
4	A	4	0	0	0	0
4	B	2	0	0	0	0
4	C	3	0	0	0	0
4	D	4	0	0	0	0
4	E	3	0	0	0	0
4	F	3	0	0	0	0
All	All	6970	0	7012	149	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 11.

All (149) 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:92:ARG:HH21	1:A:125:ASP:HB3	1.14	1.06
1:C:92:ARG:HD2	1:C:125:ASP:HA	1.54	0.89
1:B:92:ARG:HD2	1:B:125:ASP:HA	1.56	0.88
1:A:92:ARG:NH2	1:A:125:ASP:HB3	1.95	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:92:ARG:HD2	1:E:124:SER:O	1.83	0.79
1:A:150:ASN:HB3	1:A:153:LEU:HG	1.68	0.75
1:A:62:ARG:HG2	1:A:62:ARG:HH11	1.51	0.74
1:E:150:ASN:HB3	1:E:153:LEU:HG	1.72	0.71
1:D:51:LEU:HD12	1:D:136:LEU:HD22	1.72	0.71
1:F:92:ARG:HH11	1:F:92:ARG:HG2	1.54	0.71
1:E:147:VAL:HG13	1:E:148:LYS:H	1.57	0.70
1:B:101:LEU:HD21	1:B:114:VAL:HG23	1.74	0.68
1:E:7:ASN:HD22	1:E:85:LYS:HB2	1.58	0.68
1:E:30:ALA:O	1:E:34:LYS:HG3	1.94	0.68
1:A:62:ARG:HH11	1:A:62:ARG:CG	2.05	0.68
1:F:92:ARG:HD2	1:F:125:ASP:HA	1.76	0.68
1:C:92:ARG:HG2	1:C:92:ARG:HH11	1.60	0.66
1:A:90:SER:O	1:A:94:MET:HG3	1.96	0.65
1:F:17:PRO:HD3	1:F:77:VAL:HG22	1.77	0.64
1:E:139:LYS:N	1:E:139:LYS:HE3	2.13	0.62
1:B:147:VAL:HG13	1:B:148:LYS:H	1.65	0.61
1:F:38:VAL:HG11	1:F:147:VAL:HG13	1.81	0.61
1:D:10:ARG:HG2	1:D:83:GLU:HG3	1.83	0.60
1:A:56:TYR:CE2	1:A:68:LEU:HD22	2.36	0.60
1:A:139:LYS:HA	1:A:139:LYS:HE3	1.83	0.60
1:F:59:HIS:HA	1:F:62:ARG:HD2	1.84	0.60
1:F:92:ARG:HG2	1:F:92:ARG:NH1	2.17	0.60
1:C:83:GLU:OE2	1:C:147:VAL:HG11	2.01	0.59
1:A:109:ARG:NH2	1:A:119:ASN:ND2	2.51	0.59
1:F:49:LYS:HG3	1:F:50:ASP:N	2.17	0.59
1:C:138:PHE:C	1:C:139:LYS:HE2	2.23	0.58
1:A:62:ARG:HG2	1:A:62:ARG:NH1	2.19	0.58
1:E:62:ARG:HH11	1:E:62:ARG:HG2	1.69	0.58
1:D:8:LYS:O	1:D:8:LYS:HG3	2.02	0.58
1:C:147:VAL:HG13	1:C:148:LYS:H	1.68	0.57
1:C:139:LYS:HE2	1:C:139:LYS:N	2.19	0.57
1:D:92:ARG:HD2	1:D:124:SER:O	2.05	0.57
1:C:39:LEU:HD13	1:C:82:PHE:CZ	2.41	0.56
1:D:52:ALA:HB3	1:D:69:VAL:HG13	1.87	0.56
1:E:43:LYS:HG2	1:E:138:PHE:CE1	2.40	0.56
1:B:92:ARG:HD2	1:B:124:SER:O	2.04	0.56
1:E:60:LYS:HA	1:E:65:PHE:CD1	2.41	0.56
1:E:92:ARG:HD2	1:E:125:ASP:HA	1.87	0.56
1:E:92:ARG:CD	1:E:125:ASP:HA	2.36	0.55
1:D:92:ARG:HD2	1:D:125:ASP:HA	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:10:ARG:CZ	1:F:40:VAL:HG11	2.37	0.55
1:E:147:VAL:HG13	1:E:148:LYS:N	2.22	0.54
