



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 08:16 pm BST

PDB ID : 2BI1
Title : crystal structures of transition state analogue inhibitors of inosine monophosphate cyclohydrolase
Authors : Xu, L.; Chong, Y.; Hwang, I.; D'Onofrio, A.; Amore, K.; Beardsley, G.P.; Li, C.; Olson, A.J.; Boger, D.L.; Wilson, I.A.
Deposited on : 2005-09-15
Resolution : 2.02 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 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)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

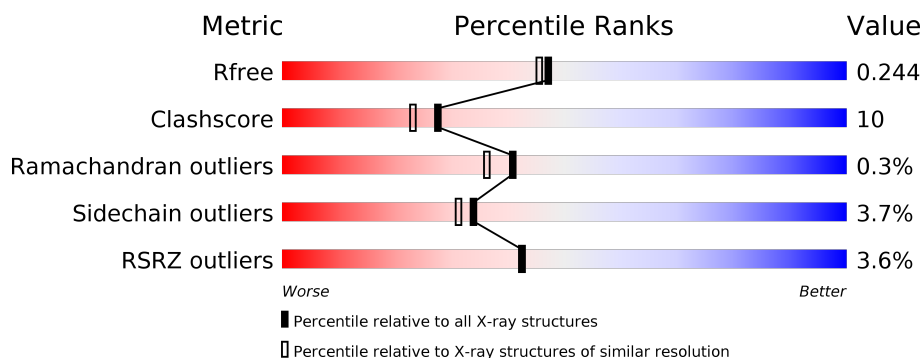
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.02 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	10434 (2.04-2.00)
Clashscore	141614	11643 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	593	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>23%</div> <div>..</div> </div> </div>
1	B	593	<div> <div>4%</div> <div> <div></div> <div>78%</div> <div>21%</div> <div>..</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9423 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.

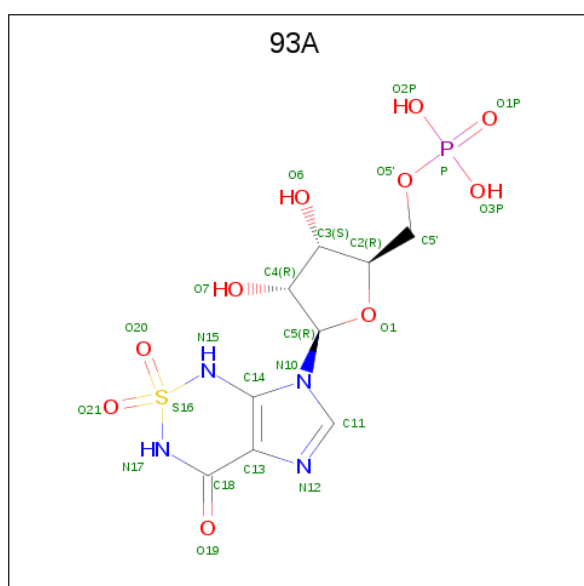
- Molecule 1 is a protein called Bifunctional purine biosynthesis protein PURH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	590	Total	C	N	O	S	0	0	0
			4511	2843	800	849	19			
1	B	590	Total	C	N	O	S	0	0	0
			4511	2843	800	849	19			

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	K	0	0
			1	1		
2	A	1	Total	K	0	0
			1	1		

- Molecule 3 is [3,4-DIHYDROXY-5R-(2,2,4-TRIOXO-1,2R,3S,4R-TETRAHYDRO-2L6-IMIDAZO[4,5-C][1,2,6]THIADIAZIN-7-YL)TETRAHYDROFURAN-2-YL]METHYL DIHYDROGEN PHOSPHATE (three-letter code: 93A) (formula: C₉H₁₃N₄O₁₀PS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			25	9	4	10	1	1		
3	A	1	Total	C	N	O	P	S	0	0
			25	9	4	10	1	1		
3	B	1	Total	C	N	O	P	S	0	0
			25	9	4	10	1	1		
3	B	1	Total	C	N	O	P	S	0	0
			25	9	4	10	1	1		

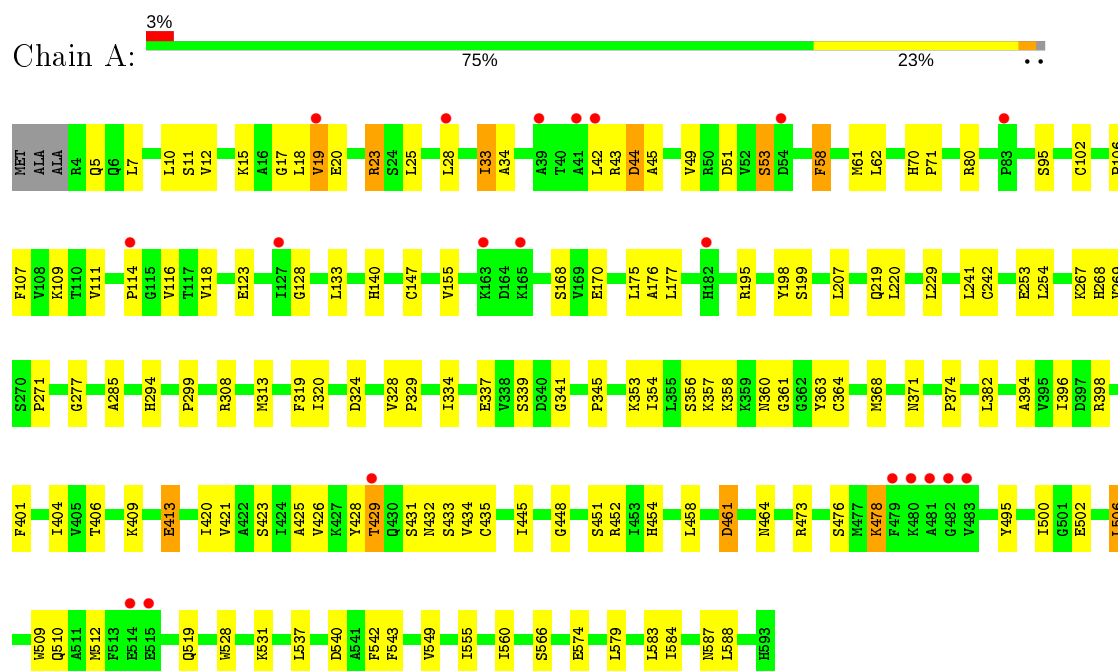
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	162	Total	O	0	0
			162	162		
4	B	137	Total	O	0	0
			137	137		

