



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 11:21 AM UTC

PDB ID : 8E4X / pdb_00008e4x
Title : Human Adenosine Deaminase Acting on dsRNA (ADAR2-R2D) bound to dsRNA containing a G:3-deaza dA pair adjacent to the target site
Authors : Fisher, A.J.; Mendoza, H.G.
Deposited on : 2022-08-19
Resolution : 2.80 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)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

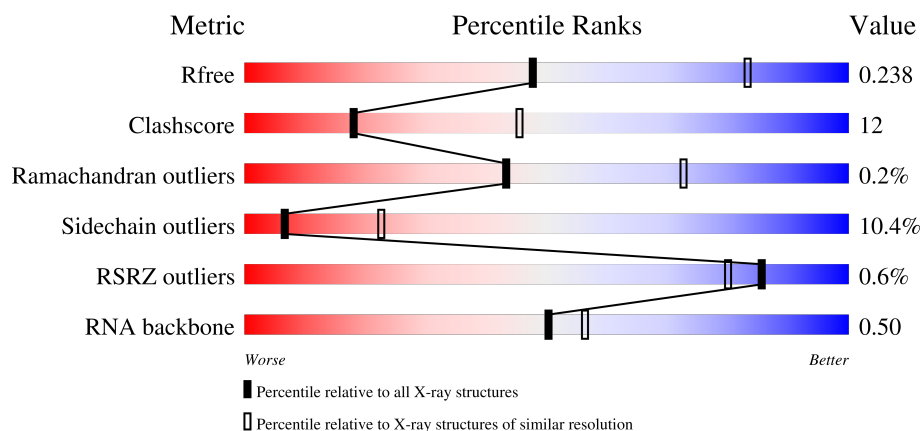
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)
RNA backbone	3983	1114 (3.00-2.60)

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 $\leq 5\%$. 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	488	58% 19% • 21%
1	B	488	60% 27% 6% 7%
2	C	32	50% 47% •
3	D	32	69% 28% •

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8029 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 Double-stranded RNA-specific editase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	384	Total	C	N	O	S	0	0	0
			3019	1904	550	554	11			
1	B	454	Total	C	N	O	S	0	0	0
			3539	2234	639	653	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	214	GLY	-	expression tag	UNP P78563
A	488	GLN	GLU	engineered mutation	UNP P78563
B	214	GLY	-	expression tag	UNP P78563
B	488	GLN	GLU	engineered mutation	UNP P78563

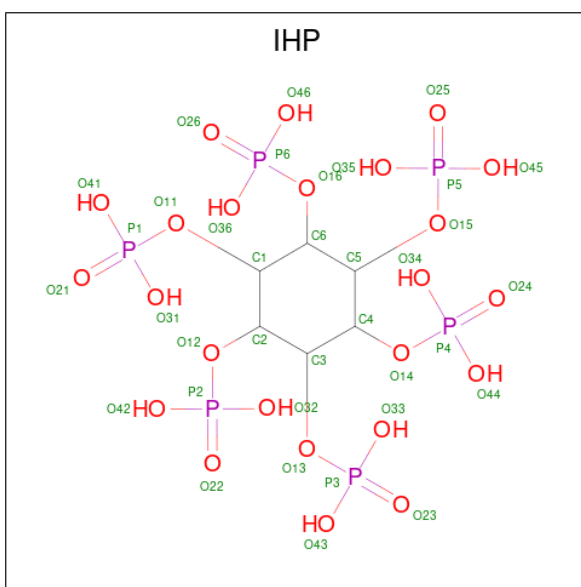
- Molecule 2 is a RNA chain called RNA (5-R(*GP*CP*UP*CP*GP*CP*GP*AP*UP*GP*CP*GP*(8AZ)P*GP*AP*GP*GP*GP*CP*UP*CP*UP*GP*AP*UP*AP*GP*CP*UP*AP*CP*G)-3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	32	Total	C	N	O	P	0	0	0
			685	305	126	223	31			

- Molecule 3 is a RNA chain called RNA(5-R(*CP*GP*UP*AP*GP*CP*UP*AP*UP*CP*AP*GP*AP*GP*CP*CP*CP*CP*CP*(4DU)*GP*CP*AP*UP*CP*GP*CP*GP*AP*GP*CP)-3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	32	Total	C	N	O	P	0	0	0
			674	304	121	218	31			

- Molecule 4 is INOSITOL HEXAKISPHOSPHATE (CCD ID: IHP) (formula: C₆H₁₈O₂₄P₆).



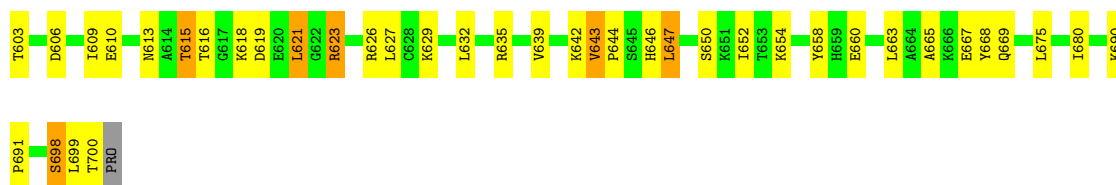
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			36	6	24	6		
4	B	1	Total	C	O	P	0	0
			36	6	24	6		

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Zn	0	0
			1	1		
5	B	1	Total	Zn	0	0
			1	1		

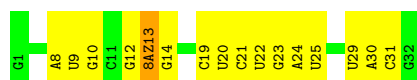
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	19	Total	O	0	0
			19	19		
6	B	9	Total	O	0	0
			9	9		
6	C	6	Total	O	0	0
			6	6		
6	D	4	Total	O	0	0
			4	4		