1:A:59:HIS:HA	1:A:62:ARG:CD	2.38	0.54
1:D:150:ASN:OD1	1:D:152:ASN:ND2	2.41	0.54
1:D:147:VAL:HG13	1:D:148:LYS:H	1.72	0.53
1:C:150:ASN:HB3	1:C:153:LEU:HG	1.89	0.53
1:C:42:LEU:HD11	1:F:42:LEU:HD21	1.90	0.53
1:B:52:ALA:HB3	1:B:69:VAL:HG13	1.91	0.53
1:B:92:ARG:CD	1:B:125:ASP:HA	2.34	0.52
1:C:10:ARG:CZ	1:C:40:VAL:HG11	2.38	0.52
1:E:52:ALA:HB3	1:E:69:VAL:HG13	1.91	0.52
1:E:39:LEU:HD21	1:E:42:LEU:HD22	1.92	0.52
1:D:7:ASN:HD22	1:D:85:LYS:HB2	1.75	0.52
1:B:83:GLU:OE2	1:B:147:VAL:HG11	2.10	0.51
1:B:54:SER:O	1:B:132:ARG:NH2	2.44	0.51
1:B:130:ALA:O	1:B:134:ILE:HG13	2.11	0.51
1:B:51:LEU:HD12	1:B:136:LEU:HD23	1.94	0.50
1:A:51:LEU:CD1	1:A:136:LEU:HD22	2.41	0.50
1:A:147:VAL:HG13	1:A:148:LYS:H	1.76	0.50
1:B:35:LYS:HG2	1:B:37:PHE:CE2	2.46	0.50
1:E:83:GLU:OE2	1:E:147:VAL:HG11	2.12	0.50
1:E:92:ARG:HH11	1:E:92:ARG:HG2	1.76	0.50
1:D:39:LEU:HD21	1:D:42:LEU:HD22	1.94	0.50
1:A:147:VAL:HG13	1:A:148:LYS:N	2.27	0.49
1:D:51:LEU:HD12	1:D:136:LEU:CD2	2.41	0.49
1:B:10:ARG:CZ	1:B:40:VAL:HG11	2.43	0.49
1:B:54:SER:OG	1:B:136:LEU:HD21	2.12	0.49
1:D:12:PHE:HE1	1:D:138:PHE:HE1	1.61	0.49
1:E:13:LEU:HD12	1:E:82:PHE:HD2	1.78	0.49
1:C:51:LEU:CD1	1:C:136:LEU:HD22	2.43	0.48
1:E:17:PRO:HD3	1:E:77:VAL:HG22	1.96	0.48
1:B:56:TYR:O	1:B:65:PHE:HE1	1.96	0.48
1:C:45:LEU:HD23	1:C:45:LEU:N	2.28	0.48
1:D:40:VAL:HG21	1:D:83:GLU:HB2	1.95	0.48
1:A:59:HIS:HA	1:A:62:ARG:HD3	1.97	0.47
1:E:19:GLY:HA2	1:E:24:LEU:HD12	1.96	0.47
1:E:10:ARG:CZ	1:E:40:VAL:HG11	2.44	0.47
1:D:139:LYS:HA	1:D:139:LYS:HE2	1.97	0.47
1:A:6:VAL:HG22	1:A:7:ASN:N	2.30	0.46
1:D:51:LEU:CD1	1:D:136:LEU:HD22	2.44	0.46
1:E:92:ARG:HG2	1:E:92:ARG:NH1	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:LEU:HD23	1:C:123:GLY:HA2	1.98	0.46
1:A:62:ARG:HH11	1:A:62:ARG:CB	2.28	0.46
1:A:101:LEU:HD11	1:F:85:LYS:HG2	1.98	0.46
1:B:45:LEU:HD23	1:B:45:LEU:N	2.31	0.46
1:E:101:LEU:HD21	1:E:114:VAL:HG23	1.97	0.46
1:B:39:LEU:HD21	1:B:42:LEU:HD22	1.97	0.45
1:F:91:ALA:O	1:F:95:ILE:HG13	2.16	0.45
1:A:105:PRO:HG2	1:F:94:MET:HA	1.99	0.45
1:D:7:ASN:ND2	1:D:85:LYS:HD3	2.32	0.45
1:D:15:VAL:HG23	1:D:80:MET:CE	2.46	0.45
1:C:62:ARG:N	1:C:62:ARG:CD	2.80	0.45
1:A:16:LYS:HB3	1:A:17:PRO:HD2	1.98	0.44
1:B:7:ASN:OD1	1:B:7:ASN:N	2.49	0.44
1:E:139:LYS:HE3	1:E:139:LYS:H	1.82	0.44