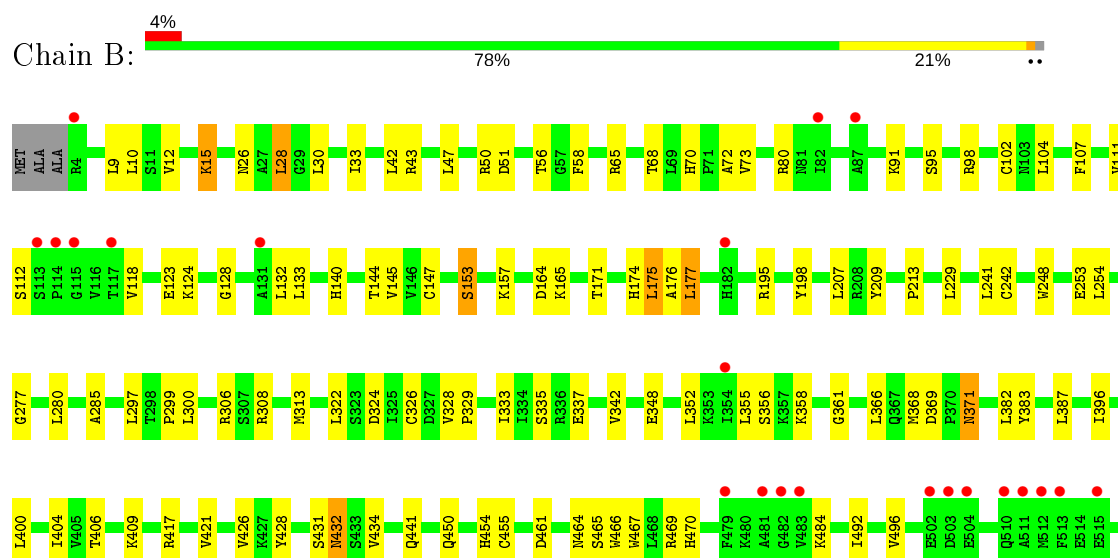
3 Residue-property plots [i](#)

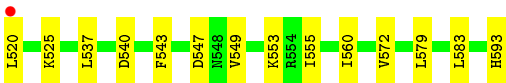
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Bifunctional purine biosynthesis protein PURH



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.00 Å 106.50 Å 101.00 Å 90.00° 91.50° 90.00°	Depositor
Resolution (Å)	38.07 – 2.02 37.99 – 2.02	Depositor EDS
% Data completeness (in resolution range)	92.3 (38.07-2.02) 89.5 (37.99-2.02)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.01 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.198 , 0.249 0.197 , 0.244	Depositor DCC
R_{free} test set	1889 reflections (2.51%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.827	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.032 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9423	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 93A, K

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	0.91	11/4595 (0.2%)	0.78	7/6230 (0.1%)
1	B	0.57	1/4595 (0.0%)	0.67	0/6230
All	All	0.76	12/9190 (0.1%)	0.73	7/12460 (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

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	23	ARG	CZ-NH1	30.59	1.72	1.33
1	A	44	ASP	CG-OD2	15.84	1.61	1.25
1	A	20	GLU	CD-OE2	15.58	1.42	1.25
1	A	20	GLU	CG-CD	14.05	1.73	1.51
1	A	20	GLU	CD-OE1	7.48	1.33	1.25
1	A	17	GLY	C-N	7.48	1.51	1.34
1	A	43	ARG	CZ-NH1	7.32	1.42	1.33
1	A	20	GLU	CA-CB	7.20	1.69	1.53
1	A	43	ARG	NE-CZ	7.04	1.42	1.33
1	B	91	LYS	CE-NZ	5.38	1.62	1.49
1	A	364	CYS	CB-SG	-5.09	1.73	1.81
1	A	20	GLU	C-N	-5.06	1.22	1.34

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	ASP	CB-CG-OD2	-11.86	107.63	118.30
1	A	44	ASP	CB-CG-OD1	-11.51	107.94	118.30
1	A	43	ARG	NE-CZ-NH2	-11.06	114.77	120.30
1	A	43	ARG	NE-CZ-NH1	10.20	125.40	120.30
1	A	44	ASP	OD1-CG-OD2	8.70	139.82	123.30
1	A	17	GLY	O-C-N	6.31	132.80	122.70
1	A	23	ARG	NE-CZ-NH2	6.07	123.33	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	44	ASP	Sidechain

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	4511	0	4560	111	1
1	B	4511	0	4559	91	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	50	0	22	3	0
3	B	50	0	22	1	0
4	A	162	0	0	5	0
4	B	137	0	0	0	0
All	All	9423	0	9163	183	1

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 10.

