



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 04:05 PM UTC

PDB ID : 7SNP / pdb_00007snp
Title : Crystal Structure of HIV-1 Reverse Transcriptase in Complex with (E)-3-(3-chloro-5-(2-(2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy)-4-(2-morpholinoethoxy)phenoxy)phenyl)acrylonitrile (JLJ530)
Authors : Frey, K.M.; Anderson, K.S.
Deposited on : 2021-10-28
Resolution : 2.89 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

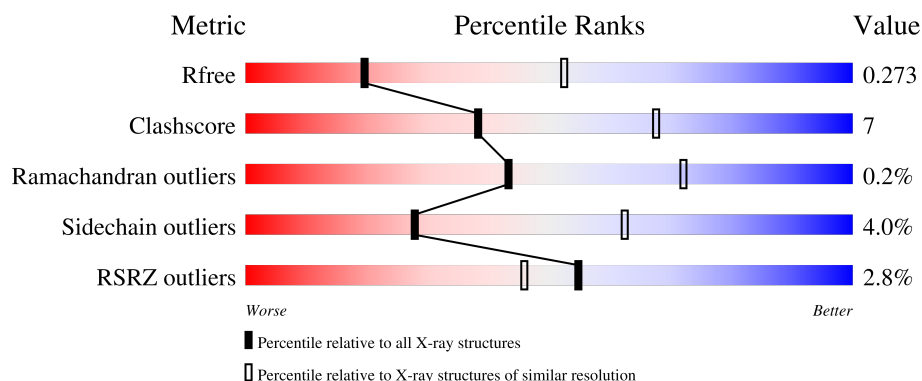
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	557	2% 77% 20% ..
2	B	428	3% 79% 19% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8076 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 p66.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	549	Total	C	N	O	S	0	0	0
			4465	2891	742	825	7			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	expression tag	UNP P03366
A	0	VAL	-	expression tag	UNP P03366
A	172	ALA	LYS	engineered mutation	UNP P03366
A	173	ALA	LYS	engineered mutation	UNP P03366
A	280	SER	CYS	engineered mutation	UNP P03366

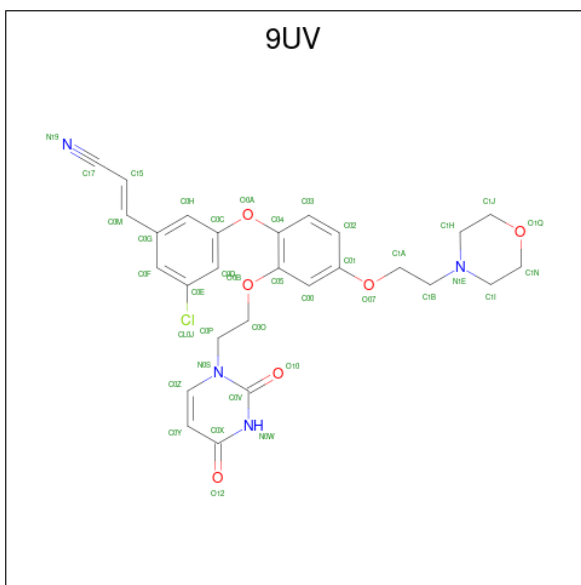
- Molecule 2 is a protein called p51 RT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	428	Total	C	N	O	S	0	0	0
			3538	2305	586	640	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	SER	CYS	conflict	UNP P03366

- Molecule 3 is (2E)-3-(3-chloro-5-{2-[2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy]-4-[2-(morpholin-4-yl)ethoxy]phenoxy}phenyl)prop-2-enenitrile (CCD ID: 9UV) (formula: C₂₇H₂₇ClN₄O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			38	27	1	4	6		