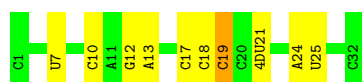
- Molecule 2: RNA (5-R(*GP*CP*UP*CP*GP*CP*GP*AP*UP*GP*CP*GP*(8AZ)P*GP*AP*GP*GP*GP*CP*UP*CP*UP*GP*AP*UP*AP*GP*CP*UP*AP*CP*G)-3)

Chain C: 50% 47%



- Molecule 3: RNA(5-R(*CP*GP*UP*AP*GP*CP*UP*AP*UP*CP*AP*GP*AP*GP*CP*CP*CP*CP*CP*(4DU)*GP*CP*AP*UP*CP*GP*CP*GP*AP*GP*C)-3)

Chain D: 69% 28%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	169.91Å 63.24Å 142.65Å 90.00° 118.07° 90.00°	Depositor
Resolution (Å)	58.27 – 2.80 58.27 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.9 (58.27-2.80) 94.9 (58.27-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.195 , 0.239 0.194 , 0.238	Depositor DCC
R_{free} test set	1582 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	81.8	Xtriage
Anisotropy	0.559	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 77.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8029	wwPDB-VP
Average B, all atoms (Å ²)	110.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 8AZ, 4DU, IHP, ZN

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	0.59	0/3081	0.96	4/4164 (0.1%)
1	B	0.45	0/3609	0.65	0/4874
2	C	0.41	0/740	0.61	0/1151
3	D	0.44	0/727	0.66	0/1128
All	All	0.50	0/8157	0.78	4/11317 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	657	VAL	N-CA-C	-5.99	99.50	108.12
1	A	594	LYS	N-CA-C	-5.24	102.63	110.23
1	A	498	SER	CA-C-N	5.19	131.31	121.97
1	A	498	SER	C-N-CA	5.19	131.31	121.97

There are no chirality outliers.

There are no planarity outliers.

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	3019	0	3051	71	0
1	B	3539	0	3567	104	0
2	C	685	0	337	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	674	0	351	8	0
4	A	36	0	6	1	0
4	B	36	0	6	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	19	0	0	1	0
6	B	9	0	0	0	0
6	C	6	0	0	0	0
6	D	4	0	0	0	0
All	All	8029	0	7318	186	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 12.

