



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 05:13 AM EDT

PDB ID : 7MXN / pdb_00007mxn
Title : PRMT5(M420T mutant):MEP50 complexed with inhibitor PF-06939999
Authors : McTigue, M.; Deng, Y.L.; Liu, W.; Brooun, A.
Deposited on : 2021-05-19
Resolution : 2.55 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

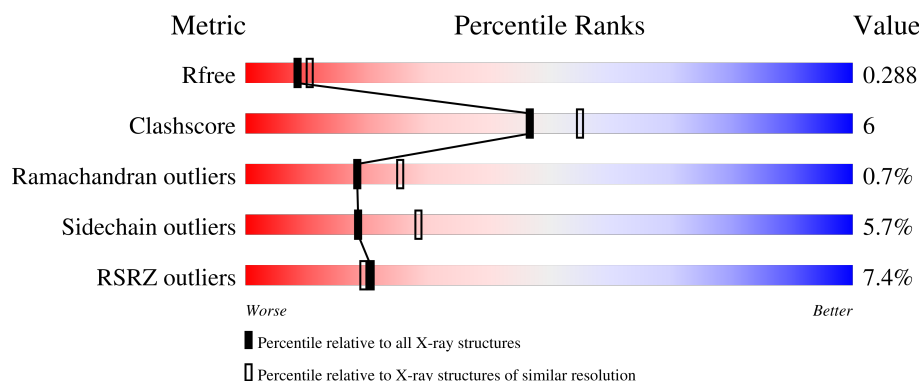
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	
2	B	366	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7580 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
			5030	3216	865	925	24			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	420	THR	MET	engineered mutation	UNP O14744

- Molecule 2 is a protein called Methylosome protein 50.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	301	Total	C	N	O	S	0	0	0
			2281	1430	390	447	14			

There are 25 discrepancies between the modelled and reference sequences:

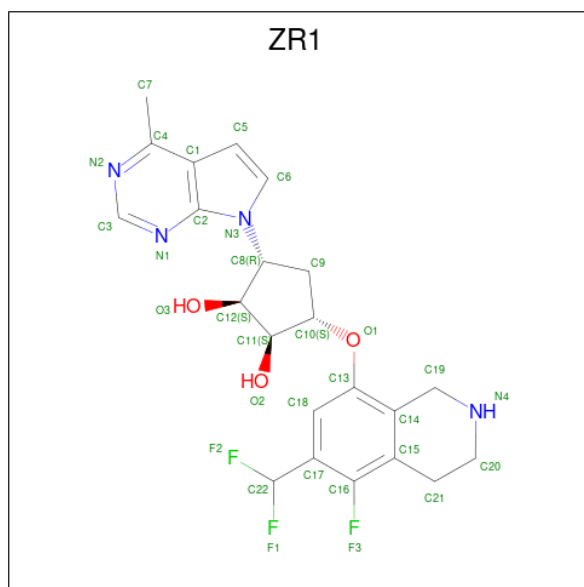
Chain	Residue	Modelled	Actual	Comment	Reference
B	-23	MET	-	expression tag	UNP Q9BQA1
B	-22	ALA	-	expression tag	UNP Q9BQA1
B	-21	SER	-	expression tag	UNP Q9BQA1
B	-20	HIS	-	expression tag	UNP Q9BQA1
B	-19	HIS	-	expression tag	UNP Q9BQA1
B	-18	HIS	-	expression tag	UNP Q9BQA1
B	-17	HIS	-	expression tag	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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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 (1S,2S,3S,5R)-3-{{[6-(difluoromethyl)-5-fluoro-1,2,3,4-tetrahydroisoquinolin-8-yl]oxy}-5-(4-methyl-7H-pyrrolo[2,3-d]pyrimidin-7-yl)cyclopentane-1,2-diol (CCD ID: ZR1) (formula: C₂₂H₂₃F₃N₄O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0
			32	22	3	4	3	

