



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:42 AM UTC

PDB ID : 3M8Q / pdb_00003m8q
Title : HIV-1 RT with AMINOPYRIMIDINE NNRTI
Authors : Harris, S.F.; Villasenor, A.
Deposited on : 2010-03-18
Resolution : 2.70 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

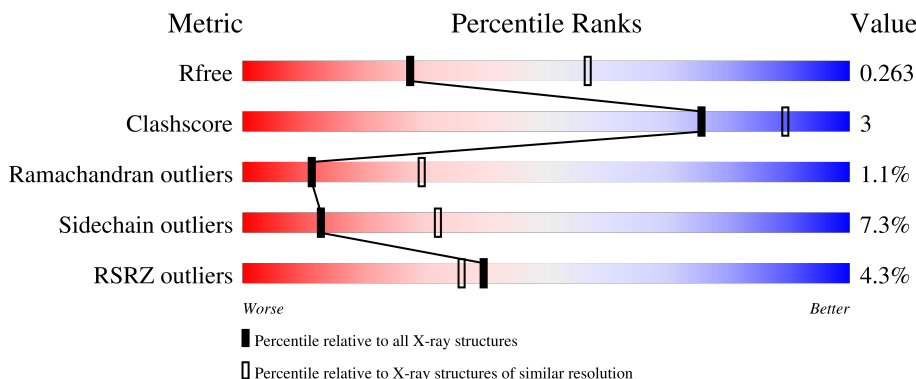
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	561	3% 83% 14% ...
2	B	440	5% 78% 11% • 10%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7926 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.

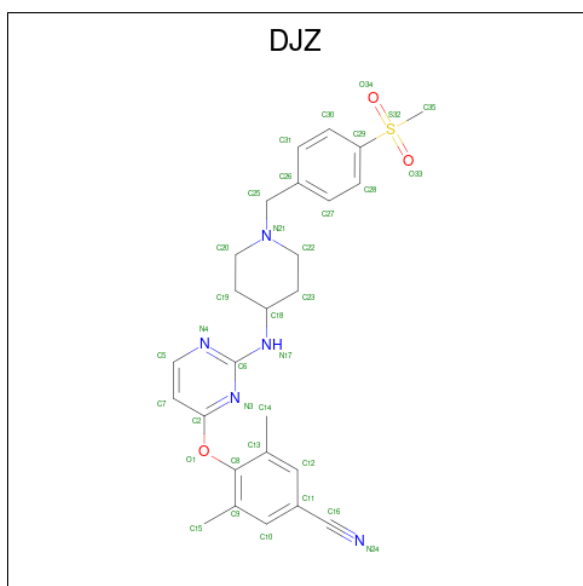
- Molecule 1 is a protein called Reverse transcriptase/ribonuclease H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	554	Total	C	N	O	S	0	1	0
			4511	2915	754	834	8			

- Molecule 2 is a protein called p51 RT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	398	Total	C	N	O	S	0	0	0
			3297	2152	539	599	7			

- Molecule 3 is 3,5-dimethyl-4-{{2-({1-[4-(methylsulfonyl)benzyl]piperidin-4-yl}amino)pyrimidin-4-yl}oxy}benzonitrile (CCD ID: DJZ) (formula: C₂₆H₂₉N₅O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			35	26	5	3	1		