All (186) 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:654:LYS:HG3	1:A:655:PRO:HD2	1.59	0.84
1:B:300:HIS:CG	1:B:301:LEU:H	1.96	0.83
1:A:620:GLU:OE2	6:A:901:HOH:O	1.97	0.81
1:B:428:LYS:HD3	1:B:428:LYS:H	1.46	0.79
1:A:554:ASP:OD1	1:A:554:ASP:N	2.14	0.75
1:B:305:PRO:HG3	1:B:462:PRO:HB2	1.65	0.75
1:B:400:ARG:HD3	1:B:523:TRP:CE2	2.22	0.74
1:B:492:PRO:O	1:B:494:ARG:NH1	2.22	0.72
1:B:319:LEU:HB3	1:B:322:VAL:HG13	1.72	0.71
1:A:327:VAL:HG21	1:A:546:ILE:HD13	1.73	0.70
1:B:239:ILE:HA	1:B:242:GLU:HG2	1.74	0.69
1:B:308:GLN:OE1	1:B:308:GLN:N	2.26	0.69
1:A:464:LEU:HD13	1:A:464:LEU:H	1.61	0.66
1:B:281:LYS:NZ	2:C:19:C:OP1	2.28	0.66
1:B:698:SER:HB2	1:B:700:THR:HG22	1.76	0.66
1:B:325:ASP:HB3	1:B:329:ARG:HH21	1.61	0.65
1:B:449:SER:HB3	1:B:455:ARG:HG3	1.78	0.64
1:A:350:LYS:NZ	1:A:591:GLN:O	2.28	0.63
1:B:317:LEU:HB3	1:B:323:LEU:HD13	1.81	0.61
2:C:24:A:H2'	2:C:25:U:H6	1.66	0.61
1:B:348:ARG:HH21	1:B:598:PHE:HE2	1.49	0.61
1:A:451:CYS:SG	1:A:483:LYS:HE3	2.41	0.60
1:B:301:LEU:O	1:B:303:GLN:N	2.34	0.60
1:A:460:HIS:C	1:A:550:LEU:HD22	2.26	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:436:LEU:HD21	1:A:540:ILE:HG12	1.84	0.59
1:A:487:GLY:O	2:C:14:G:H1'	2.02	0.58
1:A:501:THR:HG22	1:B:491:ILE:HD13	1.86	0.58
1:B:501:THR:HA	1:B:691:PRO:HB3	1.85	0.58
3:D:12:G:H2'	3:D:13:A:H8	1.68	0.58
1:A:323:LEU:O	1:A:327:VAL:HG13	2.03	0.58
1:B:430:GLU:H	1:B:430:GLU:CD	2.12	0.58
1:A:657:VAL:HG22	1:A:660:GLU:HG2	1.86	0.58
1:A:319:LEU:HB3	1:A:322:VAL:HG23	1.86	0.57
1:B:443:HIS:HB3	1:B:544:SER:HB2	1.86	0.57
1:B:392:ASP:OD2	1:B:483:LYS:NZ	2.36	0.57
1:B:247:LEU:HD21	1:B:269:VAL:HB	1.85	0.57
1:B:393:CYS:HA	1:B:398:ILE:HD11	1.87	0.57
1:B:639:VAL:O	1:B:643:VAL:HG22	2.05	0.56
1:A:430:GLU:OE1	1:A:430:GLU:N	2.25	0.56
1:B:317:LEU:HD12	1:B:363:VAL:HG13	1.88	0.56
1:B:484:ILE:HG12	1:B:484:ILE:O	2.05	0.56
1:A:462:PRO:HB3	1:A:552:HIS:CE1	2.41	0.55
1:A:654:LYS:HG2	1:A:660:GLU:HB3	1.88	0.55
1:B:516:CYS:O	1:B:520:ILE:HG13	2.06	0.55
2:C:24:A:H2'	2:C:25:U:C6	2.42	0.55
1:A:514:MET:HG3	1:A:687:TRP:CE2	2.42	0.55
1:B:535:ILE:HG23	1:B:650:SER:HB2	1.89	0.55
1:B:300:HIS:CG	1:B:301:LEU:N	2.67	0.55
1:B:300:HIS:CD2	1:B:301:LEU:H	2.25	0.54
1:B:621:LEU:HD21	1:B:623:ARG:NH2	2.23	0.54
1:A:458:SER:HB2	1:A:552:HIS:CD2	2.44	0.53
2:C:23:G:H2'	2:C:24:A:H8	1.74	0.53
3:D:12:G:H2'	3:D:13:A:C8	2.44	0.53
1:B:303:GLN:HB3	1:B:306:SER:HB2	1.91	0.52
1:A:471:HIS:HB3	1:A:474:ARG:HG3	1.91	0.52
1:A:430:GLU:H	1:A:430:GLU:CD	2.07	0.52
1:B:400:ARG:HD3	1:B:523:TRP:CZ2	2.44	0.52
1:B:365:ASP:OD1	1:B:365:ASP:N	2.43	0.52
3:D:19:C:H5''	3:D:19:C:H6	1.73	0.52
1:B:239:ILE:O	1:B:243:LEU:HB2	2.09	0.52
1:B:294:ALA:HA	1:B:298:ASN:O	2.10	0.52
1:B:319:LEU:O	1:B:322:VAL:HG22	2.10	0.52
1:B:644:PRO:HB2	1:B:646:HIS:ND1	2.25	0.52
1:B:422:LYS:HE3	1:B:423:ARG:NH1	2.25	0.51
1:A:316:GLN:HA	1:A:363:VAL:HG12	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:413:LEU:HD21	1:A:423:ARG:HB2	1.93	0.51
1:A:537:VAL:HG12	1:A:538:GLU:O	2.10	0.51
1:A:558:ARG:HA	1:A:562:GLN:HB2	1.92	0.51
1:B:267:VAL:HG12	1:B:274:PHE:HB2	1.92	0.51
1:B:485:GLU:HB3	1:B:510:ARG:HD2	1.93	0.51
1:B:525:VAL:O	1:B:541:TYR:OH	2.28	0.50
3:D:17:C:H2'	3:D:18:C:O4'	2.12	0.50
1:B:606:ASP:CG	1:B:626:ARG:HH12	2.19	0.50
1:A:364:LYS:HZ3	1:A:364:LYS:HB2	1.76	0.50
1:A:484:ILE:HG23	1:A:487:GLY:HA3	1.92	0.50
1:B:586:ASN:ND2	1:B:586:ASN:H	2.08	0.50
1:A:532:LEU:HB3	1:A:636:TRP:CD1	2.47	0.49
1:B:300:HIS:CE1	1:B:301:LEU:HG	2.48	0.49
1:A:394:HIS:CE1	1:A:483:LYS:HE2	2.48	0.49
1:B:430:GLU:OE2	1:B:435:ARG:HD3	2.12	0.49
1:A:448:THR:OG1	1:A:550:LEU:HD12	2.13	0.49
1:A:319:LEU:HD22	1:A:320:PRO:HD2	1.94	0.49
1:B:613:ASN:OD1	1:B:615:THR:HG23	2.13	0.48
2:C:30:A:H2'	2:C:31:C:C6	2.49	0.48
1:B:237:VAL:HG21	1:B:285:LYS:HG2	1.96	0.48
1:A:629:LYS:NZ	4:A:801:IHP:O42	2.42	0.48
1:B:417:ASN:HB2	1:B:420:ASP:OD2	2.14	0.48
1:A:419:ASP:HA	1:A:422:LYS:HE2	1.95	0.48
1:A:596:PRO:HB3	1:A:598:PHE:CE2	2.48	0.48
1:B:383:MET:HB2	1:B:511:LEU:HD11	1.96	0.47
1:A:481:ARG:HA	1:A:491:ILE:O	2.14	0.47
1:B:323:LEU:O	1:B:327:VAL:HG13	2.13	0.47
1:B:484:ILE:HD13	1:B:489:GLY:O	2.14	0.47
1:A:547:LEU:HD12	1:A:551:TYR:HA	1.97	0.47