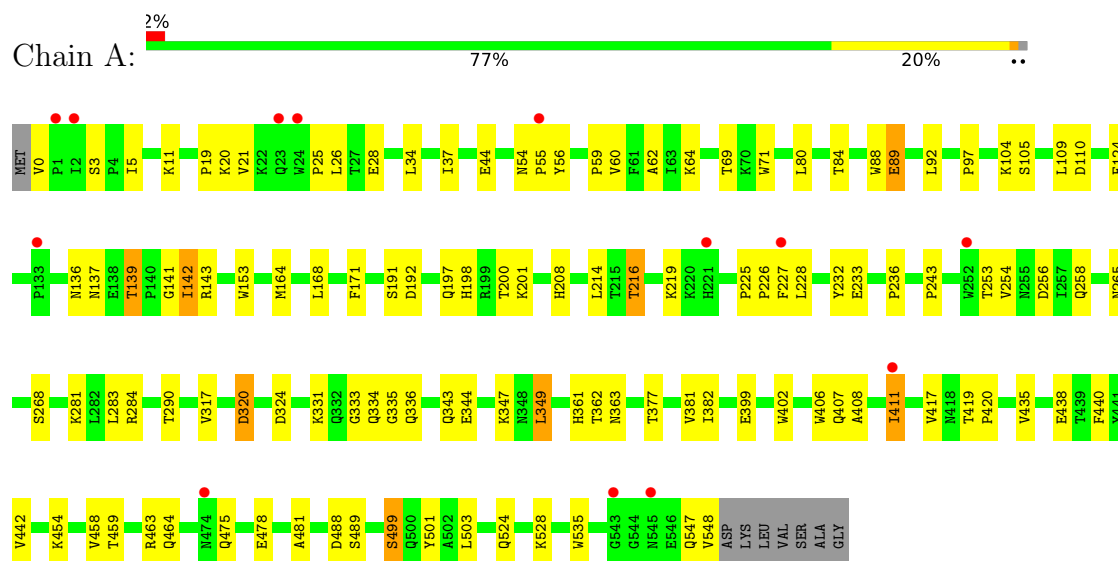
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	19	Total O 19 19	0	0
4	B	16	Total O 16 16	0	0

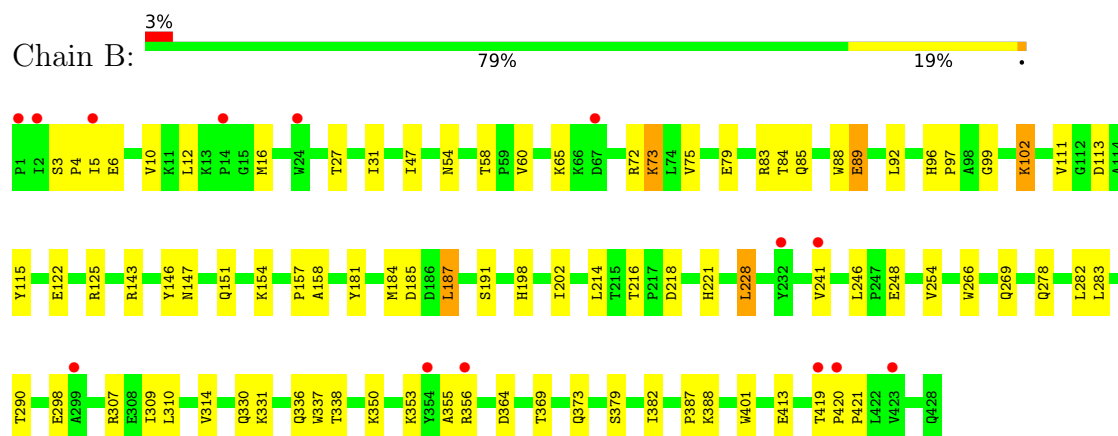
3 Residue-property plots [i](#)

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 p66



• Molecule 2: p51 RT



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	224.82Å 69.05Å 104.34Å 90.00° 106.29° 90.00°	Depositor
Resolution (Å)	36.60 – 2.89 36.60 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.5 (36.60-2.89) 99.5 (36.60-2.89)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.224 , 0.272 0.229 , 0.273	Depositor DCC
R_{free} test set	2000 reflections (5.78%)	wwPDB-VP
Wilson B-factor (Å ²)	83.8	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8076	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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: 9UV

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	0.35	0/4583	0.79	4/6231 (0.1%)
2	B	0.37	0/3643	0.85	9/4952 (0.2%)
All	All	0.36	0/8226	0.82	13/11183 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	96	HIS	CA-C-N	8.04	129.89	119.84
2	B	96	HIS	C-N-CA	8.04	129.89	119.84
2	B	419	THR	N-CA-C	7.82	122.81	112.17
1	A	335	GLY	N-CA-C	-6.39	106.31	115.64
2	B	54	ASN	CA-C-N	6.00	125.70	119.64
2	B	54	ASN	C-N-CA	6.00	125.70	119.64
1	A	320	ASP	CA-C-N	5.71	126.09	119.47
1	A	320	ASP	C-N-CA	5.71	126.09	119.47
1	A	382	ILE	N-CA-C	5.67	116.32	110.82
2	B	382	ILE	N-CA-C	5.52	115.61	110.42
2	B	420	PRO	N-CA-C	5.13	116.97	110.70
2	B	401	TRP	N-CA-C	5.06	119.55	113.17
2	B	89	GLU	N-CA-C	-5.00	107.18	113.28