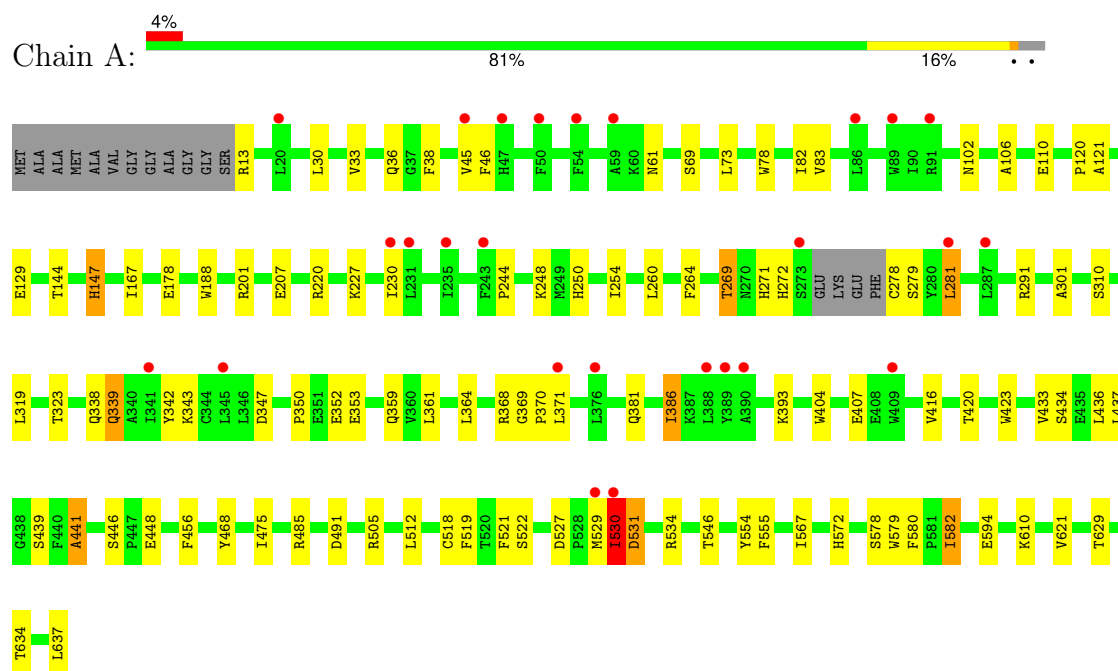
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	193	Total	O	0	0
			193	193		
4	B	44	Total	O	0	0
			44	44		

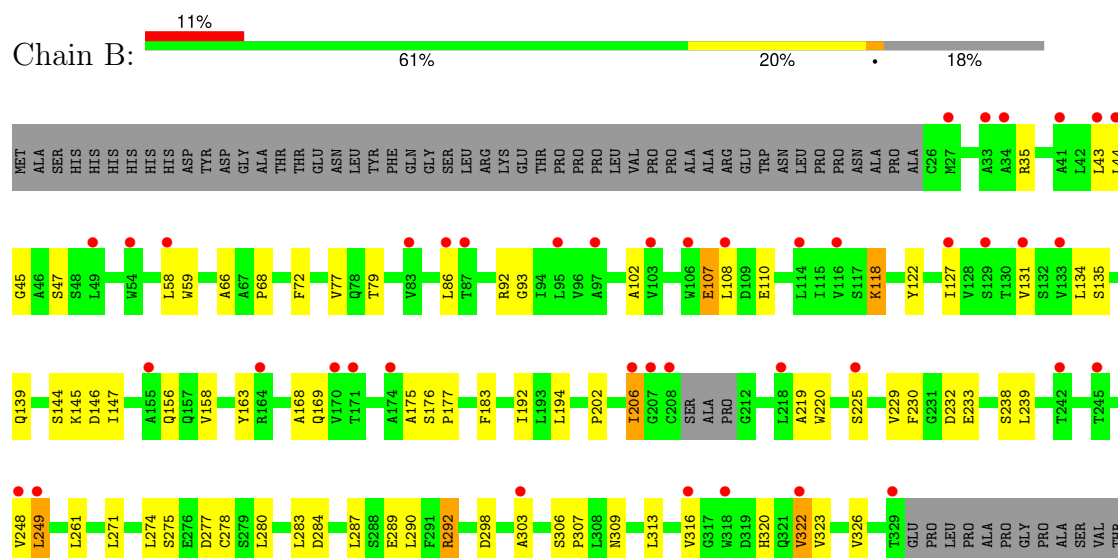
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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: Protein arginine N-methyltransferase 5