3:D:18:C:H2'	3:D:19:C:O4'	2.15	0.47
1:B:528:ILE:C	1:B:528:ILE:HD12	2.39	0.47
1:A:510:ARG:HD2	1:B:593:GLY:HA2	1.97	0.47
1:B:317:LEU:HD23	1:B:441:GLN:OE1	2.15	0.47
1:A:463:ILE:O	1:A:466:GLU:HG3	2.14	0.46
1:B:663:LEU:HD23	1:B:663:LEU:HA	1.74	0.46
1:B:317:LEU:HD12	1:B:363:VAL:HG22	1.97	0.46
1:B:308:GLN:HA	1:B:551:TYR:OH	2.16	0.46
1:B:418:LYS:HA	1:B:421:GLN:HG3	1.98	0.46
2:C:8:A:H2'	2:C:9:U:C6	2.50	0.46
1:B:319:LEU:HD23	1:B:320:PRO:HD2	1.98	0.46
1:B:348:ARG:NH2	1:B:598:PHE:HE2	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:654:LYS:HB3	1:B:654:LYS:HE2	1.76	0.46
1:B:308:GLN:NE2	1:B:321:GLN:OE1	2.50	0.45
1:B:449:SER:HA	1:B:547:LEU:HD13	1.98	0.45
1:B:256:GLY:O	1:B:261:LYS:HE3	2.17	0.45
1:B:556:LEU:HD23	1:B:581:LEU:HB3	1.98	0.45
1:B:404:LEU:HD22	1:B:533:LEU:HD12	1.98	0.45
1:A:480:LEU:O	1:A:492:PRO:HA	2.16	0.45
2:C:23:G:H2'	2:C:24:A:C8	2.52	0.45
1:A:316:GLN:HB3	1:A:364:LYS:NZ	2.31	0.45
1:A:319:LEU:HB3	1:A:322:VAL:CG2	2.46	0.45
1:A:400:ARG:HD3	1:A:523:TRP:CE2	2.52	0.45
1:B:357:MET:HE2	1:B:367:LYS:HE2	1.99	0.45
1:B:433:GLY:HA2	1:B:647:LEU:HD13	1.99	0.45
1:A:413:LEU:HD23	1:A:424:SER:HA	1.97	0.44
2:C:20:U:C2	2:C:21:C:C5	3.06	0.44
1:B:305:PRO:HG3	1:B:462:PRO:CB	2.43	0.44
1:B:532:LEU:HA	1:B:532:LEU:HD23	1.68	0.44
1:A:463:ILE:HG13	1:A:466:GLU:H	1.82	0.44
1:A:467:PRO:HB3	1:A:470:ARG:HH21	1.81	0.44
1:A:501:THR:OG1	1:B:489:GLY:HA3	2.17	0.44
1:B:259:HIS:ND1	1:B:259:HIS:N	2.64	0.44
1:A:542:PHE:O	1:A:579:PRO:HB3	2.18	0.44
1:A:316:GLN:N	1:A:316:GLN:OE1	2.51	0.43
1:A:533:LEU:HD23	1:A:533:LEU:HA	1.85	0.43
1:B:561:TYR:CD2	1:B:578:LYS:HA	2.53	0.43
1:A:632:LEU:HA	1:A:632:LEU:HD23	1.65	0.43
2:C:21:C:H2'	2:C:22:U:H6	1.83	0.43
1:A:319:LEU:O	1:A:320:PRO:C	2.61	0.43
3:D:24:A:H2'	3:D:25:U:H6	1.84	0.43
1:B:403:LEU:O	1:B:407:LEU:HG	2.18	0.43
1:B:370:SER:HB2	1:B:402:SER:HB2	2.00	0.43
1:B:390:LEU:HD21	1:B:627:LEU:HB3	2.01	0.43
1:B:606:ASP:CG	1:B:626:ARG:NH1	2.77	0.43
1:B:626:ARG:H	1:B:626:ARG:HG2	1.54	0.43
1:B:647:LEU:HD22	1:B:647:LEU:HA	1.83	0.43
1:A:348:ARG:HG2	1:A:596:PRO:HG3	2.00	0.43
1:B:269:VAL:HG11	1:B:292:ALA:HB1	2.00	0.43
1:A:485:GLU:OE2	1:B:590:ARG:NH2	2.26	0.42
1:B:477:ARG:O	1:B:479:GLN:N	2.49	0.42
1:B:699:LEU:HD23	1:B:699:LEU:HA	1.83	0.42
1:A:600:VAL:HA	1:A:610:GLU:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:LEU:HD13	1:B:317:LEU:HA	1.69	0.42
1:B:616:THR:O	1:B:618:LYS:HG2	2.20	0.42
1:A:480:LEU:HA	1:A:480:LEU:HD23	1.77	0.42
1:A:531:SER:HB3	1:A:668:TYR:CE1	2.54	0.42
1:A:481:ARG:NH1	1:A:492:PRO:HD3	2.35	0.42
1:A:509:GLU:OE2	1:B:488:GLN:N	2.48	0.42
1:B:300:HIS:NE2	1:B:301:LEU:HG	2.33	0.42
1:B:307:ARG:HA	1:B:308:GLN:OE1	2.19	0.42
1:B:349:ARG:HD2	1:B:349:ARG:C	2.44	0.42
1:A:316:GLN:HA	1:A:363:VAL:H	1.85	0.42
1:B:239:ILE:HG13	1:B:242:GLU:OE2	2.19	0.42
1:A:666:LYS:HA	1:A:666:LYS:HD3	1.85	0.42
1:B:325:ASP:HB3	1:B:329:ARG:NH2	2.32	0.42
1:B:602:TRP:HB2	1:B:609:ILE:HG22	2.02	0.42
1:A:600:VAL:HG12	1:A:611:VAL:HG22	2.01	0.42
1:B:629:LYS:HG2	1:B:658:TYR:CE2	2.55	0.42
3:D:24:A:H2'	3:D:25:U:C6	2.55	0.42
1:B:526:VAL:HB	1:B:668:TYR:CE2	2.55	0.41
1:B:635:ARG:HA	1:B:635:ARG:HD2	1.87	0.41
1:A:350:LYS:HA	1:A:350:LYS:HD3	1.73	0.41
1:A:352:LEU:HD12	1:A:372:SER:O	2.21	0.41
1:A:694:GLN:HG3	1:A:695:ASP:OD1	2.20	0.41
1:B:562:GLN:OE1	1:B:562:GLN:HA	2.20	0.41
2:C:12:G:N3	2:C:12:G:H2'	2.35	0.41
1:B:570:LEU:HD23	1:B:570:LEU:HA	1.82	0.41
1:A:639:VAL:O	1:A:643:VAL:HG23	2.20	0.41
1:A:646:HIS:CE1	1:A:647:LEU:HG	2.56	0.41
1:B:255:SER:O	1:B:261:LYS:HA	2.21	0.41
1:B:256:GLY:HA3	1:B:261:LYS:HA	2.01	0.41
1:B:665:ALA:O	1:B:669:GLN:HG2	2.21	0.41
3:D:21:4DU:H14	3:D:21:4DU:H8	1.87	0.41
1:A:485:GLU:HG2	1:A:511:LEU:H	1.86	0.40
1:A:507:GLN:HB2	1:B:455:ARG:NH1	2.36	0.40
1:B:567:ILE:HD12	1:B:675:LEU:HA	2.03	0.40
1:A:330:LEU:HB3	1:A:368:VAL:CG2	2.50	0.40
1:B:383:MET:HE3	1:B:502:TRP:CE3	2.56	0.40
1:B:610:GLU:OE2	1:B:626:ARG:HG2	2.21	0.40
1:A:630:HIS:CE1	1:A:699:LEU:HD22	2.55	0.40
1:A:556:LEU:O	1:A:560:MET:HG2	2.21	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	382/488 (78%)	370 (97%)	11 (3%)	1 (0%)	36	66
1	B	450/488 (92%)	416 (92%)	33 (7%)	1 (0%)	43	72
All	All	832/976 (85%)	786 (94%)	44 (5%)	2 (0%)	43	72