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	4465	0	4516	70	0
2	B	3538	0	3576	45	0
3	A	38	0	0	0	0
4	A	19	0	0	0	0
4	B	16	0	0	0	0
All	All	8076	0	8092	111	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 (111) 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:191:SER:HG	1:A:198:HIS:HD1	1.19	0.85
2:B:85:GLN:HA	2:B:88:TRP:HB2	1.64	0.77
1:A:84:THR:HG21	1:A:153:TRP:HE1	1.49	0.77
1:A:331:LYS:HE2	1:A:333:GLY:HA2	1.69	0.75
2:B:85:GLN:O	2:B:89:GLU:N	2.26	0.69
2:B:157:PRO:HG3	2:B:184:MET:HA	1.76	0.67
1:A:458:VAL:HG22	1:A:464:GLN:HG2	1.77	0.66
1:A:281:LYS:HG3	1:A:284:ARG:HE	1.61	0.65
2:B:125:ARG:HE	2:B:147:ASN:HA	1.62	0.65
1:A:438:GLU:OE1	1:A:463:ARG:NH2	2.31	0.63
1:A:84:THR:HG22	1:A:124:PHE:HZ	1.61	0.63
1:A:197:GLN:HA	1:A:200:THR:HG22	1.81	0.62
1:A:320:ASP:O	1:A:343:GLN:NE2	2.33	0.61
1:A:226:PRO:HA	1:A:227:PHE:HB2	1.83	0.59
2:B:122:GLU:OE1	2:B:125:ARG:NH1	2.36	0.58
1:A:19:PRO:HG3	1:A:80:LEU:HB2	1.86	0.57
1:A:164:MET:HE1	1:A:214:LEU:HD13	1.86	0.57
2:B:115:TYR:OH	2:B:184:MET:O	2.20	0.56
2:B:65:LYS:HA	2:B:72:ARG:HD2	1.86	0.56
2:B:73:LYS:NZ	2:B:146:TYR:OH	2.36	0.55
1:A:334:GLN:HB2	1:A:336:GLN:HE21	1.72	0.55
2:B:85:GLN:HG3	2:B:154:LYS:HB3	1.88	0.55
1:A:524:GLN:O	1:A:528:LYS:HG2	2.07	0.54
1:A:89:GLU:HG2	1:A:92:LEU:HD21	1.90	0.54
1:A:399:GLU:HA	1:A:402:TRP:CD1	2.43	0.54
2:B:241:VAL:HG22	2:B:350:LYS:HA	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:ASN:O	1:A:268:SER:OG	2.23	0.53
1:A:406:TRP:CH2	1:A:407:GLN:HG3	2.45	0.53
1:A:171:PHE:HB2	1:A:208:HIS:ND1	2.24	0.52
1:A:399:GLU:HA	1:A:402:TRP:NE1	2.24	0.52
1:A:253:THR:HG23	1:A:256:ASP:H	1.75	0.52
1:A:475:GLN:HB3	1:A:501:TYR:CE2	2.44	0.51
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.92	0.51
2:B:79:GLU:OE2	2:B:83:ARG:NH1	2.44	0.51
2:B:266:TRP:O	2:B:269:GLN:HG2	2.11	0.51
1:A:64:LYS:HZ3	1:A:71:TRP:CD1	2.28	0.50
1:A:362:THR:OG1	1:A:363:ASN:N	2.43	0.50
1:A:435:VAL:HG22	2:B:290:THR:HG21	1.94	0.50
1:A:478:GLU:HG2	1:A:499:SER:OG	2.12	0.49
2:B:369:THR:HG22	2:B:373:GLN:HE21	1.78	0.48
2:B:12:LEU:HD22	2:B:84:THR:HG22	1.96	0.48
2:B:221:HIS:CE1	2:B:228:LEU:HD22	2.48	0.48
2:B:84:THR:OG1	2:B:85:GLN:N	2.46	0.48
1:A:84:THR:HG21	1:A:153:TRP:NE1	2.25	0.48
1:A:137:ASN:HD22	1:A:139:THR:HG22	1.78	0.47
1:A:54:ASN:N	1:A:55:PRO:HD3	2.29	0.47
1:A:105:SER:HB2	1:A:198:HIS:CE1	2.49	0.47
2:B:92:LEU:HB3	2:B:158:ALA:HB1	1.95	0.47
2:B:278:GLN:HG3	2:B:298:GLU:HG2	1.96	0.47
2:B:111:VAL:HG11	2:B:187:LEU:HD22	1.96	0.47
2:B:151:GLN:HB3	2:B:185:ASP:OD2	2.15	0.46
1:A:343:GLN:HG3	1:A:349:LEU:HD21	1.97	0.46
2:B:27:THR:O	2:B:31:ILE:HG13	2.16	0.46
2:B:10:VAL:HG13	2:B:88:TRP:CZ2	2.51	0.46
1:A:225:PRO:HB2	1:A:236:PRO:HD3	1.98	0.46