• Molecule 2: Methylosome protein 50



GLU

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	100.77Å 137.81Å 178.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	109.19 – 2.55 109.19 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.8 (109.19-2.55) 99.9 (109.19-2.55)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.62 (at 2.55Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.236 , 0.277 0.232 , 0.288	Depositor DCC
R_{free} test set	2010 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	56.4	Xtriage
Anisotropy	1.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 75.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7580	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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: ZR1

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	0.76	4/5170 (0.1%)	1.10	5/7034 (0.1%)
2	B	0.78	2/2333 (0.1%)	1.12	4/3183 (0.1%)
All	All	0.77	6/7503 (0.1%)	1.10	9/10217 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	530	ILE	CA-CB	5.56	1.60	1.54
1	A	518	CYS	CA-C	5.55	1.56	1.52
1	A	368	ARG	CA-C	5.52	1.60	1.52
2	B	322	VAL	CA-C	5.50	1.59	1.52
2	B	168	ALA	CA-C	5.07	1.59	1.52
1	A	582	ILE	CA-C	5.02	1.58	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	529	MET	CA-C-N	7.62	130.75	120.77
1	A	529	MET	C-N-CA	7.62	130.75	120.77
2	B	232	ASP	CA-CB-CG	7.42	120.02	112.60
1	A	456	PHE	CA-CB-CG	6.48	120.28	113.80
2	B	277	ASP	CA-CB-CG	5.84	118.44	112.60
2	B	163	TYR	N-CA-C	5.61	117.91	107.99
2	B	275	SER	N-CA-C	5.23	116.75	108.96
1	A	264	PHE	CA-CB-CG	-5.20	108.61	113.80
1	A	407	GLU	N-CA-C	5.01	117.63	111.82

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	5030	0	4923	46	0
2	B	2281	0	2199	38	0
3	A	32	0	0	0	0
4	A	193	0	0	0	0
4	B	44	0	0	0	0
All	All	7580	0	7122	83	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 6.