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	486	SER
1	A	499	ILE

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	329/415 (79%)	313 (95%)	16 (5%)	22	55
1	B	384/415 (92%)	326 (85%)	58 (15%)	3	10
All	All	713/830 (86%)	639 (90%)	74 (10%)	7	22

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	318	HIS
1	A	327	VAL
1	A	364	LYS
1	A	428	LYS

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Mol	Chain	Res	Type
1	A	438	GLU
1	A	464	LEU
1	A	466	GLU
1	A	475	LYS
1	A	493	VAL
1	A	554	ASP
1	A	594	LYS
1	A	615	THR
1	A	620	GLU
1	A	645	SER
1	A	657	VAL
1	A	666	LYS
1	B	238	MET
1	B	239	ILE
1	B	243	LEU
1	B	247	LEU
1	B	259	HIS
1	B	268	VAL
1	B	269	VAL
1	B	272	GLN
1	B	299	LEU
1	B	302	ASP
1	B	303	GLN
1	B	304	THR
1	B	308	GLN
1	B	317	LEU
1	B	322	VAL
1	B	327	VAL
1	B	332	LEU
1	B	334	LYS
1	B	344	SER
1	B	348	ARG
1	B	350	LYS
1	B	363	VAL
1	B	375	THR
1	B	376	LYS
1	B	381	GLU
1	B	417	ASN
1	B	419	ASP
1	B	428	LYS
1	B	430	GLU
1	B	431	ARG

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Mol	Chain	Res	Type
1	B	481	ARG
1	B	483	LYS
1	B	484	ILE
1	B	485	GLU
1	B	488	GLN
1	B	491	ILE
1	B	494	ARG
1	B	501	THR
1	B	562	GLN
1	B	568	GLU
1	B	573	LEU
1	B	580	LEU
1	B	586	ASN
1	B	603	THR
1	B	615	THR
1	B	619	ASP
1	B	621	LEU
1	B	623	ARG
1	B	632	LEU
1	B	642	LYS
1	B	643	VAL
1	B	647	LEU
1	B	652	ILE
1	B	660	GLU
1	B	667	GLU
1	B	680	ILE
1	B	690	LYS
1	B	698	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	316	GLN
1	A	346	HIS
1	A	488	GLN
1	A	507	GLN
1	A	552	HIS
1	A	577	ASN
1	A	630	HIS
1	A	646	HIS
1	B	235	ASN
1	B	318	HIS

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Mol	Chain	Res	Type
1	B	391	ASN
1	B	479	GLN
1	B	488	GLN
1	B	507	GLN
1	B	601	ASN
1	B	694	GLN
1	B	696	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	31/32 (96%)	3 (9%)	0
3	D	29/32 (90%)	3 (10%)	0
All	All	60/64 (93%)	6 (10%)	0

All (6) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	10	G
2	C	13	8AZ
2	C	29	U
3	D	7	U
3	D	10	C
3	D	19	C

There are no RNA pucker outliers to report.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

Of 2 non-standard protein/DNA/RNA residues modelled in this entry, 1 could not be matched to an existing wwPDB Chemical Component Dictionary definition at this stage - leaving 1 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
2	8AZ	C	13	2,5	18,24,25	0.79	0	19,35,38	1.35	2 (10%)

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	8AZ	C	13	2,5	-	5/7/35/36	0/3/3/3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	13	8AZ	N9-N8-N7	4.00	109.84	107.27
2	C	13	8AZ	C5-C4-N9	2.19	106.00	104.60

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	13	8AZ	O4'-C4'-C5'-O5'
2	C	13	8AZ	C3'-C4'-C5'-O5'
2	C	13	8AZ	O4'-C1'-N9-N8
2	C	13	8AZ	O4'-C1'-N9-C4
2	C	13	8AZ	C2'-C1'-N9-N8