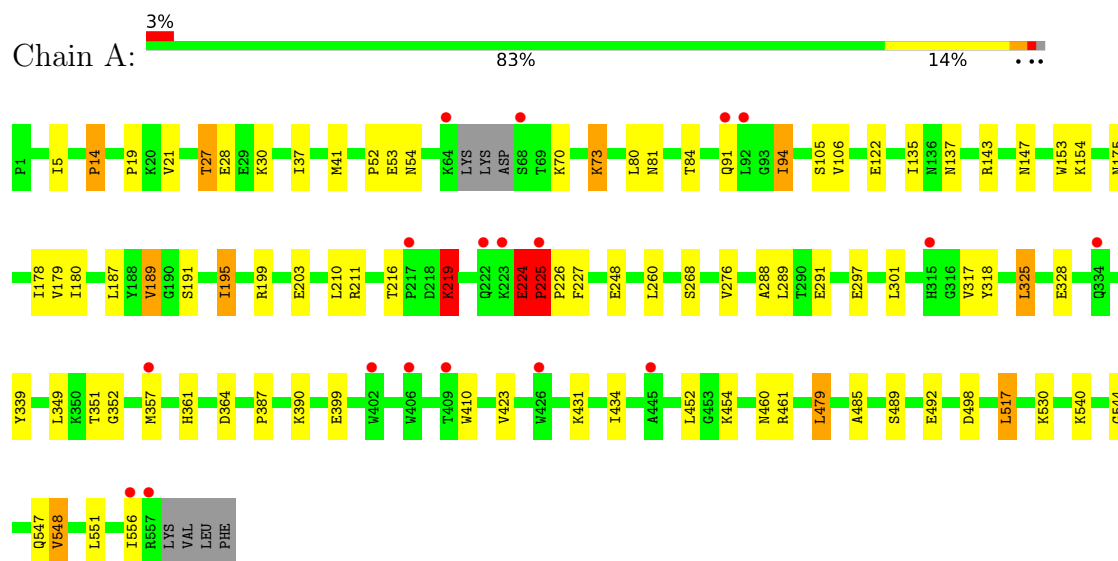
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	41	Total 41	O 41	0	0
4	B	42	Total 42	O 42	0	0

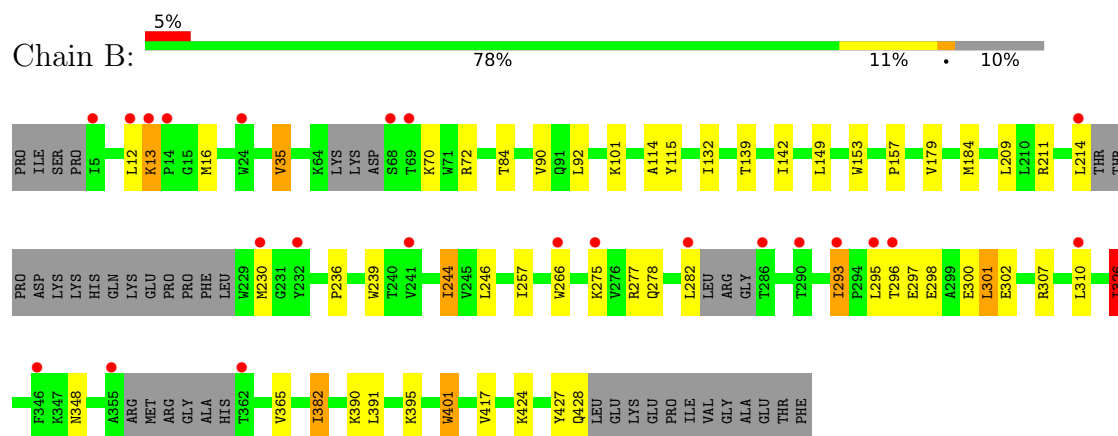
3 Residue-property plots [i](#)

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: Reverse transcriptase/ribonuclease H



- Molecule 2: p51 RT



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	118.83Å 153.10Å 154.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.09 – 2.70 47.09 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.5 (47.09-2.70) 98.5 (47.09-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.69Å)	Xtriage
Refinement program	BUSTER 2.9.2	Depositor
R, R_{free}	0.219 , 0.252 0.225 , 0.263	Depositor DCC
R_{free} test set	1976 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å ²)	62.3	Xtriage
Anisotropy	0.422	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 58.3	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	7926	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: DJZ

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	0.73	3/4628 (0.1%)	1.28	13/6289 (0.2%)
2	B	0.73	0/3390	1.28	18/4607 (0.4%)
All	All	0.73	3/8018 (0.0%)	1.28	31/10896 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	224	GLU	CA-C	5.43	1.59	1.52
1	A	225	PRO	CA-C	5.31	1.57	1.52
1	A	224	GLU	C-N	5.01	1.39	1.33