All (83) 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:147:HIS:HD2	1:A:201:ARG:HH12	0.96	0.94
1:A:147:HIS:CD2	1:A:201:ARG:HH12	1.88	0.92
2:B:107:GLU:HB2	2:B:118:LYS:HD3	1.62	0.81
2:B:134:LEU:HD12	2:B:139:GLN:HB3	1.71	0.72
1:A:144:THR:O	1:A:147:HIS:HE1	1.74	0.70
1:A:441:ALA:HB2	1:A:555:PHE:HB2	1.72	0.70
1:A:147:HIS:HD2	1:A:201:ARG:NH1	1.82	0.68
2:B:306:SER:HB3	2:B:309:ASN:O	1.93	0.68
2:B:92:ARG:HG3	2:B:108:LEU:HB2	1.76	0.67
2:B:194:LEU:HB2	2:B:206:ILE:HD11	1.77	0.67
2:B:66:ALA:HB1	2:B:72:PHE:HB3	1.76	0.66
2:B:134:LEU:HD23	2:B:175:ALA:HB1	1.80	0.63
2:B:58:LEU:HB2	2:B:77:VAL:HG12	1.81	0.62
1:A:448:GLU:HG2	1:A:637:LEU:HD21	1.83	0.60
1:A:364:LEU:HB3	1:A:420:THR:HG21	1.84	0.59
1:A:319:LEU:HD22	1:A:323:THR:HG21	1.87	0.57
2:B:144:SER:HB3	2:B:146:ASP:OD1	2.05	0.56
1:A:468:TYR:CE1	1:A:521:PHE:HB2	2.41	0.56
1:A:434:SER:HB2	1:A:436:LEU:HG	1.89	0.54
1:A:436:LEU:O	1:A:446:SER:HB2	2.07	0.54
1:A:439:SER:OG	1:A:580:PHE:O	2.25	0.52
1:A:83:VAL:HG22	1:A:121:ALA:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:VAL:HG11	1:A:423:TRP:NE1	2.25	0.52
1:A:371:LEU:HD13	1:A:433:VAL:HG12	1.93	0.51
1:A:512:LEU:HD22	1:A:546:THR:HG21	1.93	0.51
2:B:135:SER:OG	2:B:177:PRO:O	2.29	0.50
1:A:45:VAL:HG23	1:A:46:PHE:HD1	1.77	0.50
2:B:145:LYS:NZ	2:B:169:GLN:NE2	2.60	0.50
1:A:73:LEU:HB2	1:A:78:TRP:CE2	2.47	0.50
1:A:338:GLN:HE21	1:A:381:GLN:HE21	1.59	0.50
2:B:68:PRO:HG3	2:B:320:HIS:O	2.10	0.50
1:A:250:HIS:O	1:A:254:ILE:HG12	2.12	0.49
2:B:261:LEU:HD22	2:B:271:LEU:HD21	1.94	0.49
1:A:33:VAL:HG13	1:A:38:PHE:HB2	1.93	0.49
1:A:278:CYS:HB3	1:A:281:LEU:HB2	1.95	0.49
1:A:301:ALA:HB1	1:A:505:ARG:HG2	1.94	0.49
1:A:567:ILE:HG21	1:A:579:TRP:HB2	1.94	0.48
2:B:249:LEU:HD13	2:B:287:LEU:HD12	1.96	0.48
1:A:82:ILE:O	1:A:120:PRO:HD2	2.14	0.48
1:A:144:THR:O	1:A:147:HIS:CE1	2.60	0.48
1:A:244:PRO:HD2	1:A:279:SER:HB2	1.94	0.48
2:B:284:ASP:HB3	2:B:290:LEU:HD11	1.96	0.47
2:B:134:LEU:HD12	2:B:139:GLN:CB	2.43	0.47
2:B:219:ALA:HB3	2:B:229:VAL:HG23	1.97	0.47
1:A:567:ILE:HA	1:A:572:HIS:CD2	2.49	0.47
2:B:175:ALA:HA	2:B:183:PHE:HB3	1.97	0.47
2:B:192:ILE:HG22	2:B:206:ILE:HD12	1.97	0.47
1:A:269:THR:CG2	1:A:271:HIS:CE1	2.98	0.47
1:A:36:GLN:HE21	1:A:38:PHE:HE2	1.63	0.46
1:A:369:GLY:N	1:A:370:PRO:HD3	2.30	0.46
2:B:156:GLN:HB3	2:B:158:VAL:HG22	1.98	0.46
1:A:530:ILE:HG13	1:A:531:ASP:N	2.30	0.46
2:B:239:LEU:HD23	2:B:249:LEU:HB3	1.97	0.45
2:B:145:LYS:NZ	2:B:169:GLN:HE22	2.15	0.45
2:B:303:ALA:HB1	2:B:313:LEU:HD11	1.99	0.45
1:A:359:GLN:HB2	1:A:386:ILE:HG22	1.98	0.45
1:A:106:ALA:O	1:A:110:GLU:HG3	2.17	0.44
2:B:230:PHE:CE1	2:B:238:SER:HB2	2.53	0.44
1:A:207:GLU:HG3	1:A:230:ILE:HB	2.00	0.44
2:B:134:LEU:HD23	2:B:175:ALA:CB	2.47	0.44
1:A:339:GLN:HG3	1:A:342:TYR:HD2	1.83	0.43
2:B:45:GLY:HA3	2:B:86:LEU:HD21	2.00	0.43
1:A:343:LYS:O	1:A:347:ASP:OD1	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:289:GLU:OE2	2:B:292:ARG:HB2	2.19	0.43
1:A:519:PHE:CE1	1:A:534:ARG:HB3	2.53	0.43
2:B:35:ARG:HB2	2:B:43:LEU:HB3	2.00	0.42
1:A:610:LYS:HD2	1:A:634:THR:OG1	2.20	0.42
2:B:93:GLY:HA2	2:B:107:GLU:HG3	2.00	0.42
2:B:176:SER:HA	2:B:220:TRP:NE1	2.34	0.42
1:A:339:GLN:HA	1:A:342:TYR:HB3	2.01	0.42
2:B:44:LEU:HB2	2:B:59:TRP:HB2	2.02	0.42
1:A:167:ILE:HB	2:B:202:PRO:HD2	2.02	0.41
1:A:437:LEU:HD21	1:A:468:TYR:CD2	2.56	0.41
1:A:129:GLU:HA	1:A:188:TRP:CD1	2.56	0.41
2:B:306:SER:HA	2:B:307:PRO:HD3	1.93	0.41
1:A:244:PRO:HD3	1:A:272:HIS:HB3	2.02	0.41
1:A:350:PRO:HG2	1:A:353:GLU:HG3	2.03	0.41
2:B:102:ALA:HB2	2:B:122:TYR:CD2	2.56	0.40
2:B:110:GLU:H	2:B:110:GLU:CD	2.29	0.40
2:B:219:ALA:HB3	2:B:229:VAL:CG2	2.52	0.40
2:B:278:CYS:SG	2:B:298:ASP:O	2.79	0.40
2:B:316:VAL:HG12	2:B:322:VAL:HG22	2.03	0.40
1:A:554:TYR:HB3	1:A:567:ILE:HG13	2.04	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	618/637 (97%)	593 (96%)	23 (4%)	2 (0%)	36	45
2	B	297/366 (81%)	274 (92%)	19 (6%)	4 (1%)	9	12
All	All	915/1003 (91%)	867 (95%)	42 (5%)	6 (1%)	18	25