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
4	IHP	B	801	-	36,36,36	0.81	1 (2%)	60,60,60	0.82	1 (1%)
4	IHP	A	801	-	36,36,36	0.79	0	60,60,60	0.89	1 (1%)

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
4	IHP	B	801	-	-	2/30/54/54	0/1/1/1
4	IHP	A	801	-	-	0/30/54/54	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	801	IHP	P4-O14	2.10	1.63	1.59

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	801	IHP	O12-P2-O22	-2.43	100.66	109.33
4	B	801	IHP	P2-O12-C2	2.02	128.84	123.43

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	801	IHP	C1-O11-P1-O21
4	B	801	IHP	C5-O15-P5-O25

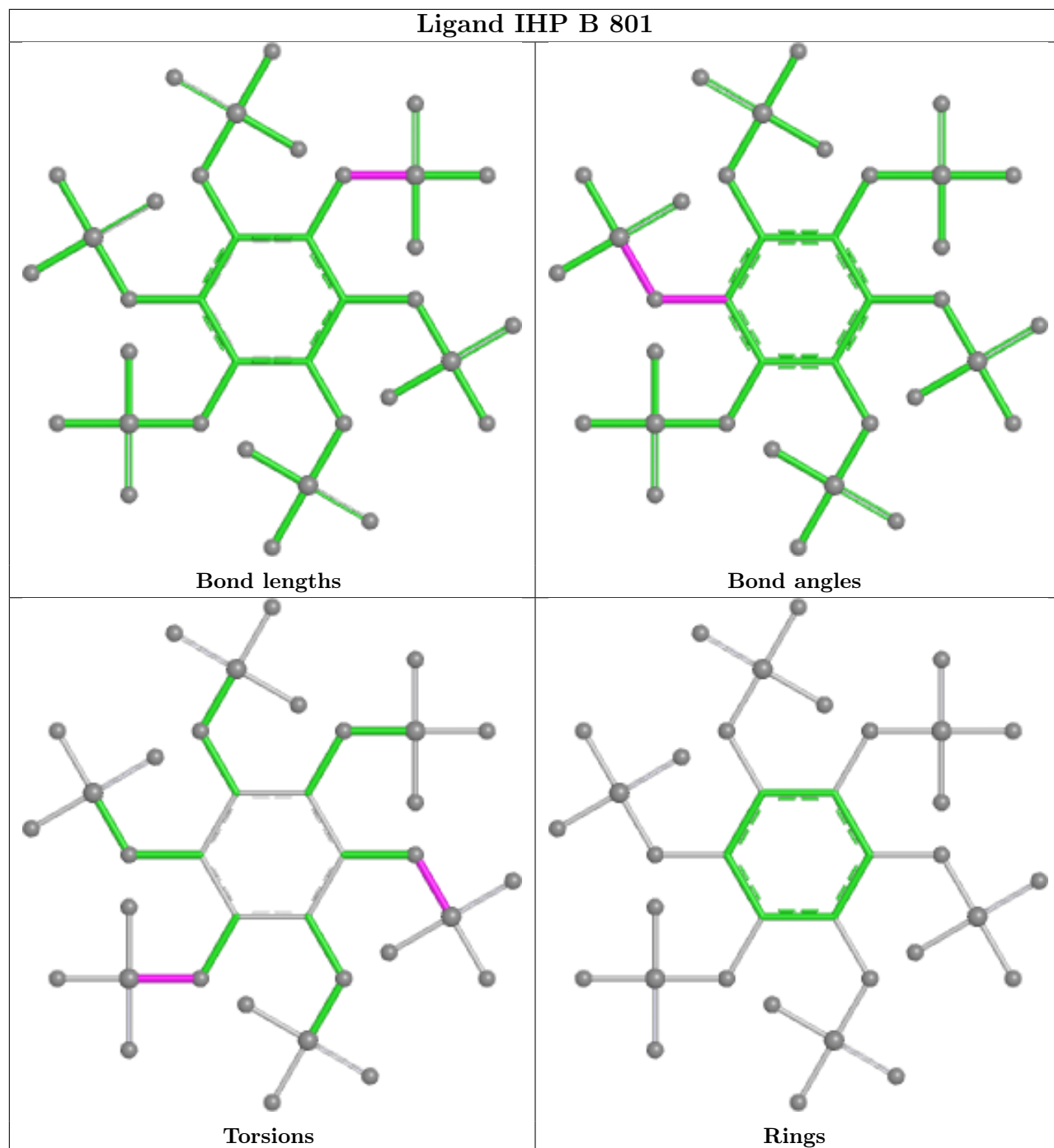
There are no ring outliers.