1:A:19:PRO:O	1:A:56:TYR:HB3	2.15	0.46
2:B:216:THR:C	2:B:218:ASP:H	2.24	0.46
2:B:97:PRO:HG3	2:B:181:TYR:HB2	1.98	0.46
1:A:80:LEU:O	1:A:84:THR:HG23	2.15	0.46
1:A:104:LYS:HB2	1:A:192:ASP:HA	1.98	0.46
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.98	0.45
2:B:47:ILE:HG22	2:B:146:TYR:HA	1.97	0.45
2:B:191:SER:OG	2:B:198:HIS:ND1	2.28	0.45
2:B:338:THR:HG22	2:B:353:LYS:HB2	1.97	0.45
1:A:201:LYS:HD3	1:A:201:LYS:HA	1.72	0.45
1:A:317:VAL:HG11	1:A:347:LYS:HD2	1.99	0.45
1:A:478:GLU:HG2	1:A:499:SER:CB	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:440:PHE:CE1	1:A:489:SER:HB3	2.52	0.45
2:B:248:GLU:HA	2:B:307:ARG:HH22	1.82	0.45
1:A:88:TRP:NE1	2:B:143:ARG:HD2	2.31	0.44
2:B:336:GLN:HG2	2:B:355:ALA:HB2	2.00	0.44
1:A:324:ASP:O	1:A:343:GLN:HG2	2.17	0.44
1:A:417:VAL:HG12	1:A:419:THR:H	1.83	0.44
1:A:21:VAL:HB	1:A:59:PRO:HD3	2.00	0.43
1:A:25:PRO:HA	1:A:26:LEU:HA	1.76	0.43
1:A:377:THR:O	1:A:381:VAL:HG23	2.18	0.43
1:A:34:LEU:HD21	1:A:62:ALA:HB2	2.00	0.43
1:A:139:THR:O	1:A:141:GLY:N	2.52	0.43
2:B:10:VAL:HG22	2:B:88:TRP:CH2	2.53	0.43
1:A:454:LYS:NZ	1:A:548:VAL:HG23	2.34	0.43
2:B:198:HIS:O	2:B:202:ILE:HG12	2.18	0.43
1:A:408:ALA:HB1	2:B:364:ASP:HB3	2.00	0.43
2:B:388:LYS:HG3	2:B:413:GLU:HB2	2.00	0.43
2:B:99:GLY:HA2	2:B:102:LYS:HG3	2.01	0.42
2:B:331:LYS:HB2	2:B:337:TRP:CZ3	2.53	0.42
1:A:25:PRO:HB2	1:A:26:LEU:HB2	2.02	0.42
1:A:228:LEU:HD12	1:A:233:GLU:HG2	2.01	0.42
1:A:258:GLN:HG2	1:A:283:LEU:HD21	2.01	0.42
1:A:233:GLU:OE2	1:A:243:PRO:HD3	2.20	0.42
2:B:282:LEU:HD12	2:B:282:LEU:HA	1.93	0.42
2:B:246:LEU:HD12	2:B:307:ARG:HG2	2.01	0.42
2:B:379:SER:CB	2:B:387:PRO:HD3	2.50	0.42
1:A:344:GLU:HB2	1:A:347:LYS:HG3	2.02	0.42
1:A:442:VAL:HB	1:A:481:ALA:HB1	2.01	0.42
1:A:419:THR:HA	1:A:420:PRO:HD3	1.92	0.41
1:A:478:GLU:HG2	1:A:499:SER:HB3	2.01	0.41
1:A:0:VAL:HG13	1:A:44:GLU:HA	2.03	0.41
1:A:109:LEU:HB3	1:A:216:THR:HG22	2.03	0.41
2:B:254:VAL:HG13	2:B:283:LEU:HD22	2.03	0.41
1:A:55:PRO:HD2	1:A:143:ARG:HH12	1.85	0.41
1:A:28:GLU:CD	1:A:28:GLU:H	2.29	0.41
1:A:411:ILE:HD12	1:A:411:ILE:O	2.21	0.41
1:A:142:ILE:H	1:A:142:ILE:HG13	1.59	0.40
1:A:281:LYS:HA	1:A:284:ARG:HG3	2.01	0.40
1:A:137:ASN:HB3	1:A:139:THR:HG23	2.03	0.40
1:A:435:VAL:HG13	2:B:290:THR:HG21	2.03	0.40
2:B:65:LYS:HE2	2:B:65:LYS:HB2	1.96	0.40
1:A:97:PRO:HG2	1:A:232:TYR:CD2	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:3:SER:HA	2:B:4:PRO:HD2	1.93	0.40
1:A:3:SER:HB3	1:A:5:ILE:HG13	2.04	0.40
1:A:463:ARG:NH1	1:A:488:ASP:O	2.55	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	547/557 (98%)	515 (94%)	31 (6%)	1 (0%)	43	72
2	B	426/428 (100%)	404 (95%)	21 (5%)	1 (0%)	43	72
All	All	973/985 (99%)	919 (94%)	52 (5%)	2 (0%)	43	72

