



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 8, 2021 – 11:32 AM EST

PDB ID : 7MXA
Title : PRMT5:MEP50 complexed with inhibitor PF-06855800
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Deposited on : 2021-05-18
Resolution : 2.71 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 : 2.23.2
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) : 2.23.2

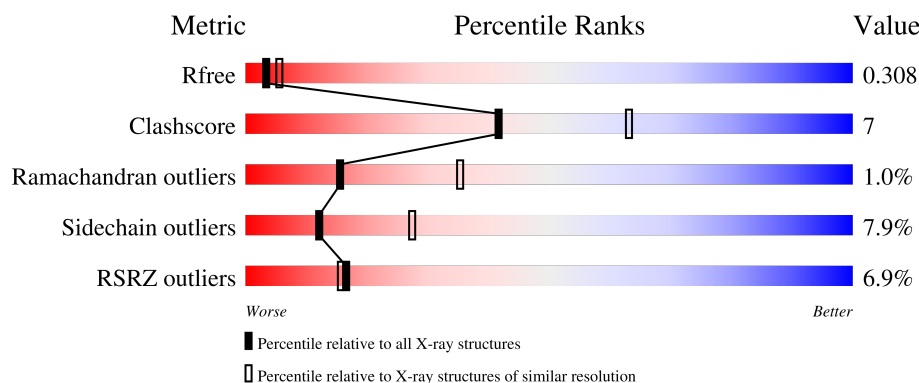
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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	3359 (2.74-2.70)
Clashscore	141614	3686 (2.74-2.70)
Ramachandran outliers	138981	3622 (2.74-2.70)
Sidechain outliers	138945	3623 (2.74-2.70)
RSRZ outliers	127900	3276 (2.74-2.70)

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 $\leq 5\%$. 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	637	<div> <div>4%</div> <div>77%</div> <div>19%</div> <div>..</div> </div>
2	B	366	<div> <div>11%</div> <div>60%</div> <div>22%</div> <div>•</div> <div>16%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7537 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 arginine N-methyltransferase 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	621	Total	C	N	O	S	0	1	0
			5031	3217	865	924	25			

- Molecule 2 is a protein called Methylosome protein 50.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	307	Total	C	N	O	S	0	0	0
			2317	1454	396	453	14			

There are 25 discrepancies between the modelled and reference sequences:

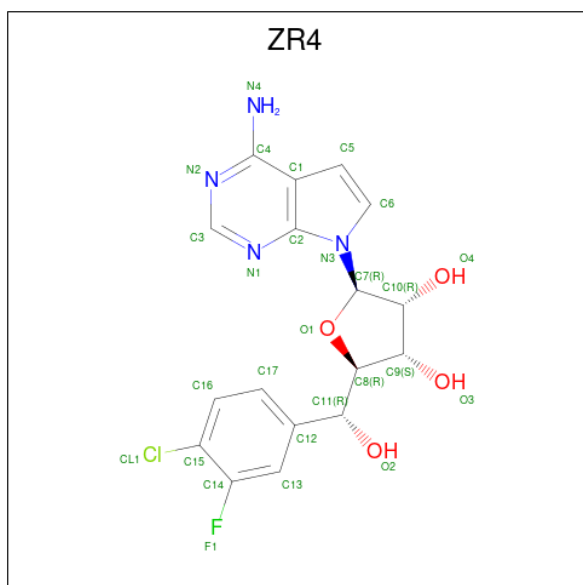
Chain	Residue	Modelled	Actual	Comment	Reference
B	-23	MET	-	expression tag	UNP Q9BQA1
B	-22	ALA	-	expression tag	UNP Q9BQA1
B	-21	SER	-	expression tag	UNP Q9BQA1
B	-20	HIS	-	expression tag	UNP Q9BQA1
B	-19	HIS	-	expression tag	UNP Q9BQA1
B	-18	HIS	-	expression tag	UNP Q9BQA1
B	-17	HIS	-	expression tag	UNP Q9BQA1
B	-16	HIS	-	expression tag	UNP Q9BQA1
B	-15	HIS	-	expression tag	UNP Q9BQA1
B	-14	ASP	-	expression tag	UNP Q9BQA1
B	-13	TYR	-	expression tag	UNP Q9BQA1
B	-12	ASP	-	expression tag	UNP Q9BQA1
B	-11	GLY	-	expression tag	UNP Q9BQA1
B	-10	ALA	-	expression tag	UNP Q9BQA1
B	-9	THR	-	expression tag	UNP Q9BQA1
B	-8	THR	-	expression tag	UNP Q9BQA1
B	-7	GLU	-	expression tag	UNP Q9BQA1
B	-6	ASN	-	expression tag	UNP Q9BQA1
B	-5	LEU	-	expression tag	UNP Q9BQA1
B	-4	TYR	-	expression tag	UNP Q9BQA1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	PHE	-	expression tag	UNP Q9BQA1
B	-2	GLN	-	expression tag	UNP Q9BQA1
B	-1	GLY	-	expression tag	UNP Q9BQA1
B	0	SER	-	expression tag	UNP Q9BQA1
B	1	LEU	-	expression tag	UNP Q9BQA1