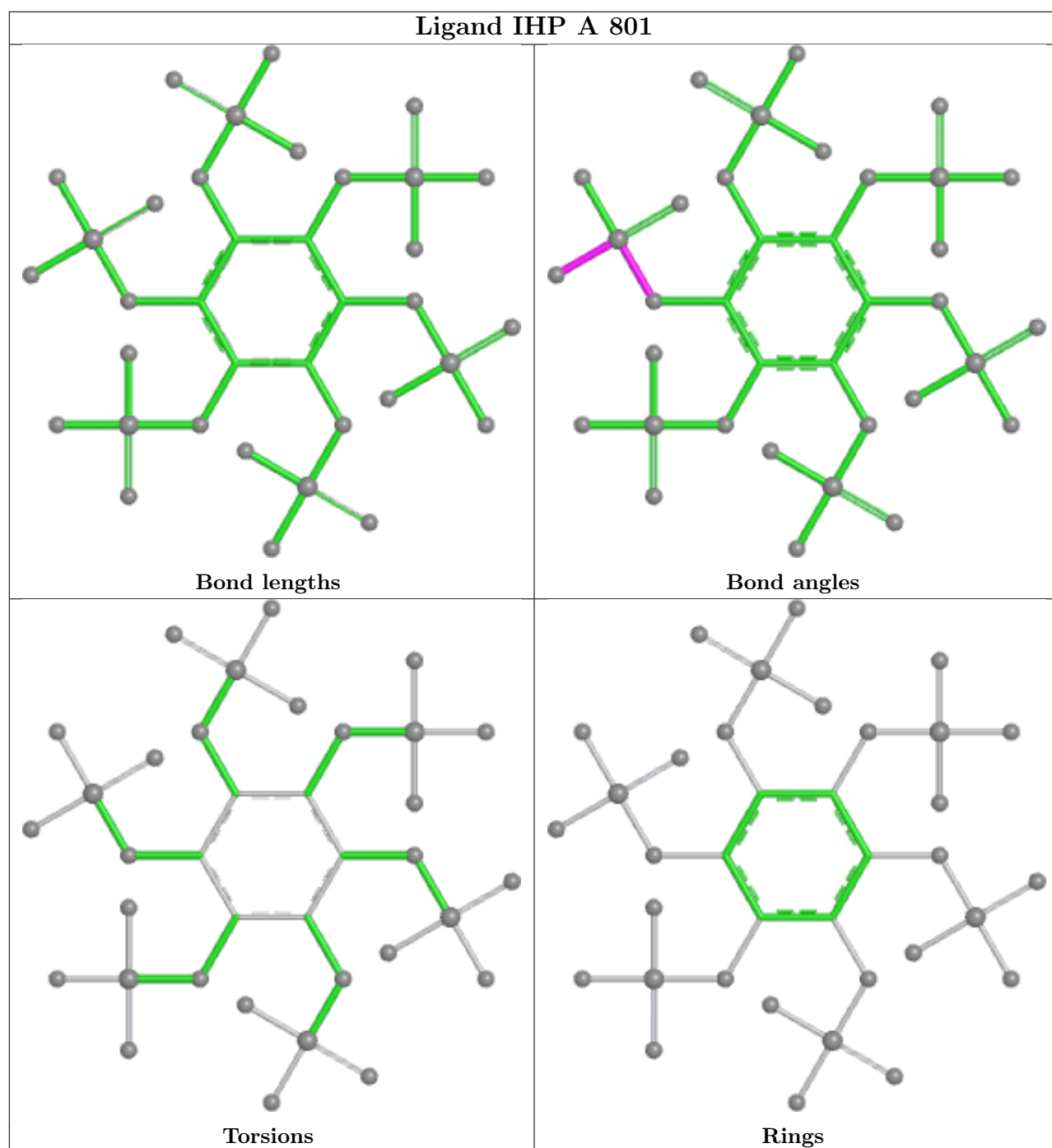
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	801	IHP	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 IHP B 801





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	384/488 (78%)	-0.07	4 (1%) 79 72	73, 88, 135, 204	0
1	B	454/488 (93%)	-0.05	1 (0%) 91 88	80, 107, 176, 213	0
2	C	31/32 (96%)	-0.56	0 100 100	81, 114, 182, 188	0
3	D	31/32 (96%)	-0.42	0 100 100	80, 130, 180, 193	0
All	All	900/1040 (86%)	-0.09	5 (0%) 85 80	73, 99, 174, 213	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	593	GLY	2.6
1	A	366	ALA	2.6
1	A	497	ALA	2.2
1	A	664	ALA	2.2
1	B	499	ILE	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
2	8AZ	C	13	22/23	0.96	0.07	72,76,83,83	0
3	4DU	D	21	21/22	0.97	0.07	79,82,90,95	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

