



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:45 PM GMT

PDB ID : 4W5T
Title : The Crystal Structure of Human Argonaute2 Bound to a Guide and Target
RNA Containing Seed Pairing from 2-7
Authors : Schirle, N.T.; MacRae, I.J.
Deposited on : 2014-08-18
Resolution : 2.50 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

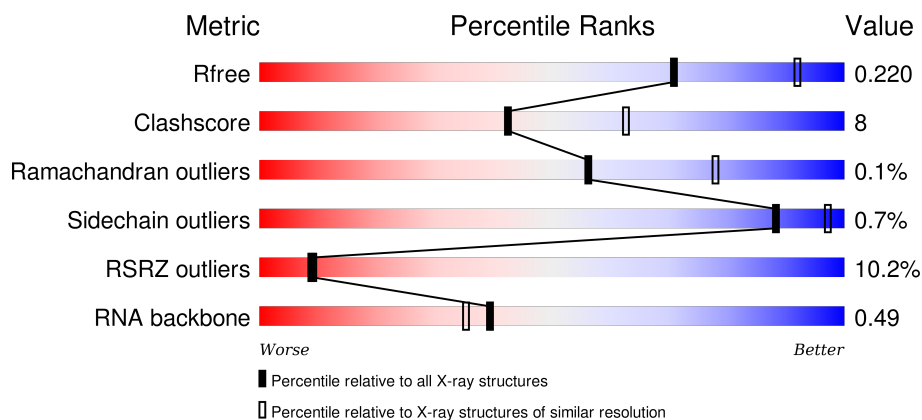
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)
RNA backbone	2183	1172 (3.00-2.00)

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. 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	859	<div> <div>9%</div> <div>78%</div> <div>15%</div> <div>6%</div> </div>
2	B	21	<div> <div>19%</div> <div>29%</div> <div>33%</div> <div>14%</div> <div>24%</div> </div>
3	D	11	<div> <div>9%</div> <div>36%</div> <div>36%</div> <div>18%</div> <div>9%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	IPH	A	902	-	-	-	X
5	IPH	A	904	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7184 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 argonaute-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	807	Total	C	N	O	S	0	0	0
			6471	4119	1165	1147	40			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	387	ASP	SER	engineered mutation	UNP Q9UKV8

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	16	Total	C	N	O	P	0	0	0
			332	149	52	115	16			

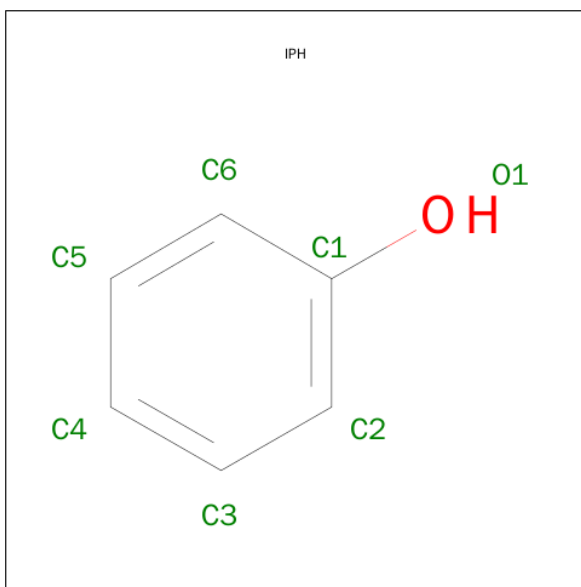
- Molecule 3 is a RNA chain called RNA (5'-R(*AP*CP*AP*UP*GP*UP*GP*AP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	10	Total	C	N	O	P	0	0	0
			195	87	37	62	9			

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		

- Molecule 5 is PHENOL (three-letter code: IPH) (formula: C₆H₆O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	6	1		
5	A	1	Total	C	O	0	0
			7	6	1		
5	A	1	Total	C	O	0	0
			7	6	1		
5	A	1	Total	C	O	0	0
			7	6	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	141	Total	O	0	0
			141	141		
6	B	10	Total	O	0	0
			10	10		
6	D	5	Total	O	0	0
			5	5		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.64Å 116.47Å 69.82Å 90.00° 92.33° 90.00°	Depositor
Resolution (Å)	35.29 – 2.50 35.29 – 2.50	Depositor EDS
% Data completeness (in resolution range)	95.6 (35.29-2.50) 95.6 (35.29-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.173 , 0.215 0.187 , 0.220	Depositor DCC
R_{free} test set	1431 reflections (5.11%)	DCC
Wilson B-factor (Å ²)	39.6	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.9	EDS
Estimated twinning fraction	0.035 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 29414 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7184	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: IPH, MG