- Molecule 3 is 7-[(5R)-5-C-(4-chloro-3-fluorophenyl)-beta-D-ribofuranosyl]-7H-pyrrolo[2,3-d]pyrimidin-4-amine (three-letter code: ZR4) (formula: C₁₇H₁₆ClFN₄O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	F	N	O	0	0
			27	17	1	1	4	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	130	Total	O	0	0
			130	130		
4	B	32	Total	O	0	0
			32	32		

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	102.17Å 138.02Å 178.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	82.12 – 2.71 82.12 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.8 (82.12-2.71) 99.8 (82.12-2.71)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.66 (at 2.73Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.232 , 0.307 0.233 , 0.308	Depositor DCC
R_{free} test set	1668 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	64.6	Xtriage
Anisotropy	1.023	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 62.7	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	0.94	EDS
Total number of atoms	7537	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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: ZR4

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	0.44	0/5171	0.62	1/7034 (0.0%)
2	B	0.40	0/2372	0.63	0/3239
All	All	0.42	0/7543	0.63	1/10273 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	529	MET	C-N-CA	5.21	134.72	121.70

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	5031	0	4925	64	0
2	B	2317	0	2233	39	0
3	A	27	0	0	3	0
4	A	130	0	0	0	0
4	B	32	0	0	0	0
All	All	7537	0	7158	102	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 7.