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	147	ILE
2	B	323	VAL
1	A	531	ASP
2	B	248	VAL
1	A	441	ALA
2	B	127	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	555/562 (99%)	523 (94%)	32 (6%)	18	27
2	B	257/310 (83%)	243 (95%)	14 (5%)	20	30
All	All	812/872 (93%)	766 (94%)	46 (6%)	18	28

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	ARG
1	A	30	LEU
1	A	61	ASN
1	A	69	SER
1	A	102	ASN
1	A	147	HIS
1	A	178	GLU
1	A	220	ARG
1	A	227	LYS
1	A	248	LYS
1	A	260	LEU
1	A	269	THR
1	A	281	LEU
1	A	291	ARG
1	A	310	SER
1	A	339	GLN
1	A	352	GLU
1	A	361	LEU

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Mol	Chain	Res	Type
1	A	386	ILE
1	A	393	LYS
1	A	404	TRP
1	A	475	ILE
1	A	485	ARG
1	A	491	ASP
1	A	522	SER
1	A	527	ASP
1	A	530	ILE
1	A	578	SER
1	A	582	ILE
1	A	594	GLU
1	A	621	VAL
1	A	629	THR
2	B	47	SER
2	B	79	THR
2	B	107	GLU
2	B	118	LYS
2	B	131	VAL
2	B	206	ILE
2	B	225	SER
2	B	233	GLU
2	B	249	LEU
2	B	274	LEU
2	B	280	LEU
2	B	283	LEU
2	B	292	ARG
2	B	326	VAL

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

Mol	Chain	Res	Type
1	A	36	GLN
1	A	102	ASN
1	A	131	ASN
1	A	147	HIS
1	A	263	GLN
1	A	339	GLN
1	A	381	GLN
1	A	405	GLN
1	A	572	HIS
1	A	588	GLN

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Mol	Chain	Res	Type
2	B	69	ASN
2	B	169	GLN
2	B	254	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	ZR1	A	701	-	35,36,36	1.00	2 (5%)	47,54,54	1.08	3 (6%)

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	ZR1	A	701	-	-	6/12/35/35	0/5/5/5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	ZR1	C19-N4	4.79	1.52	1.46
3	A	701	ZR1	C17-C16	3.26	1.42	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	ZR1	F3-C16-C17	-3.23	116.47	119.67
3	A	701	ZR1	F2-C22-C17	2.97	113.79	110.64
3	A	701	ZR1	C2-N3-C8	2.38	130.46	126.01

There are no chirality outliers.