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	225	PRO	N-CA-C	12.26	125.66	110.70
2	B	211	ARG	N-CA-C	-8.54	96.19	109.76
1	A	349	LEU	N-CA-C	-6.38	103.62	111.40
2	B	401	TRP	N-CA-C	6.37	121.34	113.50
2	B	132	ILE	N-CA-C	-5.78	101.83	107.55
1	A	147	ASN	N-CA-C	-5.69	106.68	113.97
1	A	153	TRP	CA-C-N	5.58	128.03	120.38
1	A	153	TRP	C-N-CA	5.58	128.03	120.38
2	B	296	THR	CA-C-N	5.49	128.42	120.79
2	B	296	THR	C-N-CA	5.49	128.42	120.79
2	B	348	ASN	N-CA-C	5.42	118.03	109.96
2	B	293	ILE	CB-CA-C	5.41	115.37	110.13
1	A	94	ILE	CB-CA-C	5.32	116.06	110.37
1	A	498	ASP	CA-CB-CG	5.29	117.89	112.60
2	B	382	ILE	N-CA-C	5.25	115.91	110.82
2	B	35	VAL	N-CA-CB	5.24	117.67	110.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	14	PRO	N-CA-C	5.22	123.22	112.47
2	B	139	THR	CB-CA-C	5.22	116.72	108.84
2	B	395	LYS	CA-C-N	5.20	127.19	120.44
2	B	395	LYS	C-N-CA	5.20	127.19	120.44
1	A	297	GLU	CA-C-N	5.18	127.22	120.28
1	A	297	GLU	C-N-CA	5.18	127.22	120.28
2	B	391	LEU	N-CA-C	5.18	117.08	109.42
2	B	301	LEU	CA-C-N	5.17	127.52	120.54
2	B	301	LEU	C-N-CA	5.17	127.52	120.54
1	A	52	PRO	CA-C-N	5.15	127.18	120.28
1	A	52	PRO	C-N-CA	5.15	127.18	120.28
2	B	326	ILE	CB-CA-C	5.13	118.27	110.62
1	A	225	PRO	CB-CA-C	5.09	117.14	110.92
2	B	302	GLU	CA-C-N	5.09	127.06	120.44
2	B	302	GLU	C-N-CA	5.09	127.06	120.44

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	4511	0	4545	32	0
2	B	3297	0	3313	16	0
3	A	35	0	29	3	0
4	A	41	0	0	0	0
4	B	42	0	0	0	0
All	All	7926	0	7887	47	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 3.