All (102) 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:567:ILE:HA	1:A:572:HIS:HD2	1.28	0.97
1:A:147:HIS:CD2	1:A:201:ARG:HH12	1.84	0.96
1:A:147:HIS:HD2	1:A:201:ARG:HH12	1.08	0.92
1:A:567:ILE:HA	1:A:572:HIS:CD2	2.05	0.90
1:A:190:HIS:CE1	1:A:547:VAL:HG21	2.12	0.84
1:A:212:LEU:HG	1:A:250:HIS:CE1	2.15	0.82
1:A:21:ASN:HD21	1:A:85:LYS:HZ1	1.35	0.74
1:A:434:SER:HB2	1:A:436:LEU:HG	1.70	0.72
2:B:145:LYS:HA	2:B:169:GLN:HB2	1.74	0.70
1:A:441:ALA:HB2	1:A:555:PHE:HB2	1.73	0.69
1:A:190:HIS:HE1	1:A:547:VAL:HG21	1.57	0.69
1:A:25:GLU:HB3	1:A:28:ASP:HB2	1.75	0.68
1:A:147:HIS:HD2	1:A:201:ARG:NH1	1.88	0.67
1:A:599:CYS:HB3	1:A:619:ALA:HB3	1.78	0.65
1:A:21:ASN:HD21	1:A:85:LYS:NZ	1.94	0.65
1:A:319:LEU:HD22	1:A:323:THR:HG21	1.80	0.63
2:B:262:VAL:HG12	2:B:305:TRP:HE1	1.62	0.63
1:A:330:ASP:OD2	1:A:578:SER:HB2	2.00	0.62
1:A:568:ARG:HG2	1:A:570:GLU:HG2	1.82	0.61
2:B:176:SER:HA	2:B:220:TRP:NE1	2.17	0.60
1:A:436:LEU:O	1:A:446:SER:HB2	2.03	0.59
1:A:21:ASN:ND2	1:A:85:LYS:NZ	2.51	0.59
1:A:54:PHE:H	1:A:102:ASN:HD21	1.51	0.58
2:B:58:LEU:HB2	2:B:77:VAL:HG12	1.87	0.57
1:A:45:VAL:HG23	1:A:46:PHE:HD1	1.69	0.56
1:A:416:VAL:HG11	1:A:423:TRP:CD1	2.41	0.56
1:A:465:PRO:HB3	1:A:559:LEU:HD23	1.87	0.56
2:B:104:GLU:HG2	2:B:120:CYS:HA	1.87	0.56
2:B:175:ALA:HA	2:B:183:PHE:HB3	1.88	0.56
2:B:312:LEU:HD21	2:B:324:HIS:HB3	1.86	0.56
1:A:468:TYR:CE1	1:A:521:PHE:HB2	2.42	0.55
1:A:519:PHE:CE1	1:A:534:ARG:HB3	2.42	0.55
1:A:553:GLY:HA3	1:A:582:ILE:HG22	1.88	0.55
2:B:27:MET:H	2:B:70:GLU:HG2	1.72	0.55
2:B:134:LEU:HD23	2:B:175:ALA:HB1	1.87	0.55
1:A:23:VAL:HG22	1:A:29:THR:HG21	1.89	0.54
1:A:493:ASP:HB3	1:A:496:ALA:HB2	1.88	0.54
1:A:327:PHE:CB	3:A:701:ZR4:CL1	2.94	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:437:LEU:HD21	1:A:468:TYR:CD2	2.44	0.53
2:B:303:ALA:HB1	2:B:313:LEU:HD11	1.91	0.53
2:B:262:VAL:HG12	2:B:305:TRP:NE1	2.24	0.53
2:B:170:VAL:HA	2:B:187:SER:HB2	1.91	0.52
1:A:348:ARG:NH1	1:A:460:ASP:OD1	2.42	0.52
1:A:68:ARG:NH1	2:B:54:TRP:CD2	2.79	0.51
1:A:50:PHE:CE1	1:A:62:ARG:NH1	2.78	0.51
2:B:37:ARG:O	2:B:307:PRO:HG3	2.10	0.51
2:B:282:VAL:HB	2:B:291:PHE:HB3	1.92	0.51
2:B:93:GLY:HA2	2:B:107:GLU:HA	1.94	0.50
2:B:175:ALA:HA	2:B:183:PHE:CB	2.41	0.50
2:B:249:LEU:HD13	2:B:287:LEU:HD12	1.94	0.50
1:A:327:PHE:HB3	3:A:701:ZR4:CL1	2.49	0.49
1:A:327:PHE:HB2	3:A:701:ZR4:CL1	2.50	0.49
1:A:362:MET:SD	1:A:456:PHE:CE1	3.05	0.49
2:B:47:SER:HB3	2:B:83:VAL:H	1.78	0.49
2:B:31:LEU:HB2	2:B:317:GLY:O	2.12	0.48
1:A:301:ALA:HB1	1:A:505:ARG:HG2	1.95	0.48
1:A:437:LEU:HD21	1:A:468:TYR:CE2	2.48	0.48
1:A:349:VAL:HG12	1:A:384:ARG:HH11	1.78	0.48
1:A:433:VAL:HG22	1:A:464:ILE:HB	1.94	0.48
2:B:219:ALA:HB3	2:B:229:VAL:HG23	1.96	0.48
2:B:44:LEU:HD21	2:B:316:VAL:HG11	1.96	0.47
2:B:147:ILE:HG23	2:B:168:ALA:O	2.13	0.47
1:A:41:LEU:HG	1:A:43:MET:HG2	1.96	0.47
2:B:156:GLN:HB3	2:B:158:VAL:HG22	1.97	0.47
1:A:232:PRO:O	1:A:235:ILE:HG12	2.15	0.47
2:B:107:GLU:HB2	2:B:118:LYS:HD3	1.96	0.47
1:A:371:LEU:HD13	1:A:433:VAL:HG12	1.95	0.47
1:A:45:VAL:HG23	1:A:46:PHE:CD1	2.50	0.46
1:A:362:MET:SD	1:A:456:PHE:CD1	3.09	0.46
1:A:150:MET:SD	1:A:201:ARG:HA	2.56	0.45
1:A:214:SER:HB2	1:A:217:VAL:HG23	1.98	0.45
1:A:512:LEU:HD22	1:A:546:THR:HG21	1.98	0.45
1:A:610:LYS:HD2	1:A:634:THR:CB	2.47	0.45
2:B:43:LEU:HB2	2:B:88:TRP:CZ2	2.52	0.45
1:A:21:ASN:ND2	1:A:85:LYS:HZ3	2.15	0.44
2:B:186:CYS:HB3	2:B:192:ILE:HG23	1.99	0.44
2:B:284:ASP:HB3	2:B:290:LEU:HD11	1.99	0.44
1:A:47:HIS:HB3	1:A:50:PHE:HB2	2.00	0.44
1:A:610:LYS:HD2	1:A:634:THR:HB	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:SER:HB3	1:A:89:TRP:CD1	2.53	0.44
2:B:295:ALA:HB1	2:B:323:VAL:HG21	2.00	0.44
1:A:318:ASN:HD21	1:A:395:PRO:HD2	1.82	0.44
2:B:186:CYS:HB2	2:B:215:PRO:HB2	2.00	0.44
1:A:430:ASP:HA	1:A:458:LYS:HD2	1.99	0.44
1:A:465:PRO:HA	1:A:559:LEU:HA	2.00	0.44
1:A:359:GLN:HB2	1:A:386:ILE:HG22	2.00	0.43
2:B:322:VAL:HG12	2:B:322:VAL:O	2.18	0.43
1:A:592:VAL:HG21	1:A:598:ILE:HD11	2.00	0.43
1:A:610:LYS:HD2	1:A:634:THR:OG1	2.18	0.43
1:A:55:ILE:HD12	1:A:102:ASN:HD22	1.84	0.42
2:B:278:CYS:SG	2:B:298:ASP:O	2.77	0.42
2:B:135:SER:OG	2:B:178:HIS:HA	2.19	0.42
2:B:184:LEU:HB3	2:B:218:LEU:HD13	2.02	0.42
1:A:284:LEU:HA	1:A:287:LEU:HD12	2.02	0.42
1:A:73:LEU:HB2	1:A:78:TRP:CE2	2.54	0.42
1:A:558:VAL:HG22	1:A:564:THR:HG22	2.02	0.42
2:B:28:GLU:HB3	2:B:48:SER:HB3	2.01	0.42
2:B:171:THR:HG21	2:B:216:THR:HA	2.03	0.41
1:A:62:ARG:HA	1:A:63:PRO:HD3	1.92	0.41
2:B:189:ASP:OD1	2:B:191:ARG:HB2	2.21	0.41
2:B:280:LEU:HB3	2:B:293:SER:HB3	2.03	0.41
2:B:222:PRO:HG2	2:B:269:PRO:HG2	2.02	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	618/637 (97%)	580 (94%)	35 (6%)	3 (0%)	29	53
2	B	303/366 (83%)	279 (92%)	18 (6%)	6 (2%)	7	17