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	0.39	1/6624 (0.0%)	0.55	5/8964 (0.1%)
2	B	1.72	6/367 (1.6%)	1.00	3/563 (0.5%)
3	D	1.13	2/218 (0.9%)	0.85	1/339 (0.3%)
All	All	0.58	9/7209 (0.1%)	0.60	9/9866 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	U	OP3-P	-18.61	1.38	1.61
2	B	4	A	O3'-P	-9.06	1.50	1.61
2	B	3	C	O3'-P	-7.93	1.51	1.61
2	B	5	C	O3'-P	-7.85	1.51	1.61
2	B	2	U	O3'-P	-6.94	1.52	1.61
3	D	3	A	O3'-P	-6.04	1.53	1.61
2	B	6	A	O3'-P	-6.00	1.53	1.61
3	D	5	G	O3'-P	-5.35	1.54	1.61
1	A	609	PRO	N-CD	5.05	1.54	1.47

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	109	GLY	N-CA-C	-11.99	83.11	113.10
1	A	110	ARG	N-CA-C	7.65	131.65	111.00
2	B	1	U	P-O3'-C3'	6.65	127.68	119.70
1	A	248	LYS	C-N-CD	6.20	141.42	128.40
1	A	488	GLN	C-N-CD	5.78	140.54	128.40
1	A	606	GLY	N-CA-C	5.44	126.69	113.10
3	D	5	G	O5'-P-OP1	-5.40	100.84	105.70
2	B	4	A	O5'-P-OP2	-5.17	101.05	105.70
2	B	11	C	O3'-P-O5'	5.15	113.79	104.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6471	0	6534	108	0
2	B	332	0	172	4	0
3	D	195	0	98	5	0
4	A	1	0	0	0	0
4	D	1	0	0	0	0
5	A	28	0	24	2	0
6	A	141	0	0	5	0
6	B	10	0	0	1	0
6	D	5	0	0	0	0
All	All	7184	0	6828	114	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 8.