All (183) 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:23:ARG:CZ	1:A:23:ARG:NH1	1.72	1.48
1:B:434:VAL:CG1	1:B:537:LEU:HD11	1.92	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:MET:CE	1:B:454:HIS:HE1	1.79	0.94
1:A:429:THR:HG21	1:A:445:ILE:HD11	1.50	0.92
1:A:313:MET:HE2	1:B:454:HIS:HE1	1.44	0.82
1:B:434:VAL:HG12	1:B:537:LEU:HD11	1.63	0.80
1:A:170:GLU:HG3	4:A:1059:HOH:O	1.82	0.78
1:B:464:ASN:HD22	1:B:555:ILE:HD13	1.52	0.75
1:A:18:LEU:HG	1:A:42:LEU:HD21	1.69	0.74
1:A:313:MET:HE3	1:B:454:HIS:HE1	1.49	0.73
1:B:153:SER:O	1:B:157:LYS:HG2	1.89	0.71
1:B:356:SER:O	1:B:361:GLY:HA2	1.91	0.71
1:B:465:SER:O	1:B:469:ARG:HG3	1.91	0.71
1:B:277:GLY:HA3	1:B:299:PRO:O	1.90	0.71
1:A:268:HIS:HD2	1:B:593:HIS:NE2	1.89	0.70
1:B:30:LEU:HD21	1:B:98:ARG:HE	1.59	0.68
1:A:313:MET:HE3	1:B:454:HIS:CE1	2.28	0.66
1:A:308:ARG:CZ	1:A:337:GLU:HG3	2.25	0.66
1:A:313:MET:CE	1:B:454:HIS:CE1	2.71	0.66
1:B:33:ILE:HG22	1:B:50:ARG:HB3	1.78	0.65
1:B:102:CYS:O	1:B:147:CYS:HA	1.95	0.65
1:A:308:ARG:NE	1:A:337:GLU:HG3	2.11	0.65
1:A:195:ARG:HG3	1:B:177:LEU:HD21	1.79	0.65
1:A:10:LEU:HD12	1:A:42:LEU:HD11	1.79	0.64
1:B:432:ASN:C	1:B:432:ASN:HD22	2.02	0.63
1:A:425:ALA:O	1:A:429:THR:CG2	2.46	0.63
1:A:133:LEU:HD22	1:A:147:CYS:HB3	1.79	0.63
1:A:313:MET:HE2	1:B:454:HIS:CE1	2.30	0.62
1:B:280:LEU:HD11	1:B:306:ARG:HG3	1.80	0.62
1:A:7:LEU:HD11	1:A:95:SER:HB2	1.81	0.61
1:A:308:ARG:CD	1:A:337:GLU:HG2	2.31	0.61
1:A:354:ILE:O	1:A:357:LYS:HG2	1.99	0.61
1:A:7:LEU:HB3	1:A:33:ILE:HD11	1.83	0.60
1:B:431:SER:HB3	1:B:432:ASN:HA	1.84	0.60
1:A:253:GLU:OE2	1:A:428:TYR:OH	2.13	0.60
1:B:464:ASN:ND2	1:B:555:ILE:HD13	2.15	0.60
1:A:7:LEU:HD13	1:A:33:ILE:HD11	1.82	0.59
1:A:242:CYS:HA	1:B:382:LEU:CD2	2.33	0.59
1:B:306:ARG:HB3	1:B:441:GLN:HG3	1.86	0.58
1:B:171:THR:HG22	1:B:175:LEU:HD22	1.86	0.57
1:A:80:ARG:HD2	1:B:65:ARG:CZ	2.35	0.57
1:B:396:ILE:HG23	1:B:400:LEU:HD23	1.87	0.57
1:B:426:VAL:HG13	1:B:540:ASP:HB3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:VAL:HG11	1:A:198:TYR:OH	2.04	0.56
1:A:426:VAL:HG13	1:A:540:ASP:HB3	1.86	0.56
1:B:434:VAL:HG12	1:B:537:LEU:CD1	2.33	0.56
1:B:229:LEU:HD11	1:B:366:LEU:HB3	1.87	0.56
1:A:458:LEU:HD12	4:A:983:HOH:O	2.04	0.56
1:A:308:ARG:NE	1:A:337:GLU:CG	2.68	0.56
1:A:241:LEU:HD12	1:B:387:LEU:HD21	1.88	0.56
1:A:229:LEU:HD13	1:A:368:MET:CE	2.36	0.56
1:B:396:ILE:HG23	1:B:400:LEU:CD2	2.36	0.55
1:A:396:ILE:HD13	1:A:423:SER:HB3	1.88	0.55
1:A:102:CYS:O	1:A:147:CYS:HA	2.07	0.54
1:B:297:LEU:HD22	1:B:333:ILE:HD11	1.89	0.54
1:A:242:CYS:HA	1:B:382:LEU:HD21	1.88	0.54
1:A:543:PHE:HE2	1:A:560:ILE:HG21	1.72	0.54
1:B:254:LEU:HD21	1:B:421:VAL:HA	1.89	0.54
1:B:431:SER:CB	1:B:432:ASN:HA	2.38	0.54
1:A:220:LEU:HD22	1:A:241:LEU:HD13	1.90	0.54
1:B:335:SER:HA	1:B:358:LYS:HD2	1.88	0.54
1:A:111:VAL:HA	1:A:116:VAL:HG11	1.88	0.53
1:A:425:ALA:O	1:A:429:THR:HG22	2.09	0.53
1:A:500:ILE:HD12	1:A:519:GLN:NE2	2.23	0.53
1:A:155:VAL:HG11	1:A:175:LEU:HD21	1.89	0.53
1:A:549:VAL:HG12	1:A:579:LEU:HD12	1.91	0.53
1:A:177:LEU:HD11	1:B:195:ARG:HG3	1.90	0.53
1:A:537:LEU:HD23	1:A:560:ILE:HG23	1.91	0.53
1:A:382:LEU:CD2	1:B:242:CYS:HA	2.39	0.52
1:A:70:HIS:HB3	4:A:1024:HOH:O	2.09	0.52
1:B:553:LYS:HD3	1:B:579:LEU:HB3	1.91	0.52
1:B:280:LEU:HD23	1:B:285:ALA:HA	1.91	0.52
1:A:254:LEU:HD21	1:A:421:VAL:HA	1.91	0.51
1:A:473:ARG:HD2	1:A:495:TYR:OH	2.09	0.51
1:A:431:SER:HB3	1:A:432:ASN:HA	1.91	0.51
1:A:61:MET:HG3	1:A:62:LEU:HG	1.93	0.51
1:A:271:PRO:HD2	1:A:429:THR:HB	1.92	0.51
1:B:555:ILE:HG13	1:B:555:ILE:O	2.10	0.51
1:A:425:ALA:O	1:A:429:THR:HG23	2.11	0.50
1:B:118:VAL:HG11	1:B:198:TYR:OH	2.11	0.50
1:A:371:ASN:ND2	4:A:1004:HOH:O	2.44	0.50
1:B:369:ASP:OD2	1:B:371:ASN:OD1	2.30	0.50
1:A:268:HIS:HE1	3:A:701:93A:N15	2.09	0.50
1:A:495:TYR:CE1	1:A:519:GLN:HA	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:THR:HG22	1:B:72:ALA:HB3	1.93	0.49
1:B:406:THR:HG23	1:B:583:LEU:HB3	1.94	0.49
1:A:53:SER:HA	1:A:58:PHE:O	2.11	0.49
4:A:1031:HOH:O	1:B:174:HIS:HB3	2.12	0.49
1:A:406:THR:HG23	1:A:583:LEU:HB3	1.95	0.49
1:A:454:HIS:HE1	1:B:313:MET:SD	2.35	0.49
1:A:542:PHE:CD1	1:A:566:SER:HB2	2.47	0.49