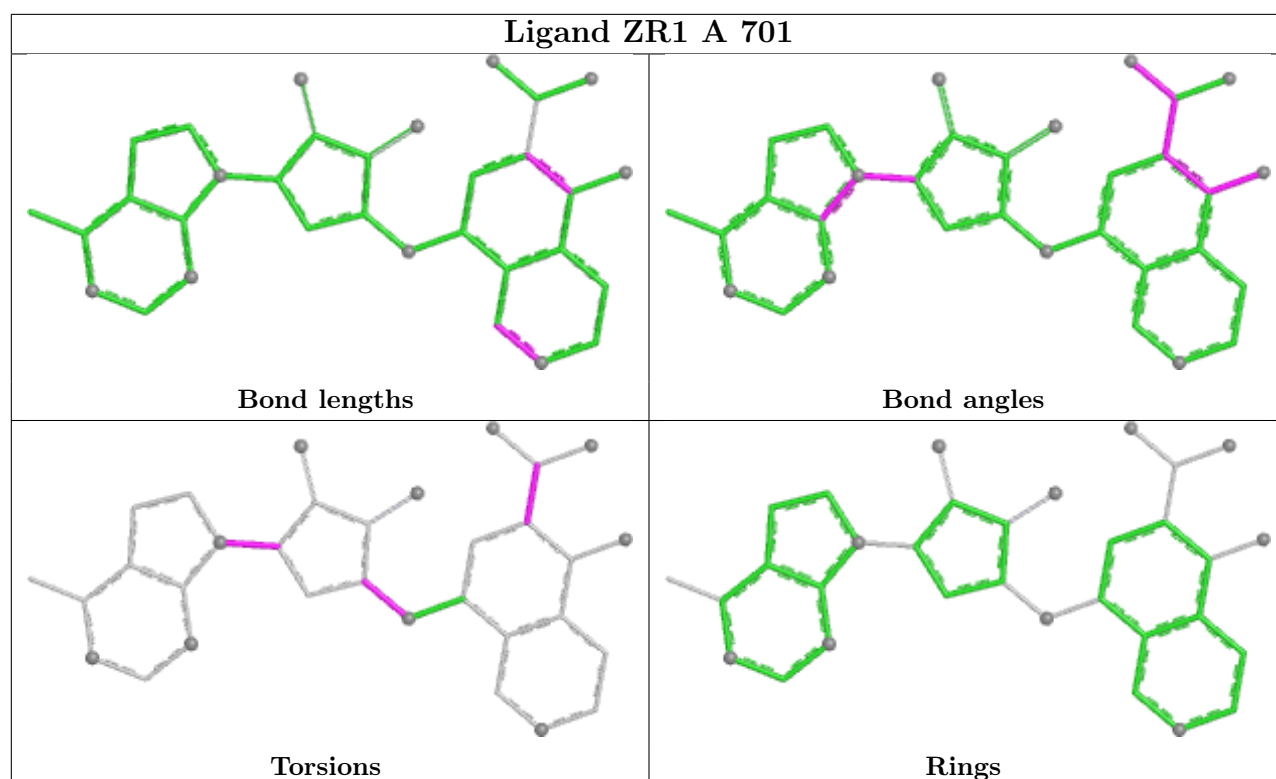
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	701	ZR1	C9-C10-O1-C13
3	A	701	ZR1	C11-C10-O1-C13
3	A	701	ZR1	C18-C17-C22-F1
3	A	701	ZR1	C9-C8-N3-C2
3	A	701	ZR1	C12-C8-N3-C2
3	A	701	ZR1	C12-C8-N3-C6