1:D:59:HIS:N	1:D:59:HIS:ND1	2.65	0.44
1:F:7:ASN:HD22	1:F:85:LYS:HB2	1.82	0.44
1:A:10:ARG:CZ	1:A:40:VAL:HG11	2.47	0.44
1:A:109:ARG:NE	1:A:119:ASN:HB2	2.33	0.44
1:B:140:PRO:HD2	1:B:141:GLU:OE2	2.18	0.44
1:F:52:ALA:HB3	1:F:69:VAL:HG13	1.99	0.44
1:A:62:ARG:HH11	1:A:62:ARG:HB3	1.83	0.43
1:D:54:SER:O	1:D:132:ARG:NH2	2.43	0.43
1:E:59:HIS:O	1:E:65:PHE:HB2	2.18	0.43
1:B:100:PRO:HB2	1:B:113:GLY:C	2.39	0.43
1:C:42:LEU:O	1:F:43:LYS:HA	2.19	0.43
1:C:6:VAL:HG13	1:C:8:LYS:H	1.82	0.43
1:A:146:GLU:CD	1:A:146:GLU:H	2.21	0.43
1:A:29:ILE:O	1:A:33:GLU:HG3	2.19	0.43
1:A:55:HIS:CD2	1:A:56:TYR:CE1	3.07	0.43
1:B:47:PRO:HB3	1:B:51:LEU:HD23	2.00	0.43
1:E:62:ARG:HG2	1:E:62:ARG:NH1	2.34	0.43
1:A:49:LYS:O	1:A:53:GLU:HG3	2.19	0.42
1:B:51:LEU:HD12	1:B:136:LEU:CD2	2.49	0.42
1:D:57:ALA:O	1:D:60:LYS:HG2	2.19	0.42
1:F:39:LEU:HD21	1:F:42:LEU:HD22	2.00	0.42
1:B:59:HIS:O	1:B:65:PHE:HB2	2.18	0.42
1:B:154:TYR:CE1	1:D:114:VAL:CG2	3.03	0.42
1:F:16:LYS:HB3	1:F:17:PRO:HD2	2.01	0.42
1:E:15:VAL:HG23	1:E:80:MET:CE	2.49	0.42
1:E:16:LYS:NZ	1:E:119:ASN:OD1	2.52	0.42
1:C:59:HIS:O	1:C:65:PHE:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12:PHE:CZ	1:B:14:ALA:HB2	2.55	0.41
1:B:60:LYS:HB3	1:B:60:LYS:HE3	1.66	0.41
1:B:43:LYS:NZ	1:B:137:TRP:O	2.53	0.41
1:F:60:LYS:HA	1:F:65:PHE:CD2	2.55	0.41
1:B:16:LYS:HB3	1:B:17:PRO:HD2	2.03	0.41
1:A:59:HIS:HA	1:A:62:ARG:HD2	2.01	0.41
1:C:147:VAL:HG13	1:C:148:LYS:N	2.34	0.41
1:A:42:LEU:HD11	1:D:42:LEU:HD21	2.02	0.41
1:A:51:LEU:HD12	1:A:136:LEU:HD22	2.03	0.41
1:C:99:ASN:HA	1:C:100:PRO:HD2	1.95	0.41
1:A:138:PHE:CB	1:A:143:LEU:HD21	2.51	0.41
1:C:100:PRO:HB2	1:C:113:GLY:C	2.41	0.41
1:D:37:PHE:CD1	1:D:84:GLY:HA3	2.56	0.41
1:E:58:GLU:HG3	1:E:59:HIS:CD2	2.55	0.41
1:F:57:ALA:O	1:F:60:LYS:HG2	2.21	0.41
1:E:11:THR:HB	1:E:88:VAL:HG22	2.03	0.40
1:A:94:MET:HA	1:E:105:PRO:HG2	2.02	0.40
1:F:100:PRO:HB2	1:F:113:GLY:C	2.41	0.40
1:B:13:LEU:HD12	1:B:82:PHE:CD2	2.57	0.40
1:B:60:LYS:HA	1:B:65:PHE:CG	2.57	0.40
1:C:17:PRO:HD3	1:C:77:VAL:HG22	2.03	0.40
1:A:56:TYR:CD2	1:A:68:LEU:HD13	2.57	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	148/155 (96%)	138 (93%)	9 (6%)	1 (1%)	22	46
1	B	148/155 (96%)	138 (93%)	8 (5%)	2 (1%)	11	28
1	C	148/155 (96%)	141 (95%)	5 (3%)	2 (1%)	11	28