1:A:140:HIS:HB3	1:A:176:ALA:HB2	1.94	0.49
1:B:253:GLU:OE2	1:B:428:TYR:OH	2.22	0.49
1:B:432:ASN:HD21	1:B:455:CYS:HB2	1.78	0.48
1:B:10:LEU:HD13	1:B:42:LEU:HD11	1.95	0.48
1:B:10:LEU:CD1	1:B:42:LEU:HD11	2.42	0.48
1:A:320:ILE:HD13	1:A:334:ILE:HG12	1.96	0.48
1:B:58:PHE:CD2	1:B:70:HIS:HE1	2.30	0.48
1:A:528:TRP:O	1:A:531:LYS:HB2	2.14	0.48
1:B:297:LEU:CD2	1:B:333:ILE:HD11	2.44	0.48
1:B:417:ARG:O	1:B:421:VAL:HG23	2.13	0.48
1:B:43:ARG:HH22	1:B:51:ASP:CG	2.17	0.48
1:A:394:ALA:CB	1:A:588:LEU:HD11	2.43	0.47
1:B:300:LEU:HD11	1:B:322:LEU:HB3	1.96	0.47
1:B:434:VAL:CG1	1:B:537:LEU:CD1	2.81	0.47
1:A:319:PHE:CE2	1:A:341:GLY:HA3	2.49	0.47
1:A:133:LEU:CD2	1:A:147:CYS:HB3	2.44	0.47
1:B:543:PHE:HE2	1:B:560:ILE:HG21	1.79	0.47
1:A:495:TYR:HE1	1:A:519:GLN:HA	1.80	0.47
1:B:207:LEU:HD21	1:B:241:LEU:CD1	2.45	0.47
1:A:285:ALA:HB2	1:A:294:HIS:CD2	2.49	0.47
1:B:426:VAL:CG1	1:B:540:ASP:HB3	2.44	0.46
1:A:345:PRO:HA	1:A:368:MET:O	2.15	0.46
1:B:467:TRP:O	1:B:470:HIS:HB2	2.14	0.46
1:A:267:LYS:HG2	1:B:450:GLN:HB3	1.98	0.46
1:A:229:LEU:HD13	1:A:368:MET:HE2	1.96	0.46
1:A:509:TRP:O	1:A:512:MET:HB2	2.16	0.46
1:A:7:LEU:CD1	1:A:95:SER:HB2	2.46	0.46
1:A:268:HIS:CD2	1:B:593:HIS:NE2	2.77	0.45
1:A:25:LEU:HA	1:A:28:LEU:HD12	1.97	0.45
1:A:426:VAL:HA	1:A:429:THR:HG23	1.99	0.45
1:A:7:LEU:CB	1:A:33:ILE:HD11	2.45	0.45
1:A:358:LYS:HB3	1:A:363:TYR:HB2	1.98	0.45
1:A:328:VAL:HB	1:A:329:PRO:HD3	1.99	0.45
1:A:476:SER:O	1:A:478:LYS:HD2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:THR:HG23	1:B:175:LEU:HD23	1.99	0.45
1:B:248:TRP:CE2	1:B:368:MET:HG2	2.52	0.45
1:B:28:LEU:HG	1:B:165:LYS:HB3	1.98	0.45
1:B:492:ILE:O	1:B:496:VAL:HG22	2.17	0.45
1:B:9:LEU:HD22	1:B:73:VAL:HG13	1.99	0.44
1:A:42:LEU:HB2	1:A:49:VAL:HG11	1.99	0.44
1:A:431:SER:O	1:A:448:GLY:HA2	2.18	0.44
1:A:426:VAL:HG23	1:A:435:CYS:CB	2.48	0.44
1:B:328:VAL:HB	1:B:329:PRO:HD3	1.98	0.44
1:B:520:LEU:HB2	1:B:525:LYS:HE3	2.00	0.44
1:B:434:VAL:HG13	1:B:537:LEU:HD11	1.92	0.44
1:A:431:SER:CB	1:A:432:ASN:HA	2.46	0.44
1:A:11:SER:O	1:A:102:CYS:HA	2.18	0.43
1:A:168:SER:HB2	1:A:170:GLU:OE2	2.19	0.43
1:B:308:ARG:CZ	1:B:337:GLU:HB3	2.48	0.43
1:A:23:ARG:HG3	1:A:23:ARG:HH11	1.83	0.43
1:A:537:LEU:HD21	1:A:543:PHE:HZ	1.81	0.43
1:B:549:VAL:HG21	1:B:572:VAL:HG13	2.01	0.43
1:B:33:ILE:HA	1:B:50:ARG:O	2.19	0.43
1:A:207:LEU:HG	1:A:219:GLN:HA	2.00	0.43
1:B:12:VAL:HG23	1:B:15:LYS:HD2	2.00	0.43
1:A:374:PRO:HD3	1:B:383:TYR:CE1	2.54	0.43
1:A:106:PRO:HB2	1:A:109:LYS:HB3	1.99	0.43
1:A:229:LEU:HD13	1:A:368:MET:HE1	2.00	0.43
1:B:26:ASN:O	1:B:28:LEU:O	2.36	0.42
1:A:71:PRO:HG3	1:B:70:HIS:CE1	2.55	0.42
1:A:268:HIS:CE1	3:A:701:93A:N15	2.86	0.42
1:A:401:PHE:HA	1:A:584:ILE:HG21	1.99	0.42
1:A:128:GLY:HA3	3:A:600:93A:C18	2.50	0.42
1:B:461:ASP:O	1:B:464:ASN:HB2	2.19	0.42
1:B:342:VAL:HG11	1:B:355:LEU:HD13	2.01	0.42
1:A:12:VAL:HG23	1:A:15:LYS:HD3	2.00	0.42
1:B:140:HIS:HB3	1:B:176:ALA:HB2	2.01	0.42
1:A:432:ASN:HB2	1:A:452:ARG:NH2	2.34	0.42
1:A:277:GLY:HA3	1:A:299:PRO:O	2.19	0.42
1:A:398:ARG:HH22	1:A:413:GLU:CD	2.23	0.42
1:B:145:VAL:O	1:B:175:LEU:HG	2.20	0.42
1:B:466:TRP:CE2	1:B:470:HIS:HE1	2.36	0.42
1:A:308:ARG:HD3	1:A:337:GLU:HG2	2.00	0.41
1:B:209:TYR:CG	1:B:213:PRO:HA	2.56	0.41
1:B:107:PHE:CZ	1:B:111:VAL:HG11	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:68:THR:OG1	3:B:601:93A:N12	2.43	0.41
1:A:10:LEU:O	1:A:34:ALA:HA	2.20	0.41
1:A:308:ARG:CD	1:A:337:GLU:CG	2.98	0.41
1:A:356:SER:O	1:A:361:GLY:HA2	2.21	0.41
1:A:107:PHE:O	1:A:111:VAL:HG22	2.21	0.41
1:A:461:ASP:O	1:A:464:ASN:HB2	2.21	0.41
1:A:506:LEU:HD12	1:A:510:GLN:HG3	2.03	0.41
1:A:434:VAL:CG1	1:A:537:LEU:HD11	2.51	0.41
1:A:404:ILE:HG21	1:A:409:LYS:HA	2.03	0.40
1:A:464:ASN:HD22	1:A:555:ILE:HD13	1.85	0.40
1:B:326:CYS:O	1:B:348:GLU:HG3	2.22	0.40
1:A:19:VAL:HG12	1:A:45:ALA:HB2	2.02	0.40
1:A:401:PHE:HE2	1:A:420:ILE:HG13	1.87	0.40
1:A:574:GLU:HA	1:A:574:GLU:OE1	2.22	0.40
1:B:128:GLY:O	1:B:132:LEU:HG	2.22	0.40
1:B:104:LEU:HD21	1:B:133:LEU:HD12	2.02	0.40
1:B:404:ILE:HG21	1:B:409:LYS:HA	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:LYS:NZ	1:A:587:ASN:OD1[1_655]	2.14	0.06