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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%)	0.49	26 (4%) 40 41	29, 71, 103, 114	1 (0%)
2	B	301/366 (82%)	1.13	42 (13%) 6 5	66, 88, 116, 124	0
All	All	922/1003 (91%)	0.70	68 (7%) 20 19	29, 79, 108, 124	1 (0%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	273	SER	4.6
2	B	248	VAL	4.0
2	B	329	THR	3.5
2	B	245	THR	3.4
2	B	218	LEU	3.1
1	A	390	ALA	3.1
1	A	20	LEU	3.1
1	A	243	PHE	3.1
2	B	249	LEU	3.1
1	A	231	LEU	3.0
2	B	318	TRP	3.0
2	B	41	ALA	2.9
2	B	97	ALA	2.9
2	B	225	SER	2.9
2	B	103	VAL	2.8
1	A	45	VAL	2.8
2	B	164	ARG	2.8
2	B	208	CYS	2.8
2	B	43	LEU	2.8
2	B	44	LEU	2.8
1	A	91	ARG	2.7
1	A	59	ALA	2.6
1	A	388	LEU	2.6
2	B	129	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	281	LEU	2.6
2	B	242	THR	2.5
2	B	207	GLY	2.5
1	A	50	PHE	2.5
2	B	95	LEU	2.5
1	A	345	LEU	2.4
2	B	171	THR	2.4
1	A	530	ILE	2.4
1	A	389	TYR	2.4
2	B	114	LEU	2.4
2	B	27	MET	2.4
2	B	116	VAL	2.4
2	B	133	VAL	2.4
2	B	303	ALA	2.4
2	B	206	ILE	2.4
2	B	83	VAL	2.4
2	B	131	VAL	2.4
1	A	54	PHE	2.3
1	A	86	LEU	2.3
2	B	33	ALA	2.2
2	B	87	THR	2.2
1	A	230	ILE	2.2
1	A	371	LEU	2.2
2	B	127	ILE	2.2
2	B	174	ALA	2.2
1	A	341	ILE	2.2
2	B	58	LEU	2.2
1	A	89	TRP	2.2
2	B	34	ALA	2.2
2	B	86	LEU	2.2
2	B	108	LEU	2.2
2	B	54	TRP	2.2
2	B	170	VAL	2.1
2	B	106	TRP	2.1
2	B	155	ALA	2.1
1	A	235	ILE	2.1
1	A	529	MET	2.1
1	A	47	HIS	2.1
2	B	322	VAL	2.1
1	A	376	LEU	2.1
2	B	49	LEU	2.1
2	B	316	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	287	LEU	2.0
1	A	409	TRP	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

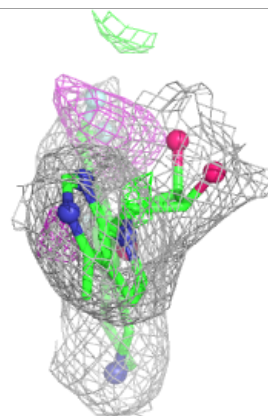
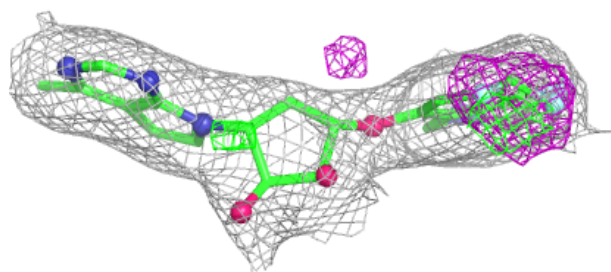
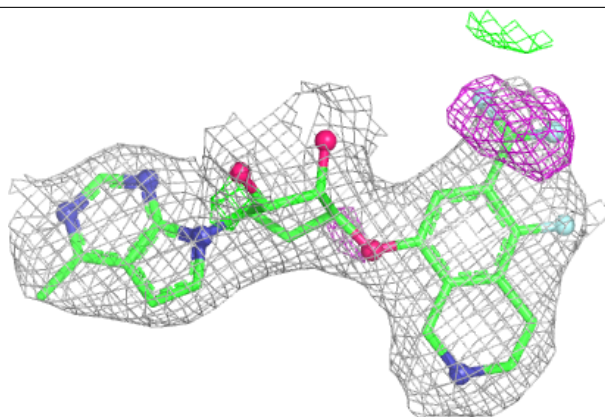
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ZR1	A	701	32/32	0.90	0.12	56,61,70,71	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 ZR1 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.