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	148/155 (96%)	141 (95%)	6 (4%)	1 (1%)	22	46
1	E	148/155 (96%)	141 (95%)	5 (3%)	2 (1%)	11	28
1	F	148/155 (96%)	139 (94%)	7 (5%)	2 (1%)	11	28
All	All	888/930 (96%)	838 (94%)	40 (4%)	10 (1%)	14	34

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	120	ILE
1	A	120	ILE
1	E	148	LYS
1	D	148	LYS
1	F	148	LYS
1	B	148	LYS
1	C	120	ILE
1	C	148	LYS
1	E	120	ILE
1	F	120	ILE

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	122/127 (96%)	108 (88%)	14 (12%)	5	13
1	B	122/127 (96%)	115 (94%)	7 (6%)	20	44
1	C	122/127 (96%)	117 (96%)	5 (4%)	30	59
1	D	122/127 (96%)	116 (95%)	6 (5%)	25	52
1	E	122/127 (96%)	111 (91%)	11 (9%)	9	22
1	F	122/127 (96%)	116 (95%)	6 (5%)	25	52
All	All	732/762 (96%)	683 (93%)	49 (7%)	16	37

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	45	LEU
1	A	50	ASP
1	A	60	LYS
1	A	61	GLU
1	A	62	ARG
1	A	68	LEU
1	A	72	ILE
1	A	80	MET
1	A	92	ARG
1	A	136	LEU
1	A	139	LYS
1	A	145	THR
1	A	146	GLU
1	B	6	VAL
1	B	7	ASN
1	B	43	LYS
1	B	50	ASP
1	B	62	ARG
1	B	80	MET
1	B	129	SER
1	C	6	VAL
1	C	50	ASP
1	C	90	SER
1	C	136	LEU
1	C	148	LYS
1	D	49	LYS
1	D	50	ASP
1	D	58	GLU
1	D	59	HIS
1	D	72	ILE
1	D	136	LEU
1	E	43	LYS
1	E	50	ASP
1	E	61	GLU
1	E	80	MET
1	E	115	ASP
1	E	116	VAL
1	E	136	LEU
1	E	139	LYS
1	E	141	GLU
1	E	145	THR
1	E	146	GLU

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Mol	Chain	Res	Type
1	F	7	ASN
1	F	17	PRO
1	F	61	GLU
1	F	116	VAL
1	F	136	LEU
1	F	148	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	44	GLN
1	A	99	ASN
1	B	59	HIS
1	B	99	ASN
1	B	122	HIS
1	C	7	ASN
1	D	7	ASN
1	D	99	ASN
1	E	7	ASN
1	F	7	ASN
1	F	99	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
3	POP	F	660	-	6,8,8	1.40	0	13,13,13	0.89	1 (7%)
2	FUP	A	160	-	23,25,25	1.08	1 (4%)	35,38,38	1.93	8 (22%)
3	POP	D	460	-	6,8,8	1.38	0	13,13,13	0.90	1 (7%)
3	POP	E	560	-	6,8,8	1.43	0	13,13,13	0.89	0
3	POP	C	360	-	6,8,8	1.70	2 (33%)	13,13,13	0.84	0
3	POP	B	260	-	6,8,8	1.31	0	13,13,13	1.08	1 (7%)

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	POP	F	660	-	-	3/6/6/6	-
2	FUP	A	160	-	-	4/16/28/28	0/2/2/2
3	POP	D	460	-	-	0/6/6/6	-
3	POP	E	560	-	-	2/6/6/6	-
3	POP	C	360	-	-	0/6/6/6	-
3	POP	B	260	-	-	0/6/6/6	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	360	POP	P2-O5	-2.06	1.46	1.54
3	C	360	POP	P1-O3	2.04	1.62	1.54
2	A	160	FUP	C2-N1	2.04	1.41	1.38

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	160	FUP	C5'-C4'-C3'	-5.33	109.81	114.70
2	A	160	FUP	F3'-C3'-C2'	5.21	120.00	109.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	160	FUP	C3'-C2'-C1'	4.14	106.03	102.36
2	A	160	FUP	O4'-C1'-C2'	-3.45	99.73	106.25
2	A	160	FUP	C2'-C1'-N1	3.03	120.75	113.77
2	A	160	FUP	O2-C2-N3	-2.76	116.35	121.50
3	B	260	POP	O2-P1-O1	2.64	121.03	110.68
2	A	160	FUP	N3-C2-N1	2.42	118.10	114.89
3	D	460	POP	O2-P1-O1	2.17	119.17	110.68
3	F	660	POP	O2-P1-O1	2.12	118.97	110.68
2	A	160	FUP	O2-C2-N1	2.07	125.54	122.79

There are no chirality outliers.