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	588/593 (99%)	570 (97%)	15 (3%)	3 (0%)	29	22
1	B	588/593 (99%)	568 (97%)	20 (3%)	0	100	100
All	All	1176/1186 (99%)	1138 (97%)	35 (3%)	3 (0%)	41	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	53	SER
1	A	114	PRO
1	A	269	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	484/485 (100%)	466 (96%)	18 (4%)	34	31
1	B	484/485 (100%)	466 (96%)	18 (4%)	34	31
All	All	968/970 (100%)	932 (96%)	36 (4%)	34	31

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	19	VAL
1	A	33	ILE
1	A	51	ASP
1	A	58	PHE
1	A	123	GLU
1	A	199	SER
1	A	324	ASP
1	A	339	SER
1	A	360	ASN
1	A	413	GLU
1	A	429	THR
1	A	433	SER
1	A	451	SER
1	A	461	ASP
1	A	478	LYS
1	A	502	GLU
1	A	506	LEU
1	B	15	LYS
1	B	28	LEU

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Mol	Chain	Res	Type
1	B	47	LEU
1	B	80	ARG
1	B	95	SER
1	B	112	SER
1	B	123	GLU
1	B	124	LYS
1	B	153	SER
1	B	164	ASP
1	B	175	LEU
1	B	177	LEU
1	B	324	ASP
1	B	352	LEU
1	B	371	ASN
1	B	432	ASN
1	B	484	LYS
1	B	547	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	HIS
1	A	86	ASN
1	A	90	ASN
1	A	268	HIS
1	A	371	ASN
1	A	464	ASN
1	A	490	ASN
1	A	494	GLN
1	A	558	GLN
1	B	70	HIS
1	B	371	ASN
1	B	432	ASN
1	B	464	ASN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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	93A	A	701	-	20,27,27	4.06	8 (40%)	23,43,43	4.10	8 (34%)
3	93A	A	600	-	20,27,27	4.09	10 (50%)	23,43,43	5.38	11 (47%)
3	93A	B	700	-	20,27,27	4.06	9 (45%)	23,43,43	4.63	8 (34%)
3	93A	B	601	-	20,27,27	4.32	10 (50%)	23,43,43	5.25	10 (43%)

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	93A	A	701	-	-	0/6/40/40	0/2/3/3
3	93A	A	600	-	-	0/6/40/40	0/2/3/3
3	93A	B	700	-	-	0/6/40/40	0/2/3/3
3	93A	B	601	-	-	5/6/40/40	0/2/3/3

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	600	93A	O21-S16	13.06	1.62	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	93A	O21-S16	13.00	1.62	1.43
3	B	601	93A	O21-S16	12.55	1.61	1.43
3	B	700	93A	O21-S16	12.52	1.61	1.43
3	B	601	93A	O20-S16	10.74	1.59	1.43
3	A	600	93A	O20-S16	8.67	1.56	1.43
3	A	701	93A	O20-S16	8.66	1.56	1.43
3	B	700	93A	O20-S16	8.52	1.55	1.43
3	A	701	93A	O5'-C5'	-4.96	1.25	1.44
3	B	700	93A	C11-N12	4.92	1.43	1.35
3	B	700	93A	O5'-C5'	-4.92	1.25	1.44
3	B	601	93A	P-O1P	4.78	1.66	1.50
3	A	600	93A	P-O1P	4.54	1.65	1.50
3	A	701	93A	C13-C18	4.21	1.56	1.47
3	B	700	93A	C13-C18	4.21	1.56	1.47
3	A	600	93A	P-O2P	3.76	1.69	1.54
3	A	701	93A	C11-N12	3.72	1.41	1.35
3	B	601	93A	P-O2P	3.68	1.69	1.54
3	A	600	93A	C11-N12	3.64	1.41	1.35
3	B	601	93A	O1-C2	-3.36	1.37	1.45
3	A	600	93A	O5'-C5'	-3.13	1.32	1.44
3	B	601	93A	O5'-C5'	-3.10	1.32	1.44
3	B	601	93A	C13-C18	3.03	1.53	1.47
3	B	601	93A	C11-N12	2.98	1.40	1.35
3	A	600	93A	C4-C5	-2.75	1.49	1.53
3	B	700	93A	C5'-C2	-2.67	1.43	1.51
3	A	701	93A	C5'-C2	-2.59	1.43	1.51
3	A	701	93A	P-O1P	2.55	1.58	1.50
3	B	700	93A	O1-C2	-2.48	1.39	1.45
3	B	601	93A	C5'-C2	-2.44	1.44	1.51
3	B	601	93A	C4-C5	-2.41	1.50	1.53
3	B	700	93A	P-O1P	2.39	1.58	1.50
3	A	600	93A	C13-C18	2.35	1.52	1.47
3	A	701	93A	O1-C2	-2.34	1.39	1.45
3	A	600	93A	O1-C2	-2.31	1.39	1.45
3	A	600	93A	O1-C5	2.12	1.44	1.41
3	B	700	93A	C3-C4	-2.10	1.47	1.53

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	601	93A	O20-S16-O21	-17.97	96.28	118.85
3	A	600	93A	O20-S16-O21	-16.54	98.08	118.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	700	93A	O20-S16-O21	-16.15	98.57	118.85
3	A	600	93A	P-O5'-C5'	14.84	159.17	118.30
3	A	701	93A	O20-S16-O21	-12.51	103.14	118.85
3	B	601	93A	P-O5'-C5'	11.93	151.16	118.30
3	B	700	93A	P-O5'-C5'	9.54	144.56	118.30
3	B	601	93A	O5'-C5'-C2	9.16	140.53	108.99
3	A	701	93A	P-O5'-C5'	8.54	141.82	118.30
3	A	701	93A	O5'-C5'-C2	8.42	137.96	108.99
3	A	600	93A	O5'-C5'-C2	8.20	137.21	108.99
3	B	700	93A	O5'-C5'-C2	8.11	136.90	108.99
3	A	600	93A	C11-N12-C13	-5.16	93.16	102.99
3	B	700	93A	C11-N12-C13	-4.61	94.21	102.99
3	A	701	93A	C11-N12-C13	-4.45	94.52	102.99
3	B	601	93A	O6-C3-C2	-4.14	99.07	111.05
3	A	600	93A	O3P-P-O5'	-4.12	95.78	106.73
3	B	601	93A	C11-N12-C13	-4.02	95.34	102.99
3	B	700	93A	C5'-C2-C3	-3.83	100.83	115.18
3	A	701	93A	O1-C2-C5'	-3.71	97.16	109.37
3	A	600	93A	O19-C18-C13	-3.56	117.41	124.37
3	B	601	93A	O19-C18-C13	-3.48	117.58	124.37
3	A	701	93A	C4-C3-C2	-3.43	95.97	102.64
3	A	701	93A	C5'-C2-C3	-3.38	102.53	115.18
3	A	600	93A	O19-C18-N17	3.23	125.19	120.82
3	A	600	93A	C5'-C2-C3	-3.13	103.47	115.18
3	B	601	93A	O19-C18-N17	3.04	124.92	120.82
3	A	600	93A	O1-C2-C3	2.92	110.90	105.11
3	B	700	93A	O1-C2-C5'	-2.90	99.83	109.37
3	A	701	93A	O5'-P-O1P	-2.70	98.91	106.47
3	A	600	93A	O6-C3-C2	-2.35	104.24	111.05
3	B	601	93A	C5'-C2-C3	-2.28	106.65	115.18
3	B	700	93A	O5'-P-O1P	2.25	112.80	106.47
3	B	601	93A	O3P-P-O5'	-2.21	100.84	106.73
3	B	700	93A	C4-C3-C2	-2.10	98.56	102.64
3	A	600	93A	O3P-P-O1P	2.09	118.85	110.68
3	B	601	93A	O1-C2-C3	2.03	109.13	105.11

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	601	93A	O1-C2-C5'-O5'
3	B	601	93A	C3-C2-C5'-O5'