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	921/1003 (92%)	859 (93%)	53 (6%)	9 (1%)	15	35

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	531	ASP
1	A	578	SER
2	B	135	SER
2	B	147	ILE
2	B	323	VAL
2	B	127	ILE
2	B	266	HIS
2	B	21	PRO
1	A	445	LEU

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	555/562 (99%)	513 (92%)	42 (8%)	13	29
2	B	260/310 (84%)	238 (92%)	22 (8%)	10	23
All	All	815/872 (94%)	751 (92%)	64 (8%)	12	27

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	20	LEU
1	A	25	GLU
1	A	30	LEU
1	A	92	PRO
1	A	102	ASN
1	A	140	ASN
1	A	146	HIS
1	A	149	SER

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Mol	Chain	Res	Type
1	A	178	GLU
1	A	183	GLU
1	A	193	ARG
1	A	234	SER
1	A	248	LYS
1	A	256	ARG
1	A	269	THR
1	A	281	LEU
1	A	286	TYR
1	A	299	LEU
1	A	309	GLN
1	A	310	SER
1	A	322	GLN
1	A	355	ASP
1	A	375	SER
1	A	386	ILE
1	A	392	GLU
1	A	393	LYS
1	A	404	TRP
1	A	405	GLN
1	A	431	ILE
1	A	444	GLU
1	A	467	GLU
1	A	482	ASN
1	A	526	ARG
1	A	529	MET
1	A	530	ILE
1	A	556	GLU
1	A	577	PHE
1	A	578	SER
1	A	582	ILE
1	A	631	ARG
1	A	632	SER
2	B	52	ARG
2	B	53	CYS
2	B	69	ASN
2	B	70	GLU
2	B	78	GLN
2	B	92	ARG
2	B	98	SER
2	B	107	GLU
2	B	114	LEU