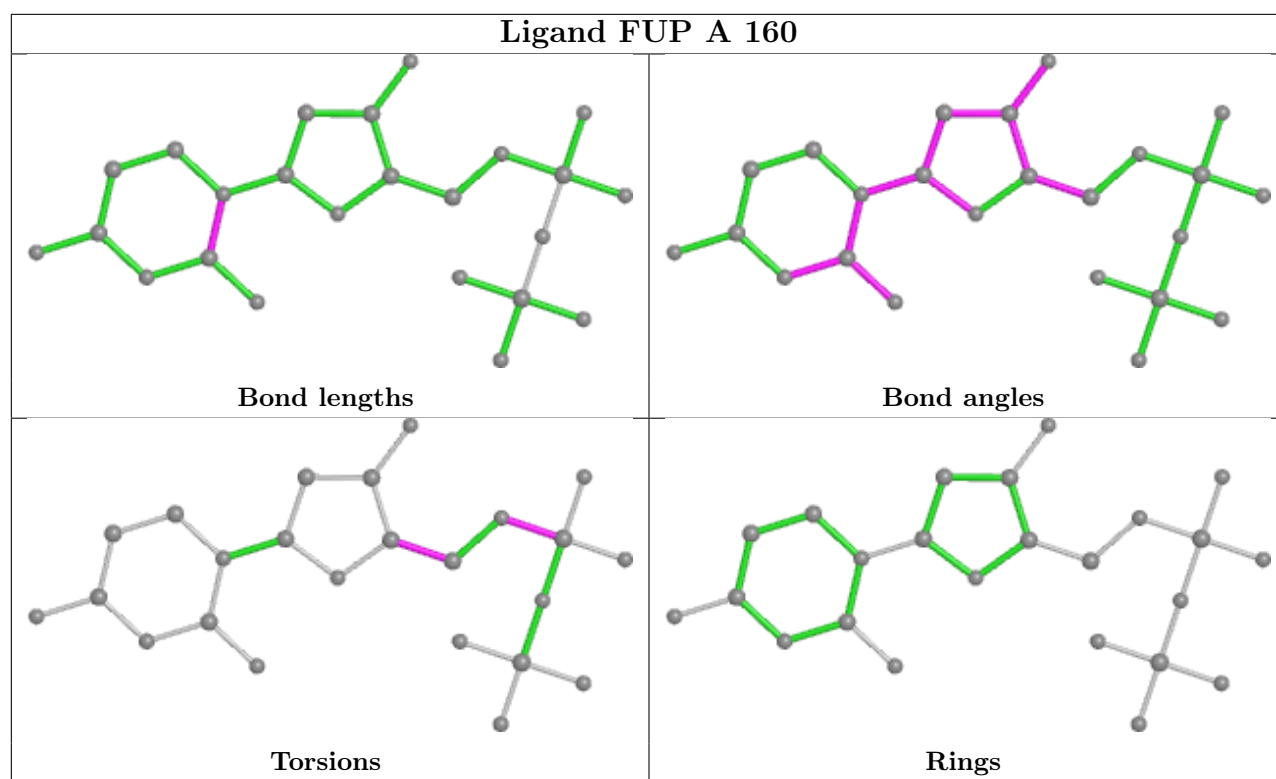
All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	160	FUP	C5'-O5'-PA-O1A
3	E	560	POP	P1-O-P2-O5
3	F	660	POP	P1-O-P2-O4
2	A	160	FUP	C5'-O5'-PA-O3A
2	A	160	FUP	O4'-C4'-C5'-O5'
2	A	160	FUP	C5'-O5'-PA-O2A
3	E	560	POP	P2-O-P1-O2
3	F	660	POP	P1-O-P2-O5
3	F	660	POP	P1-O-P2-O6

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.