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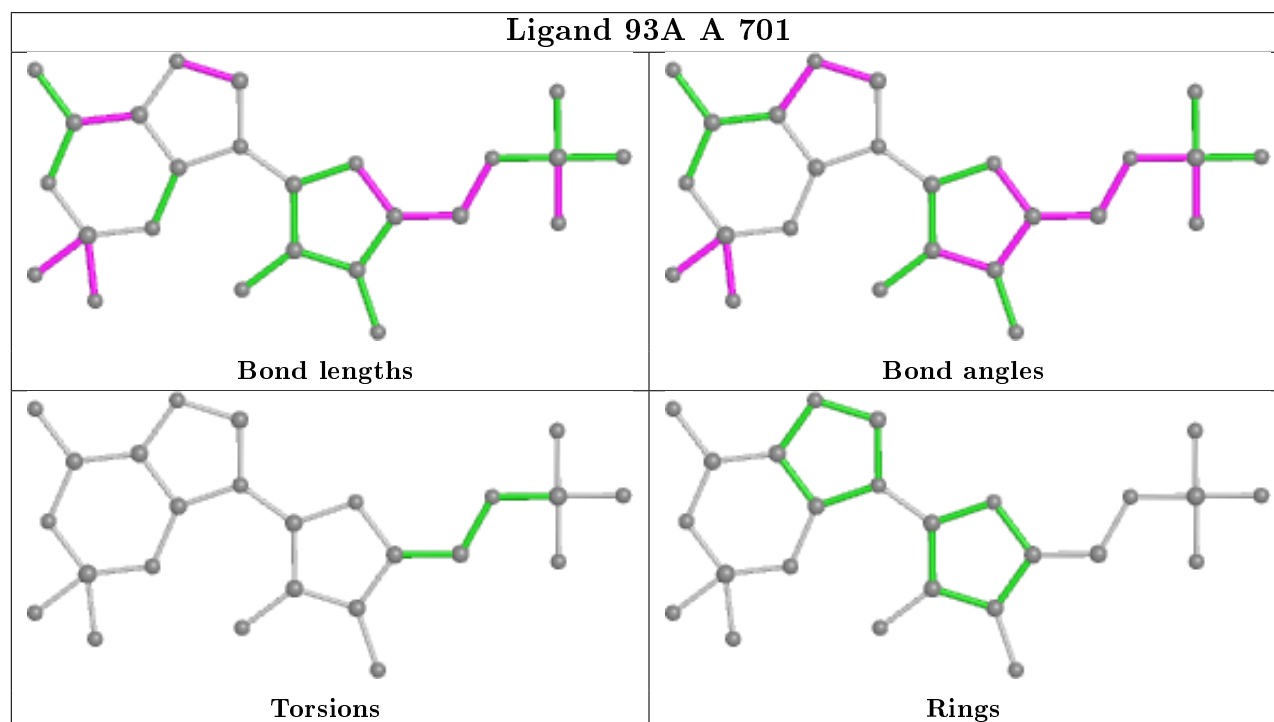
Mol	Chain	Res	Type	Atoms
3	B	601	93A	C5'-O5'-P-O2P
3	B	601	93A	C5'-O5'-P-O3P
3	B	601	93A	C5'-O5'-P-O1P

There are no ring outliers.

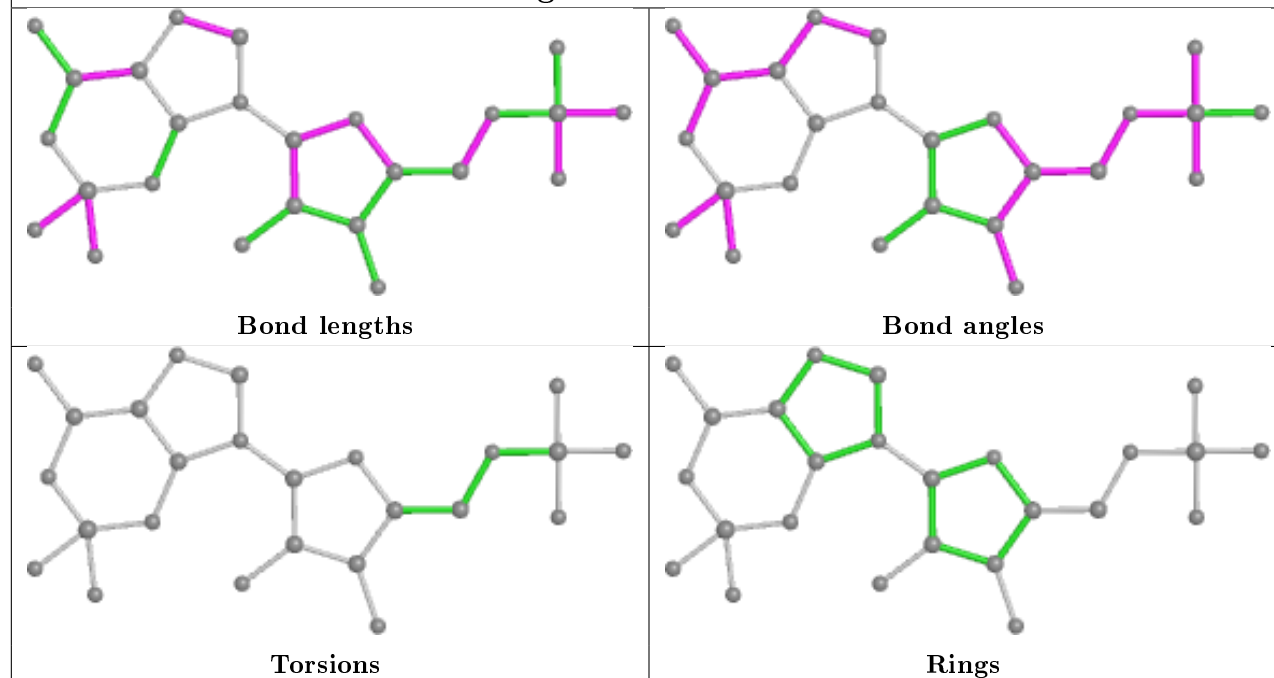
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	701	93A	2	0
3	A	600	93A	1	0
3	B	601	93A	1	0

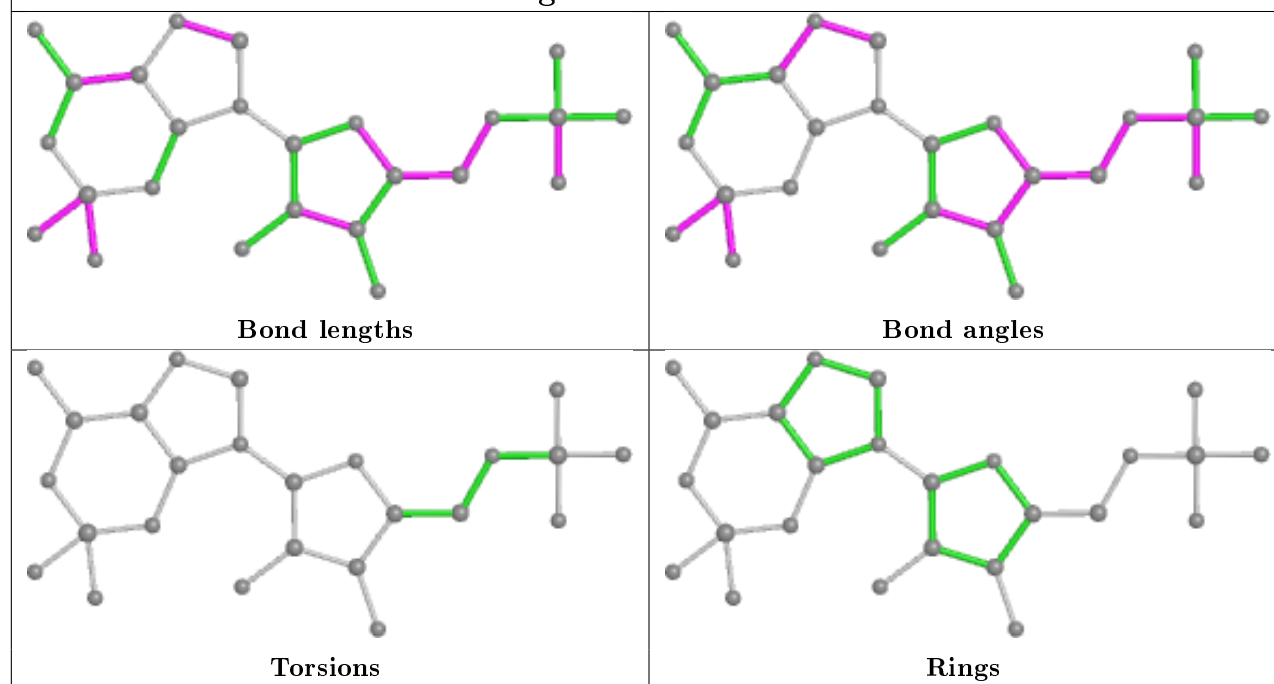
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