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Mol	Chain	Res	Type
2	B	131	VAL
2	B	187	SER
2	B	190	ASN
2	B	192	ILE
2	B	206	ILE
2	B	208	CYS
2	B	249	LEU
2	B	274	LEU
2	B	283	LEU
2	B	292	ARG
2	B	302	ASP
2	B	311	SER
2	B	320	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	102	ASN
1	A	140	ASN
1	A	147	HIS
1	A	270	ASN
1	A	339	GLN
1	A	405	GLN
1	A	572	HIS
2	B	310	HIS

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

1 ligand is 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	ZR4	A	701	-	27,30,30	0.84	1 (3%)	35,45,45	0.85	2 (5%)

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	ZR4	A	701	-	-	0/8/28/28	0/4/4/4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	ZR4	C4-C1	-3.63	1.41	1.45

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	ZR4	C1-C4-N2	-2.70	119.12	121.93
3	A	701	ZR4	O1-C8-C11	2.34	111.51	108.88

There are no chirality outliers.

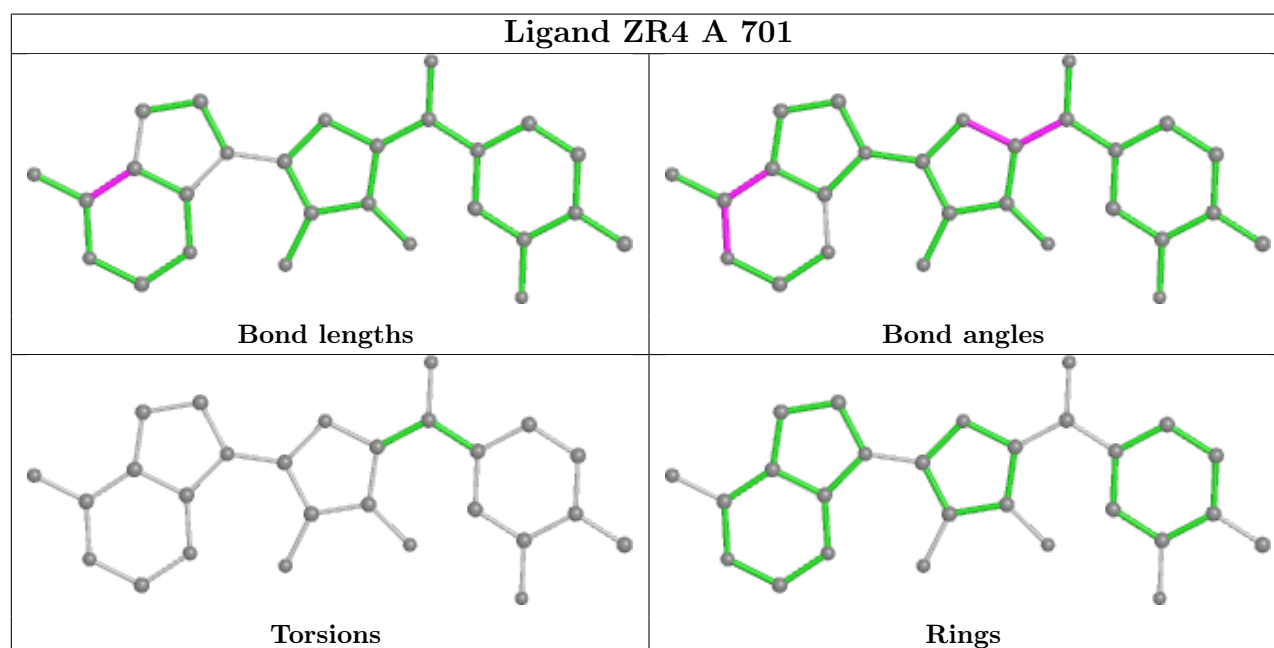
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	701	ZR4	3	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.