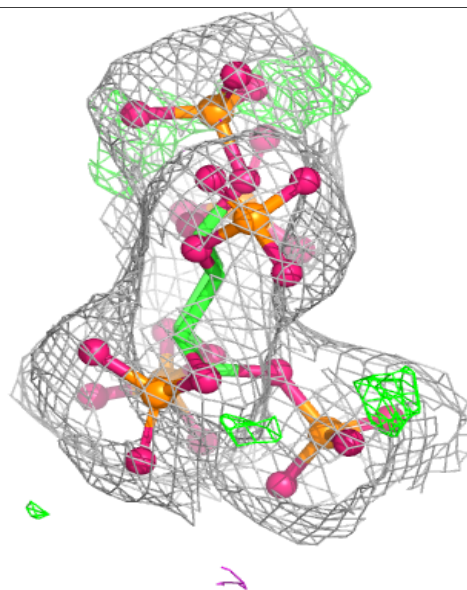
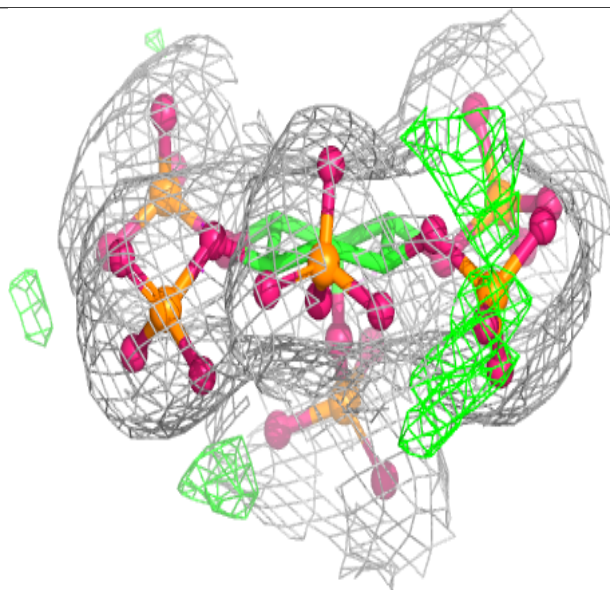
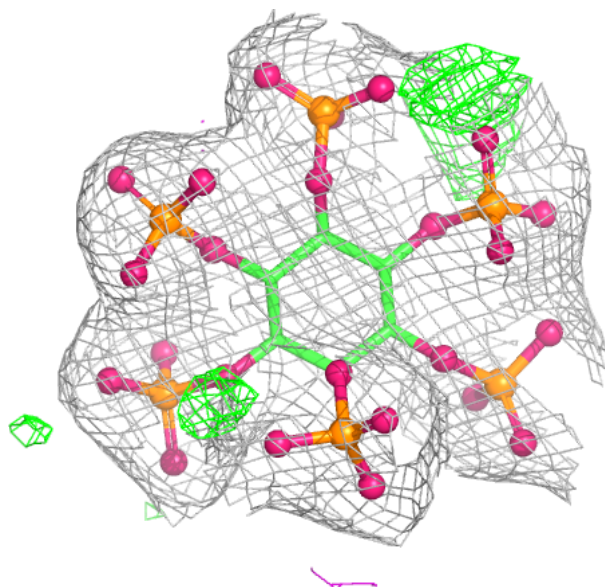
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	IHP	B	801	36/36	0.95	0.07	78,90,98,104	0
4	IHP	A	801	36/36	0.96	0.06	67,75,84,86	0
5	ZN	A	802	1/1	0.99	0.04	74,74,74,74	0
5	ZN	B	802	1/1	1.00	0.03	90,90,90,90	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.

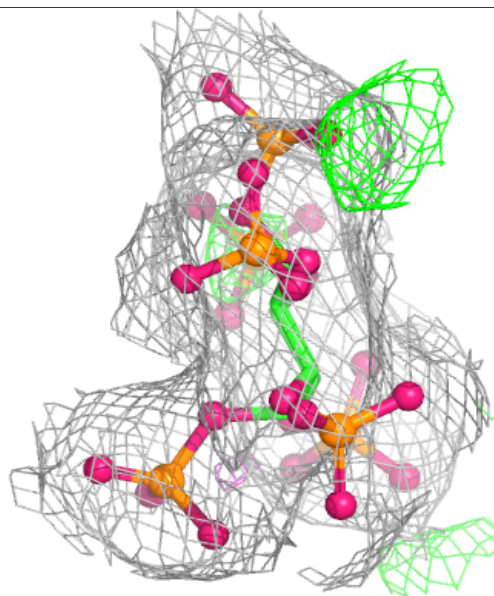
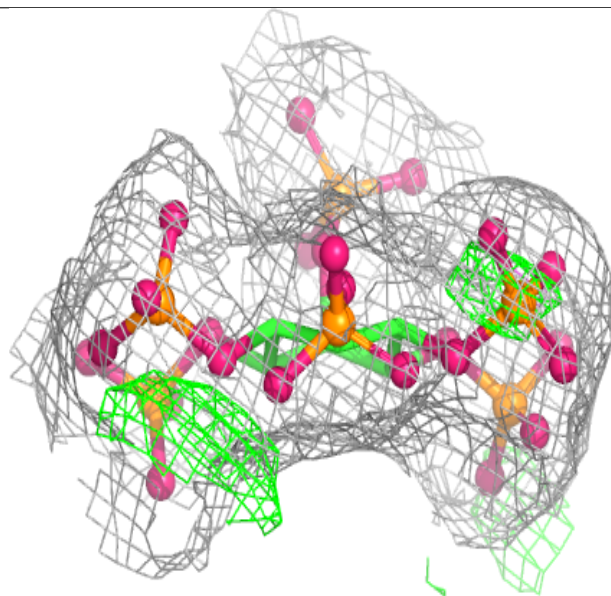
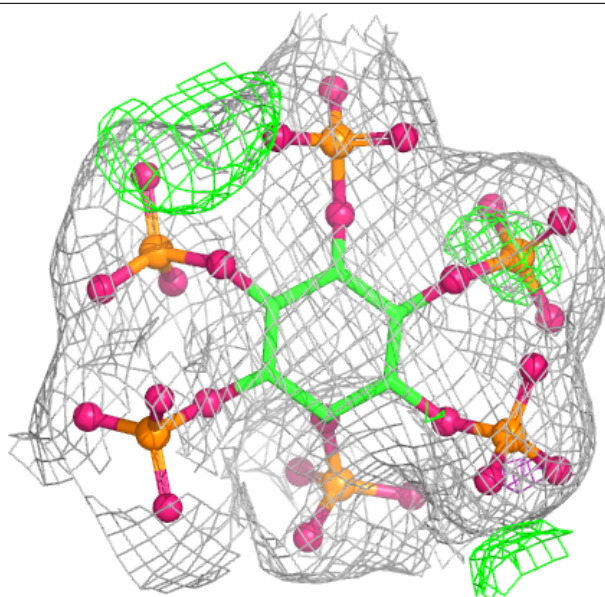
Electron density around IHP B 801:

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



Electron density around IHP A 801:

$2mF_o-DF_c$ (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.