All (47) 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:224:GLU:HB2	1:A:225:PRO:HD2	1.56	0.88
1:A:224:GLU:HB2	1:A:225:PRO:CD	2.07	0.85
1:A:106:VAL:HG21	3:A:562:DJZ:H22	1.74	0.70
1:A:37:ILE:HG22	1:A:41:MET:HE2	1.79	0.64
1:A:227:PHE:HE2	3:A:562:DJZ:H35B	1.68	0.59
1:A:224:GLU:CB	1:A:225:PRO:HD2	2.32	0.58
1:A:460:ASN:HD21	1:A:461:ARG:HH11	1.51	0.58
2:B:246:LEU:HD12	2:B:307:ARG:HG3	1.86	0.58
1:A:225:PRO:CB	1:A:226:PRO:HD3	2.33	0.57
1:A:328:GLU:HG2	1:A:390:LYS:HB2	1.86	0.56
1:A:19:PRO:HG3	1:A:80:LEU:HB2	1.87	0.56
1:A:410:TRP:HB3	2:B:365:VAL:HG23	1.89	0.54
2:B:326:ILE:HD11	2:B:390:LYS:HG3	1.93	0.50
1:A:41:MET:HE1	1:A:73:LYS:HZ2	1.77	0.50
1:A:175:ASN:HB3	1:A:178:ILE:HG12	1.94	0.50
2:B:101:LYS:HD3	2:B:382:ILE:HG23	1.95	0.49
1:A:84:THR:HG23	1:A:154:LYS:HD3	1.96	0.48
1:A:325:LEU:HB3	1:A:387:PRO:HB3	1.96	0.47
1:A:288:ALA:HB3	1:A:291:GLU:HB2	1.97	0.47
1:A:364:ASP:HB3	1:A:423:VAL:HG13	1.98	0.46
1:A:81:ASN:HA	1:A:84:THR:HG22	1.98	0.46
2:B:114:ALA:HB2	2:B:214:LEU:HD11	1.98	0.46
1:A:485:ALA:O	1:A:489:SER:HB2	2.16	0.46
2:B:12:LEU:O	2:B:13:LYS:HB2	2.16	0.45
2:B:244:ILE:HB	2:B:310:LEU:HD22	1.98	0.45
1:A:225:PRO:HB3	1:A:226:PRO:HD3	1.97	0.45
1:A:199:ARG:HG2	1:A:219:LYS:HE2	1.97	0.45
2:B:209:LEU:HB3	2:B:214:LEU:HB2	1.98	0.45
1:A:479:LEU:HB3	1:A:517:LEU:HD13	1.99	0.45
2:B:115:TYR:HB3	2:B:149:LEU:HB2	1.99	0.45
2:B:84:THR:HG21	2:B:153:TRP:HZ2	1.82	0.44
2:B:278:GLN:HG2	2:B:298:GLU:HB2	1.99	0.44
1:A:318:TYR:CZ	3:A:562:DJZ:H20	2.53	0.44
1:A:54:ASN:HB3	1:A:143:ARG:HH21	1.82	0.43
1:A:225:PRO:CB	1:A:226:PRO:CD	2.95	0.43
2:B:157:PRO:HG3	2:B:184:MET:HA	2.00	0.43
1:A:492:GLU:HG2	1:A:530:LYS:HB2	2.01	0.43
1:A:27:THR:HG22	1:A:30:LYS:H	1.84	0.43
2:B:266:TRP:HZ2	2:B:427:TYR:CZ	2.37	0.43
1:A:544:GLY:O	1:A:548:VAL:HG23	2.18	0.42
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.54	0.42
2:B:266:TRP:CZ2	2:B:427:TYR:CZ	3.08	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:GLU:HG2	1:A:135:ILE:HG12	2.02	0.42
1:A:339:TYR:CZ	1:A:352:GLY:HA3	2.55	0.42
1:A:180:ILE:HG12	1:A:189:VAL:HG13	2.02	0.42
2:B:365:VAL:HG11	2:B:401:TRP:HB2	2.02	0.40
1:A:268:SER:O	1:A:351:THR:HG22	2.22	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	551/561 (98%)	523 (95%)	22 (4%)	6 (1%)	11	29
2	B	388/440 (88%)	370 (95%)	14 (4%)	4 (1%)	12	32
All	All	939/1001 (94%)	893 (95%)	36 (4%)	10 (1%)	11	29

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	224	GLU
1	A	225	PRO
2	B	13	LYS
1	A	195	ILE
1	A	14	PRO
1	A	219	LYS
2	B	16	MET
2	B	230	MET
2	B	277	ARG
1	A	556	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	493/501 (98%)	451 (92%)	42 (8%)	10	25
2	B	363/400 (91%)	343 (94%)	20 (6%)	19	45
All	All	856/901 (95%)	794 (93%)	62 (7%)	13	32

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ILE
1	A	21	VAL
1	A	27	THR
1	A	53	GLU
1	A	70	LYS
1	A	73	LYS
1	A	91	GLN
1	A	94	ILE
1	A	105	SER
1	A	122	GLU
1	A	137	ASN
1	A	179	VAL
1	A	187	LEU
1	A	189	VAL
1	A	191	SER
1	A	195	ILE
1	A	203	GLU
1	A	210	LEU
1	A	211	ARG
1	A	216	THR
1	A	219	LYS
1	A	224	GLU
1	A	248	GLU
1	A	260	LEU
1	A	276	VAL
1	A	289	LEU
1	A	301	LEU