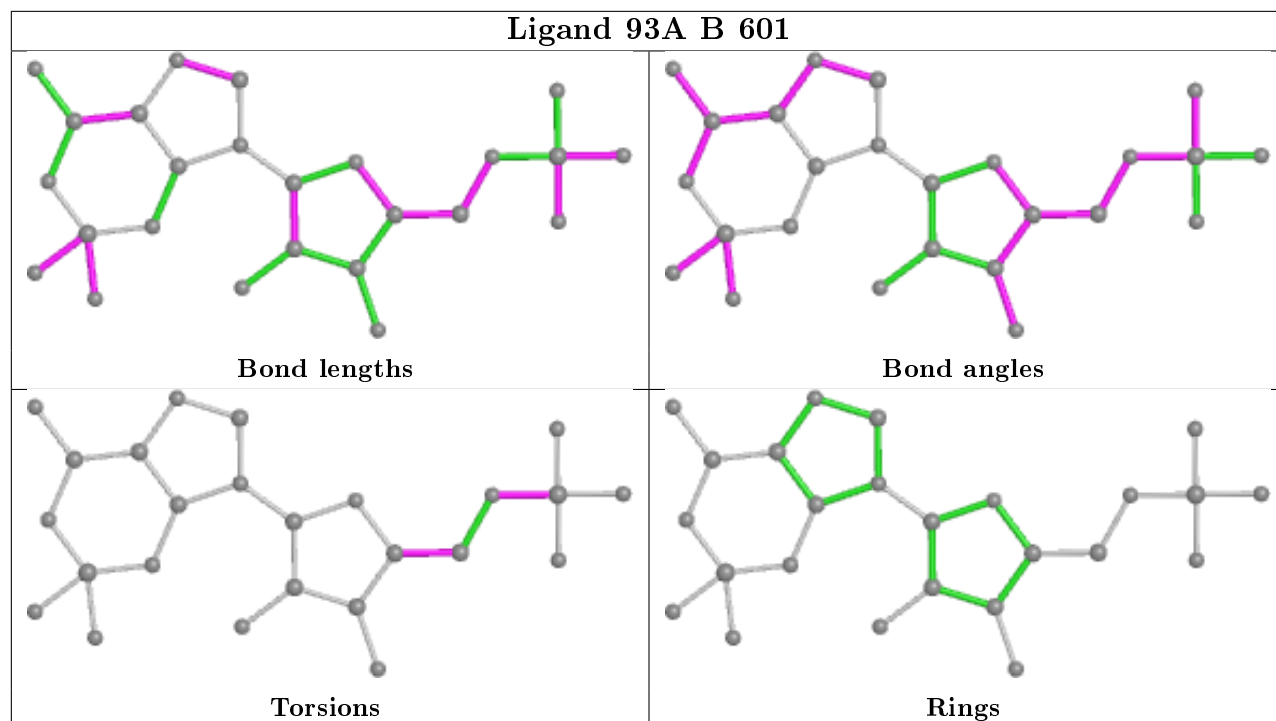


Ligand 93A A 600



Ligand 93A B 700





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	590/593 (99%)	0.28	20 (3%) 45 45	21, 40, 79, 96	0
1	B	590/593 (99%)	0.25	23 (3%) 39 39	27, 45, 74, 105	0
All	All	1180/1186 (99%)	0.26	43 (3%) 42 42	21, 42, 77, 105	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	82	ILE	4.6
1	B	481	ALA	4.5
1	B	87	ALA	4.3
1	A	114	PRO	4.2
1	B	114	PRO	4.2
1	B	482	GLY	4.2
1	B	483	VAL	4.0
1	A	163	LYS	3.9
1	B	115	GLY	3.8
1	A	481	ALA	3.6
1	A	483	VAL	3.5
1	A	514	GLU	3.5
1	B	479	PHE	3.1
1	B	512	MET	3.0
1	A	482	GLY	2.9
1	A	515	GLU	2.9
1	B	504	GLU	2.9
1	A	41	ALA	2.7
1	A	39	ALA	2.7
1	A	28	LEU	2.5
1	A	42	LEU	2.5
1	B	515	GLU	2.5
1	B	117	THR	2.4
1	A	83	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	54	ASP	2.4
1	B	520	LEU	2.3
1	B	503	ASP	2.3
1	B	131	ALA	2.3
1	A	182	HIS	2.3
1	A	19	VAL	2.2
1	A	479	PHE	2.2
1	B	513	PHE	2.2
1	B	510	GLN	2.2
1	A	127	ILE	2.1
1	A	429	THR	2.1
1	A	165	LYS	2.1
1	B	4	ARG	2.1
1	B	502	GLU	2.1
1	B	354	ILE	2.1
1	B	511	ALA	2.1
1	A	480	LYS	2.0
1	B	113	SER	2.0
1	B	182	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	93A	A	600	25/25	0.93	0.12	43,46,50,52	0
3	93A	B	601	25/25	0.96	0.11	34,38,43,44	0
3	93A	A	701	25/25	0.97	0.12	29,37,42,42	0
3	93A	B	700	25/25	0.97	0.13	38,44,48,49	0