All (114) 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:108:ILE:HD13	1:A:113:VAL:HG21	1.45	0.98
1:A:66:CYS:SG	1:A:70:VAL:HB	2.09	0.92
1:A:66:CYS:SG	1:A:70:VAL:HG11	2.11	0.91
1:A:66:CYS:SG	1:A:70:VAL:CB	2.59	0.90
1:A:836:GLY:O	1:A:840:GLN:HB2	1.76	0.85
1:A:269:ILE:HD11	1:A:339:LEU:HD22	1.60	0.84
1:A:66:CYS:SG	1:A:70:VAL:CG1	2.66	0.83
1:A:55:TYR:CZ	1:A:105:PRO:HG3	2.15	0.81
1:A:55:TYR:CE2	1:A:105:PRO:HG3	2.17	0.80
1:A:837:ARG:HH11	1:A:837:ARG:CB	1.96	0.78
1:A:111:ASP:O	1:A:113:VAL:HG23	1.88	0.74
1:A:108:ILE:HD13	1:A:113:VAL:CG2	2.17	0.74
1:A:837:ARG:HB3	1:A:837:ARG:HH11	1.53	0.72
1:A:207:ARG:NH2	1:A:673:GLU:OE1	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:811:PHE:CE1	3:D:1:A:H4'	2.24	0.72
1:A:243:ILE:HG22	1:A:244:GLU:H	1.54	0.71
1:A:293:THR:HG22	1:A:307:THR:HG22	1.73	0.70
1:A:66:CYS:SG	1:A:67:PRO:HD2	2.32	0.70
5:A:905:IPH:H6	6:A:1137:HOH:O	1.92	0.69
1:A:143:LEU:HB2	1:A:158:THR:HG21	1.74	0.68
1:A:836:GLY:O	1:A:840:GLN:CB	2.41	0.68
1:A:88:GLY:O	1:A:89:ASP:HB2	1.94	0.68
1:A:428:ALA:HB2	1:A:440:LYS:HE2	1.77	0.67
1:A:56:HIS:NE2	1:A:99:ASN:OD1	2.30	0.64
1:A:108:ILE:CD1	1:A:113:VAL:HG21	2.24	0.64
1:A:320:LEU:HD22	1:A:342:GLU:OE2	1.98	0.62
1:A:692:ILE:HD11	5:A:902:IPH:H4	1.82	0.62
1:A:243:ILE:HG22	1:A:244:GLU:N	2.14	0.62
1:A:58:GLU:OE2	1:A:98:LYS:HE2	1.99	0.61
1:A:269:ILE:HG22	1:A:270:THR:N	2.15	0.61
1:A:264:GLY:O	1:A:280:ARG:NH1	2.31	0.61
1:A:147:LEU:HD21	1:A:213:MET:HE2	1.83	0.60
1:A:449:LYS:HG2	1:A:486:GLN:OE1	2.01	0.59
1:A:269:ILE:HD11	1:A:339:LEU:CD2	2.31	0.59
1:A:211:TRP:CD1	1:A:212:LYS:HG3	2.38	0.58
1:A:109:GLY:O	1:A:111:ASP:OD2	2.22	0.58
1:A:605:ASP:CB	1:A:608:LYS:HE3	2.34	0.57
1:A:605:ASP:HB3	1:A:608:LYS:HE3	1.86	0.57
1:A:820:LYS:O	1:A:821:GLU:HB2	2.02	0.57
1:A:279:TYR:HB3	1:A:330:VAL:HB	1.87	0.57
1:A:635:ARG:NH1	2:B:10:C:O2'	2.37	0.57
1:A:66:CYS:SG	1:A:70:VAL:HG21	2.45	0.57
1:A:92:PRO:HB3	1:A:100:LEU:HD21	1.85	0.57
1:A:77:HIS:CE1	1:A:119:LEU:HD12	2.40	0.56
1:A:87:PHE:O	1:A:90:ARG:HG2	2.05	0.56
3:D:1:A:H2'	3:D:2:C:C6	2.40	0.56
1:A:269:ILE:CG2	1:A:270:THR:N	2.69	0.55
1:A:756:ILE:HD11	1:A:795:ARG:HH22	1.71	0.55
1:A:374:ASP:O	1:A:378:GLU:HG2	2.05	0.55
1:A:291:HIS:O	1:A:293:THR:HG23	2.06	0.55
1:A:66:CYS:SG	1:A:70:VAL:CG2	2.95	0.54
1:A:269:ILE:CD1	1:A:339:LEU:HD22	2.36	0.54
1:A:243:ILE:CG2	1:A:244:GLU:H	2.19	0.54
1:A:680:LEU:HD22	1:A:768:LEU:HB3	1.88	0.54
1:A:86:ILE:HG22	1:A:87:PHE:N	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:PRO:HD2	1:A:128:PHE:CE1	2.44	0.52
1:A:525:LYS:HE2	3:D:3:A:OP1	2.09	0.51
2:B:10:C:O2	6:B:104:HOH:O	2.17	0.51