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Mol	Chain	Res	Type
1	A	317	VAL
1	A	325	LEU
1	A	357	MET
1	A	361	HIS
1	A	399	GLU
1	A	431	LYS
1	A	434	ILE
1	A	452	LEU
1	A	454	LYS
1	A	479	LEU
1	A	517	LEU
1	A	540	LYS
1	A	547	GLN
1	A	548	VAL
1	A	551	LEU
2	B	35	VAL
2	B	70	LYS
2	B	72	ARG
2	B	90	VAL
2	B	92	LEU
2	B	142	ILE
2	B	179	VAL
2	B	244	ILE
2	B	257	ILE
2	B	275	LYS
2	B	282	LEU
2	B	293	ILE
2	B	295	LEU
2	B	297	GLU
2	B	300	GLU
2	B	301	LEU
2	B	326	ILE
2	B	417	VAL
2	B	424	LYS
2	B	428	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
1	A	91	GLN
1	A	137	ASN

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Mol	Chain	Res	Type
1	A	147	ASN
1	A	255	ASN
1	A	334	GLN
1	A	407	GLN
1	A	500	GLN
1	A	509	GLN
2	B	151	GLN
2	B	258	GLN
2	B	269	GLN
2	B	315	HIS
2	B	332	GLN
2	B	394	GLN

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	DJZ	A	562	-	38,38,38	3.36	21 (55%)	54,54,54	3.76	23 (42%)

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	DJZ	A	562	-	-	6/20/30/30	0/4/4/4

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	562	DJZ	C29-S32	8.00	1.85	1.77
3	A	562	DJZ	C31-C26	7.00	1.52	1.38
3	A	562	DJZ	C6-N4	6.31	1.43	1.34
3	A	562	DJZ	C30-C29	6.08	1.48	1.38
3	A	562	DJZ	C28-C29	6.05	1.48	1.38
3	A	562	DJZ	C7-C5	5.65	1.49	1.38
3	A	562	DJZ	C2-N3	5.24	1.42	1.33
3	A	562	DJZ	C12-C11	4.28	1.47	1.39
3	A	562	DJZ	C10-C9	4.11	1.45	1.39
3	A	562	DJZ	C10-C11	3.59	1.45	1.39
3	A	562	DJZ	C5-N4	3.28	1.41	1.34
3	A	562	DJZ	C23-C18	3.10	1.59	1.52
3	A	562	DJZ	C8-C13	3.03	1.45	1.40
3	A	562	DJZ	C8-C9	2.82	1.45	1.40
3	A	562	DJZ	C6-N17	2.54	1.37	1.34
3	A	562	DJZ	C28-C27	2.50	1.42	1.38
3	A	562	DJZ	C15-C9	2.50	1.55	1.51
3	A	562	DJZ	C27-C26	2.39	1.43	1.38
3	A	562	DJZ	C35-S32	2.33	1.83	1.75
3	A	562	DJZ	C31-C30	2.29	1.42	1.38
3	A	562	DJZ	C6-N3	2.10	1.40	1.34

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	562	DJZ	O33-S32-C29	-14.61	96.87	108.23
3	A	562	DJZ	O34-S32-C35	12.02	126.10	108.47
3	A	562	DJZ	C6-N3-C2	8.87	125.14	115.00
3	A	562	DJZ	C27-C28-C29	6.58	125.84	119.44
3	A	562	DJZ	N4-C6-N3	-6.49	120.14	126.42
3	A	562	DJZ	C5-C7-C2	6.21	119.19	115.48
3	A	562	DJZ	C28-C27-C26	-4.02	115.71	121.00
3	A	562	DJZ	C20-C19-C18	-3.99	104.38	110.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	562	DJZ	O34-S32-O33	-3.68	111.74	117.99
3	A	562	DJZ	C23-C18-N17	-3.67	104.86	110.77
3	A	562	DJZ	C14-C13-C12	-3.45	113.49	119.57
3	A	562	DJZ	C7-C2-N3	-3.43	116.37	123.48
3	A	562	DJZ	C12-C13-C8	3.36	122.80	117.94
3	A	562	DJZ	C12-C11-C16	3.09	123.72	119.55
3	A	562	DJZ	O34-S32-C29	-2.64	106.17	108.23
3	A	562	DJZ	C10-C11-C16	-2.61	116.03	119.55
3	A	562	DJZ	C23-C18-C19	2.32	114.81	110.80
3	A	562	DJZ	C30-C29-S32	2.27	121.59	119.58
3	A	562	DJZ	C25-C26-C27	-2.24	116.62	120.75
3	A	562	DJZ	N17-C6-N4	2.16	120.17	116.75
3	A	562	DJZ	C31-C30-C29	-2.06	117.43	119.44
3	A	562	DJZ	C22-C23-C18	-2.06	107.42	110.67
3	A	562	DJZ	C11-C12-C13	-2.01	118.17	121.06