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	621/637 (97%)	0.43	23 (3%) 41 41	43, 77, 112, 133	0
2	B	307/366 (83%)	0.74	41 (13%) 3 2	71, 99, 126, 144	0
All	All	928/1003 (92%)	0.53	64 (6%) 16 16	43, 87, 121, 144	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	247	CYS	5.9
1	A	341	ILE	5.8
2	B	23	ALA	5.3
2	B	42	LEU	4.9
2	B	312	LEU	4.5
2	B	40	GLY	4.2
2	B	67	ALA	4.2
1	A	296	ALA	4.0
2	B	59	TRP	3.9
2	B	24	PRO	3.8
2	B	29	ARG	3.8
2	B	106	TRP	3.8
2	B	31	LEU	3.6
2	B	43	LEU	3.5
2	B	27	MET	3.5
2	B	115	ILE	3.5
2	B	174	ALA	3.4
2	B	314	THR	3.3
2	B	41	ALA	3.3
2	B	313	LEU	3.3
1	A	388	LEU	3.2
2	B	61	PHE	3.2
1	A	284	LEU	3.2
2	B	21	PRO	3.2

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Mol	Chain	Res	Type	RSRZ
2	B	218	LEU	3.2
1	A	378	ALA	3.2
2	B	26	CYS	3.2
2	B	316	VAL	3.1
2	B	49	LEU	3.0
2	B	44	LEU	3.0
2	B	303	ALA	3.0
2	B	33	ALA	2.9
2	B	86	LEU	2.9
1	A	361	LEU	2.9
1	A	58	PRO	2.8
1	A	360	VAL	2.8
1	A	297	TYR	2.6
2	B	25	ALA	2.6
1	A	345	LEU	2.6
1	A	358	VAL	2.6
2	B	60	LEU	2.5
2	B	20	PRO	2.5
2	B	94	ILE	2.5
2	B	133	VAL	2.4
2	B	87	THR	2.4
2	B	114	LEU	2.4
1	A	409	TRP	2.4
2	B	318	TRP	2.4
1	A	342	TYR	2.4
1	A	254	ILE	2.4
1	A	431	ILE	2.3
2	B	83	VAL	2.3
1	A	390	ALA	2.3
2	B	22	ASN	2.2
1	A	62	ARG	2.2
2	B	297	ARG	2.2
1	A	82	ILE	2.1
1	A	376	LEU	2.1
1	A	122	PHE	2.1
1	A	33	VAL	2.1
1	A	47	HIS	2.1
2	B	227	VAL	2.1
1	A	208	ILE	2.0
2	B	127	ILE	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 monosaccharides in this entry.

6.4 Ligands [i](#)

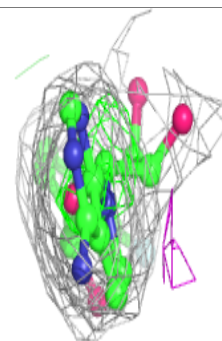
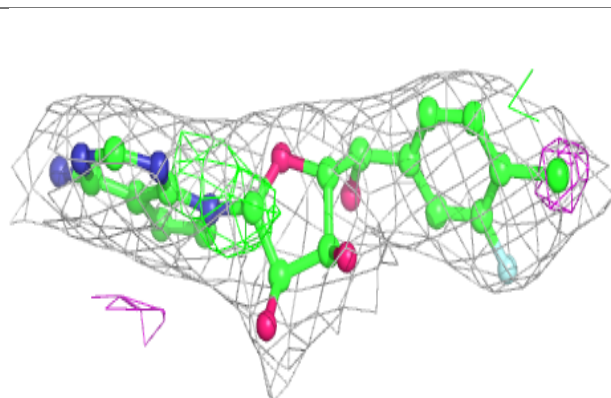
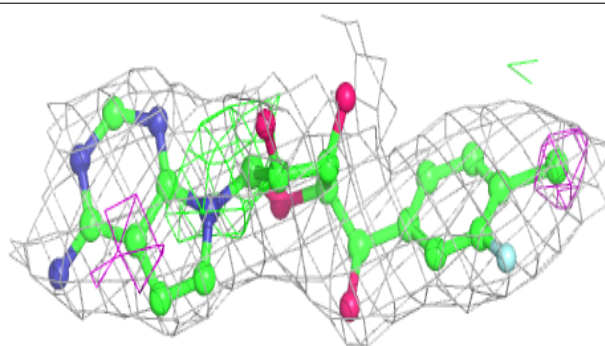
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists 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	ZR4	A	701	27/27	0.92	0.19	53,60,67,71	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 ZR4 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.