1:A:79:VAL:HA	1:A:87:PHE:CE2	2.45	0.51
1:A:77:HIS:CE1	1:A:119:LEU:CD1	2.94	0.51
1:A:61:ILE:HG12	1:A:130:VAL:HG22	1.93	0.50
1:A:521:ILE:HG12	1:A:552:VAL:HG21	1.93	0.50
1:A:332:GLN:OE1	1:A:335:LYS:HE3	2.12	0.50
1:A:636:GLN:NE2	1:A:638:ILE:O	2.40	0.50
1:A:315:ARG:HG3	1:A:316:HIS:CD2	2.47	0.49
1:A:237:VAL:CG1	1:A:261:GLU:HG3	2.42	0.49
1:A:389:ASN:O	1:A:395:ARG:HD3	2.13	0.49
2:B:9:G:H2'	2:B:10:C:O4'	2.13	0.49
1:A:117:VAL:O	1:A:127:ILE:N	2.46	0.48
1:A:237:VAL:HG11	1:A:261:GLU:HG3	1.94	0.48
1:A:167:ARG:HB2	1:A:181:PHE:HZ	1.79	0.48
1:A:55:TYR:CZ	1:A:105:PRO:CG	2.93	0.48
1:A:108:ILE:HG21	1:A:113:VAL:HG23	1.96	0.48
1:A:77:HIS:NE2	1:A:119:LEU:HD12	2.29	0.47
1:A:569:VAL:HG21	1:A:791:VAL:HB	1.96	0.47
1:A:108:ILE:HG21	1:A:113:VAL:CG2	2.43	0.47
1:A:196:ARG:NH2	1:A:261:GLU:OE1	2.46	0.47
1:A:55:TYR:CE1	1:A:105:PRO:HD3	2.49	0.47
1:A:708:GLN:HE21	1:A:710:ARG:H	1.63	0.47
1:A:55:TYR:HD2	1:A:134:TRP:CH2	2.33	0.46
1:A:244:GLU:OE1	1:A:247:GLN:HB2	2.15	0.46
1:A:235:CYS:SG	1:A:243:ILE:CD1	3.04	0.46
1:A:709:LYS:HG3	6:A:1125:HOH:O	2.16	0.46
1:A:87:PHE:O	1:A:90:ARG:CG	2.64	0.45
2:B:3:C:H2'	2:B:4:A:C8	2.51	0.45
1:A:451:TRP:CZ2	1:A:489:PRO:HD3	2.51	0.45
1:A:837:ARG:HH11	1:A:837:ARG:CG	2.29	0.45
1:A:811:PHE:CD1	3:D:1:A:H4'	2.52	0.45
1:A:45:PHE:CZ	1:A:383:MET:HG3	2.52	0.45
1:A:260:LYS:HD3	1:A:260:LYS:HA	1.87	0.45
1:A:88:GLY:O	1:A:89:ASP:CB	2.64	0.45
1:A:726:LYS:HG3	6:A:1102:HOH:O	2.17	0.45
1:A:112:LYS:HA	1:A:132:ILE:O	2.17	0.44
3:D:6:U:H2'	3:D:7:G:O4'	2.18	0.44
1:A:72:ARG:NH2	6:A:1072:HOH:O	2.30	0.44
1:A:361:THR:O	1:A:365:ILE:HG12	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:ASP:HB3	1:A:390:THR:OG1	2.18	0.43
1:A:742:HIS:CG	1:A:743:PRO:HD2	2.53	0.43
1:A:112:LYS:HG3	1:A:131:SER:HB2	2.01	0.43
1:A:210:LEU:HD23	1:A:743:PRO:HA	2.00	0.43
1:A:636:GLN:NE2	1:A:640:GLN:OE1	2.40	0.42
1:A:263:LYS:HE3	1:A:263:LYS:HB2	1.90	0.42
1:A:81:HIS:HD2	1:A:82:PHE:CE1	2.37	0.42
1:A:491:PHE:CD1	1:A:508:LEU:HD21	2.55	0.42
1:A:316:HIS:O	1:A:318:LEU:HG	2.20	0.41
1:A:487:GLY:N	6:A:1116:HOH:O	2.36	0.41
1:A:63:PRO:HD2	1:A:128:PHE:HE1	1.85	0.41
1:A:658:ARG:HA	1:A:658:ARG:HD3	1.65	0.41
1:A:218:ASP:OD1	1:A:375:ARG:NH1	2.53	0.41
1:A:235:CYS:SG	1:A:243:ILE:HD13	2.61	0.40
1:A:77:HIS:HE1	1:A:119:LEU:CD1	2.33	0.40
1:A:330:VAL:HG11	1:A:339:LEU:HD12	2.03	0.40
1:A:64:GLU:O	1:A:64:GLU:HG3	2.22	0.40
1:A:309:ALA:HB2	1:A:320:LEU:HD12	2.03	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	797/859 (93%)	769 (96%)	27 (3%)	1 (0%)	56 78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	110	ARG