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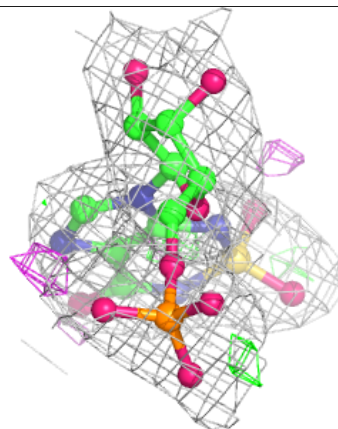
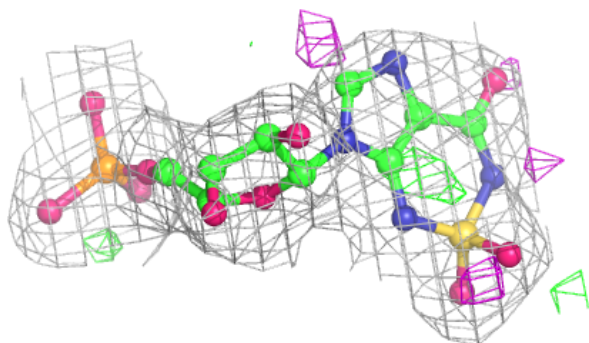
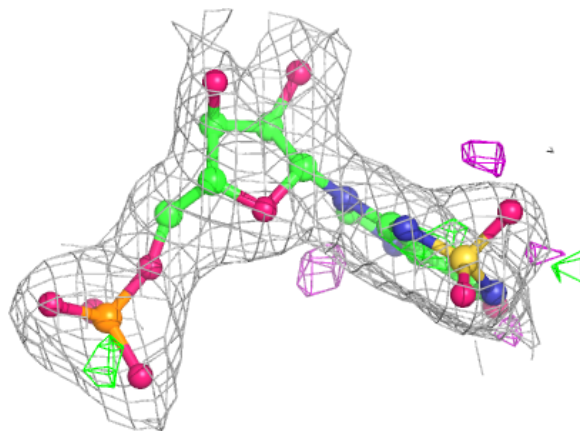
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	K	A	901	1/1	0.99	0.10	26,26,26,26	0
2	K	B	902	1/1	0.99	0.08	35,35,35,35	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

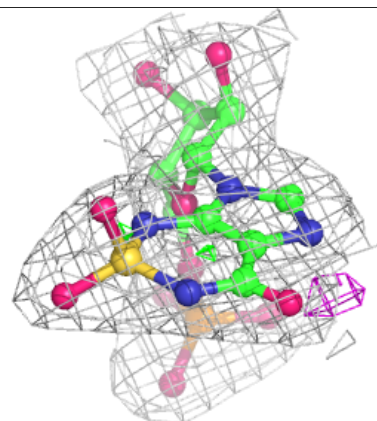
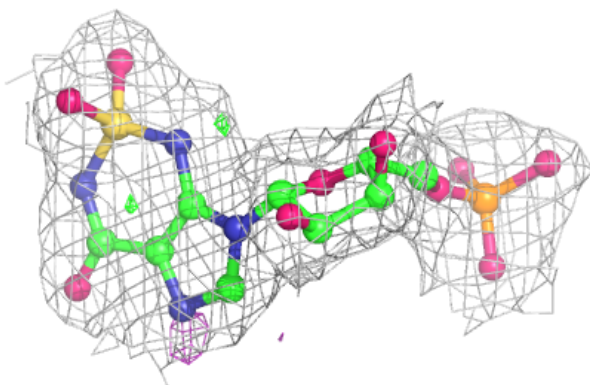
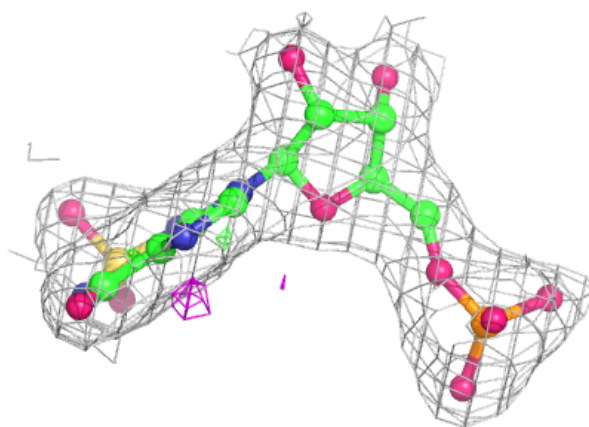
Electron density around 93A A 600:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

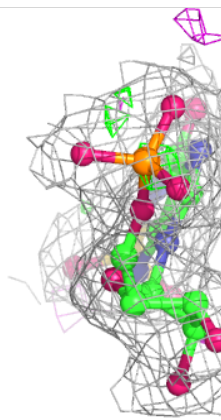
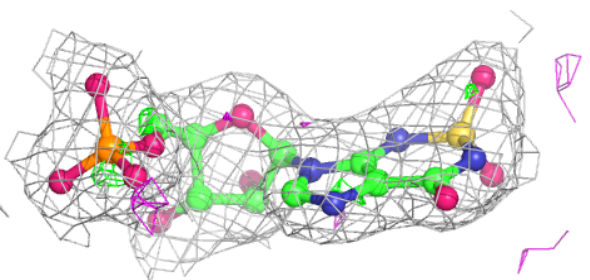
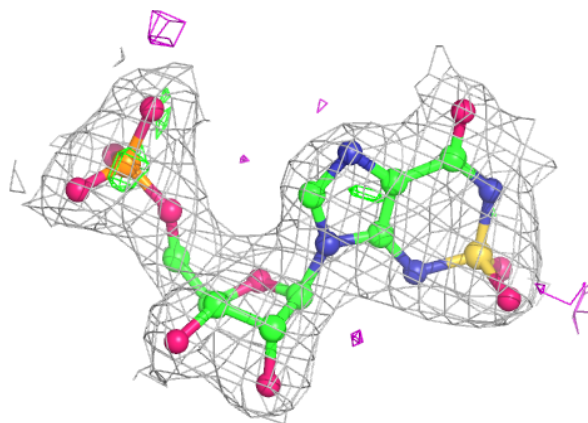


Electron density around 93A B 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

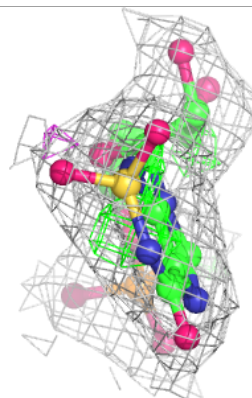
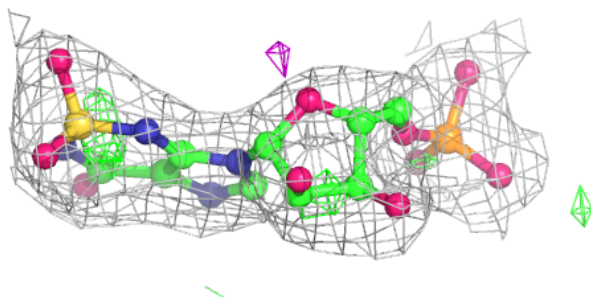
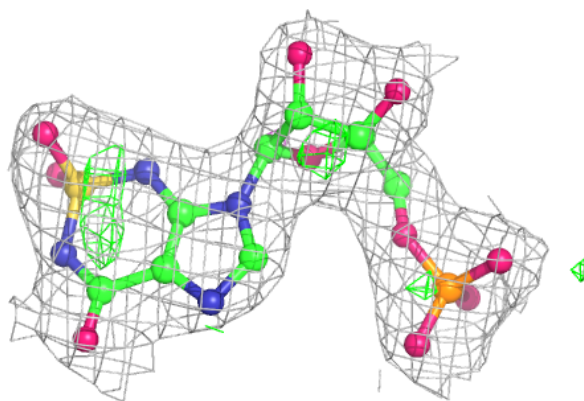
**Electron density around 93A A 701:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 93A B 700:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.