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	562	DJZ	C28-C29-S32-C35
3	A	562	DJZ	C30-C29-S32-C35
3	A	562	DJZ	C28-C29-S32-O33
3	A	562	DJZ	C30-C29-S32-O33
3	A	562	DJZ	C28-C29-S32-O34
3	A	562	DJZ	C30-C29-S32-O34

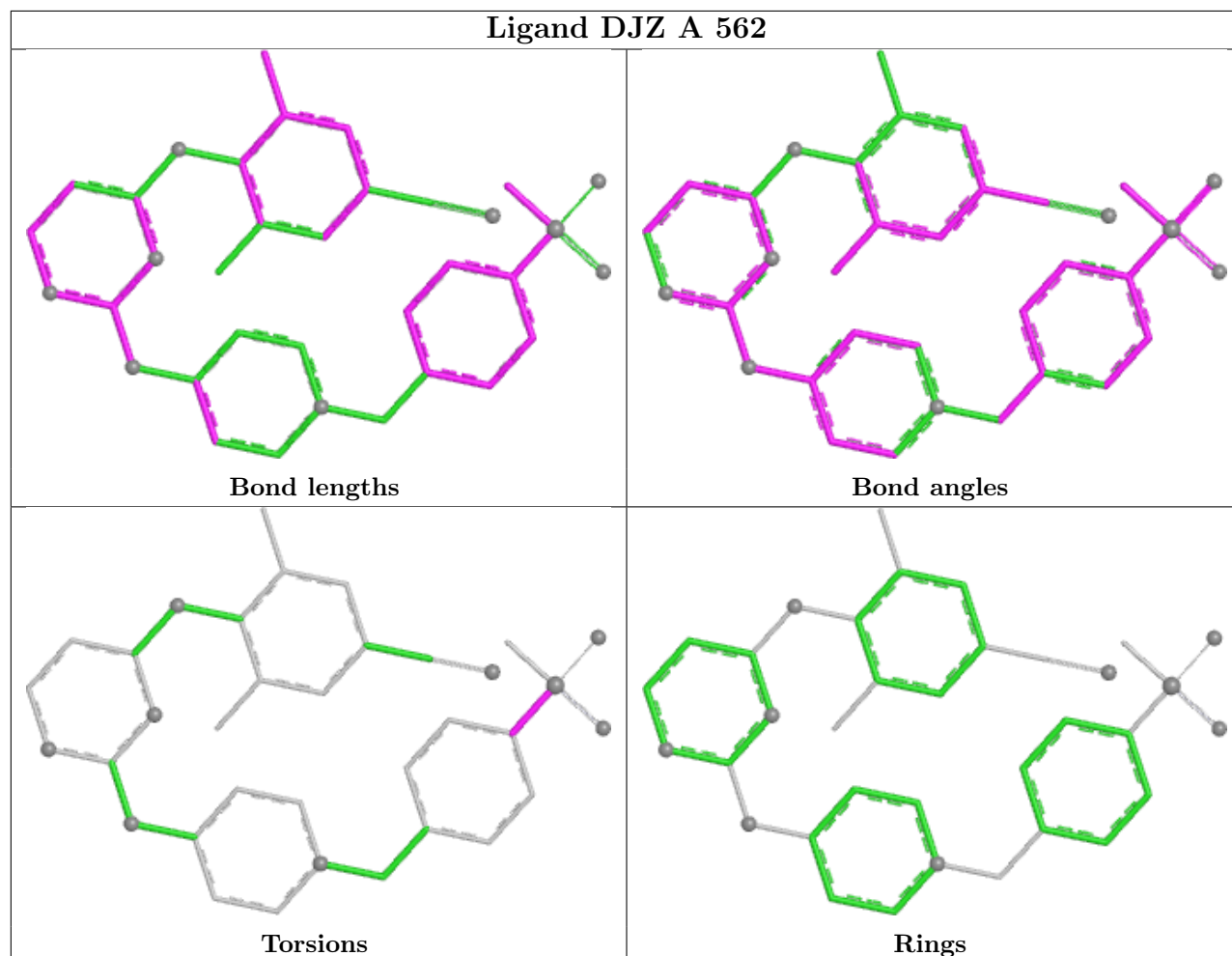
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	562	DJZ	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	554/561 (98%)	0.36	18 (3%)	50	46	31, 70, 100, 138	1 (0%)
2	B	398/440 (90%)	0.33	23 (5%)	29	25	43, 65, 105, 123	0
All	All	952/1001 (95%)	0.35	41 (4%)	40	36	31, 68, 103, 138	1 (0%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	282	LEU	4.9
2	B	266	TRP	4.2
2	B	12	LEU	4.2
2	B	5	ILE	4.0
1	A	409	THR	3.9
1	A	223	LYS	3.6
1	A	225	PRO	3.5
1	A	92	LEU	3.5
1	A	557	ARG	3.4
2	B	293	ILE	3.2
1	A	315[A]	HIS	3.1
2	B	290	THR	3.1
2	B	13	LYS	3.1
2	B	346	PHE	3.1
2	B	355	ALA	3.0
1	A	217	PRO	2.9
1	A	91	GLN	2.8
1	A	556	ILE	2.8
2	B	214	LEU	2.8
1	A	222	GLN	2.7
1	A	406	TRP	2.7
2	B	232	TYR	2.7
2	B	295	LEU	2.7
1	A	357	MET	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	362	THR	2.6
1	A	426	TRP	2.5
2	B	24	TRP	2.5
1	A	445	ALA	2.5
1	A	68	SER	2.4
2	B	14	PRO	2.4
2	B	275	LYS	2.3
2	B	286	THR	2.3
2	B	230	MET	2.3
1	A	402	TRP	2.1
1	A	334	GLN	2.1
1	A	64	LYS	2.1
2	B	296	THR	2.1
2	B	310	LEU	2.1
2	B	241	VAL	2.1
2	B	69	THR	2.0
2	B	68	SER	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