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	714/752 (95%)	709 (99%)	5 (1%)	88 97

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	172	MET
1	A	186	GLU
1	A	351	ARG
1	A	804	TYR
1	A	837	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	HIS
1	A	708	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	14/21 (66%)	2 (14%)	0
3	D	8/11 (72%)	1 (12%)	0
All	All	22/32 (68%)	3 (13%)	0

All (3) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	9	G
2	B	13	A
3	D	2	C

There are no RNA pucker outliers to report.

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	$\# Z > 2$
5	IPH	A	902	-	7,7,7	0.46	0	8,8,8	0.25	0
5	IPH	A	903	-	7,7,7	0.41	0	8,8,8	0.34	0
5	IPH	A	904	-	7,7,7	0.37	0	8,8,8	0.25	0
5	IPH	A	905	-	7,7,7	0.41	0	8,8,8	0.27	0

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
5	IPH	A	902	-	-	0/0/0/0	0/1/1/1
5	IPH	A	903	-	-	0/0/0/0	0/1/1/1
5	IPH	A	904	-	-	0/0/0/0	0/1/1/1
5	IPH	A	905	-	-	0/0/0/0	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	902	IPH	1	0
5	A	905	IPH	1	0

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	807/859 (93%)	0.41	80 (9%) 9 10	19, 42, 93, 115	0
2	B	16/21 (76%)	0.86	4 (25%) 1 1	27, 87, 113, 117	0
3	D	10/11 (90%)	-0.22	1 (10%) 9 10	44, 50, 82, 89	0
All	All	833/891 (93%)	0.41	85 (10%) 9 9	19, 43, 96, 117	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	246	GLN	7.4
1	A	108	ILE	6.9
1	A	120	PRO	6.5
1	A	119	LEU	6.0
1	A	104	MET	5.9
1	A	117	VAL	5.8
1	A	23	PHE	5.2
1	A	118	THR	5.1
1	A	86	ILE	4.8
1	A	114	GLU	4.3
1	A	486	GLN	4.2
1	A	247	GLN	4.2
1	A	63	PRO	4.2
1	A	64	GLU	4.1
1	A	422	GLY	4.0
1	A	305	GLU	4.0
1	A	89	ASP	4.0
1	A	134	TRP	3.9
1	A	836	GLY	3.8
1	A	81	HIS	3.8
1	A	604	GLY	3.7
1	A	77	HIS	3.7
1	A	423	ARG	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	69	ARG	3.6
1	A	90	ARG	3.6
1	A	65	LYS	3.5
1	A	150	ARG	3.4
1	A	87	PHE	3.3
1	A	605	ASP	3.3
2	B	10	C	3.3
1	A	244	GLU	3.3
1	A	296	LEU	3.2
1	A	55	TYR	3.2
1	A	66	CYS	3.2
1	A	151	LEU	3.2
1	A	141	GLN	3.1
1	A	606	GLY	3.1
2	B	9	G	3.1
1	A	88	GLY	3.1
1	A	270	THR	3.1
1	A	355	LYS	3.1
1	A	80	GLN	3.0
1	A	22	ALA	2.9
1	A	367	ALA	2.9
1	A	133	LYS	2.9
1	A	128	PHE	2.9
1	A	332	GLN	2.9
1	A	211	TRP	2.8
1	A	152	PRO	2.8
1	A	51	LYS	2.8
1	A	313	LYS	2.8
1	A	116	GLU	2.7
1	A	243	ILE	2.7
1	A	424	ASN	2.7
1	A	115	LEU	2.7
1	A	61	ILE	2.6
1	A	129	LYS	2.6
1	A	111	ASP	2.6
1	A	140	LEU	2.6
1	A	607	LYS	2.5
1	A	837	ARG	2.5
1	A	803	ALA	2.5
2	B	14	A	2.5
3	D	1	A	2.4
1	A	67	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	289	ALA	2.4
1	A	113	VAL	2.4
2	B	20	U	2.3
1	A	70	VAL	2.3
1	A	85	GLN	2.3
1	A	97	ARG	2.3
1	A	479	ARG	2.3
1	A	132	ILE	2.2
1	A	50	PRO	2.2
1	A	291	HIS	2.2
1	A	105	PRO	2.2
1	A	24	LYS	2.2
1	A	366	ARG	2.2
1	A	52	ILE	2.1
1	A	802	PRO	2.1
1	A	156	PHE	2.1
1	A	277	ARG	2.1
1	A	142	ALA	2.1
1	A	53	ASP	2.0
1	A	57	TYR	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
5	IPH	A	902	7/7	0.60	0.21	3.90	78,80,82,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	IPH	A	904	7/7	0.96	0.20	2.09	38,39,43,45	0
5	IPH	A	905	7/7	0.94	0.18	0.61	45,46,47,47	0
5	IPH	A	903	7/7	0.94	0.14	0.18	38,40,45,49	0
4	MG	A	901	1/1	0.93	0.07	-4.36	43,43,43,43	0
4	MG	D	101	1/1	0.93	0.13	-	58,58,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.