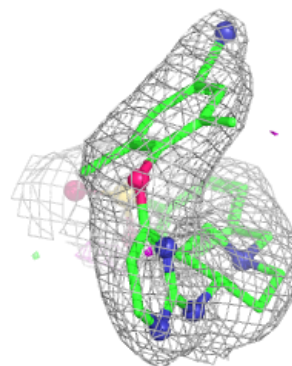
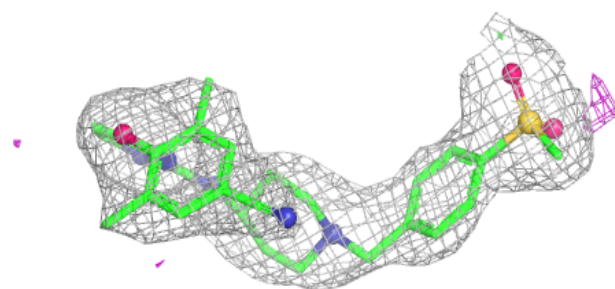
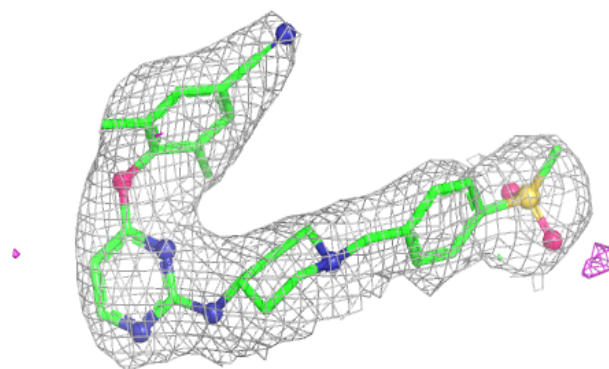
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	DJZ	A	562	35/35	0.95	0.09	49,54,70,73	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 DJZ A 562:

$2mF_o - DF_c$ (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.