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	136	ASN
2	B	421	PRO

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	489/495 (99%)	469 (96%)	20 (4%)	27	61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	390/390 (100%)	375 (96%)	15 (4%)	29	64
All	All	879/885 (99%)	844 (96%)	35 (4%)	28	62

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LYS
1	A	20	LYS
1	A	37	ILE
1	A	60	VAL
1	A	69	THR
1	A	89	GLU
1	A	110	ASP
1	A	139	THR
1	A	142	ILE
1	A	168	LEU
1	A	216	THR
1	A	219	LYS
1	A	254	VAL
1	A	290	THR
1	A	349	LEU
1	A	361	HIS
1	A	411	ILE
1	A	459	THR
1	A	499	SER
1	A	547	GLN
2	B	5	ILE
2	B	6	GLU
2	B	16	MET
2	B	58	THR
2	B	73	LYS
2	B	102	LYS
2	B	113	ASP
2	B	187	LEU
2	B	214	LEU
2	B	228	LEU
2	B	309	ILE
2	B	310	LEU
2	B	314	VAL
2	B	330	GLN
2	B	356	ARG

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	137	ASN
1	A	197	GLN
1	A	361	HIS
1	A	428	GLN
1	A	487	GLN
1	A	519	ASN
1	A	539	HIS
2	B	147	ASN
2	B	174	GLN
2	B	221	HIS
2	B	255	ASN
2	B	269	GLN
2	B	330	GLN
2	B	340	GLN
2	B	428	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	9UV	A	601	-	41,41,41	1.31	2 (4%)	53,54,54	1.69	8 (15%)

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	9UV	A	601	-	-	2/19/28/28	0/4/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	9UV	C0Z-N0S	3.85	1.42	1.37
3	A	601	9UV	C02-C01	2.51	1.43	1.38

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	9UV	O0B-C05-C04	5.80	127.75	115.75
3	A	601	9UV	O0B-C05-C00	-4.75	113.24	123.49
3	A	601	9UV	C0G-C0H-C0C	3.14	123.01	119.98
3	A	601	9UV	C04-O0A-C0C	3.09	125.43	117.95
3	A	601	9UV	C0P-N0S-C0V	2.76	121.22	118.48
3	A	601	9UV	C0O-O0B-C05	2.37	123.40	117.69
3	A	601	9UV	O0A-C0C-C0D	2.17	125.96	119.16
3	A	601	9UV	C0H-C0G-C0F	-2.00	116.39	119.11

There are no chirality outliers.

All (2) torsion outliers are listed below:

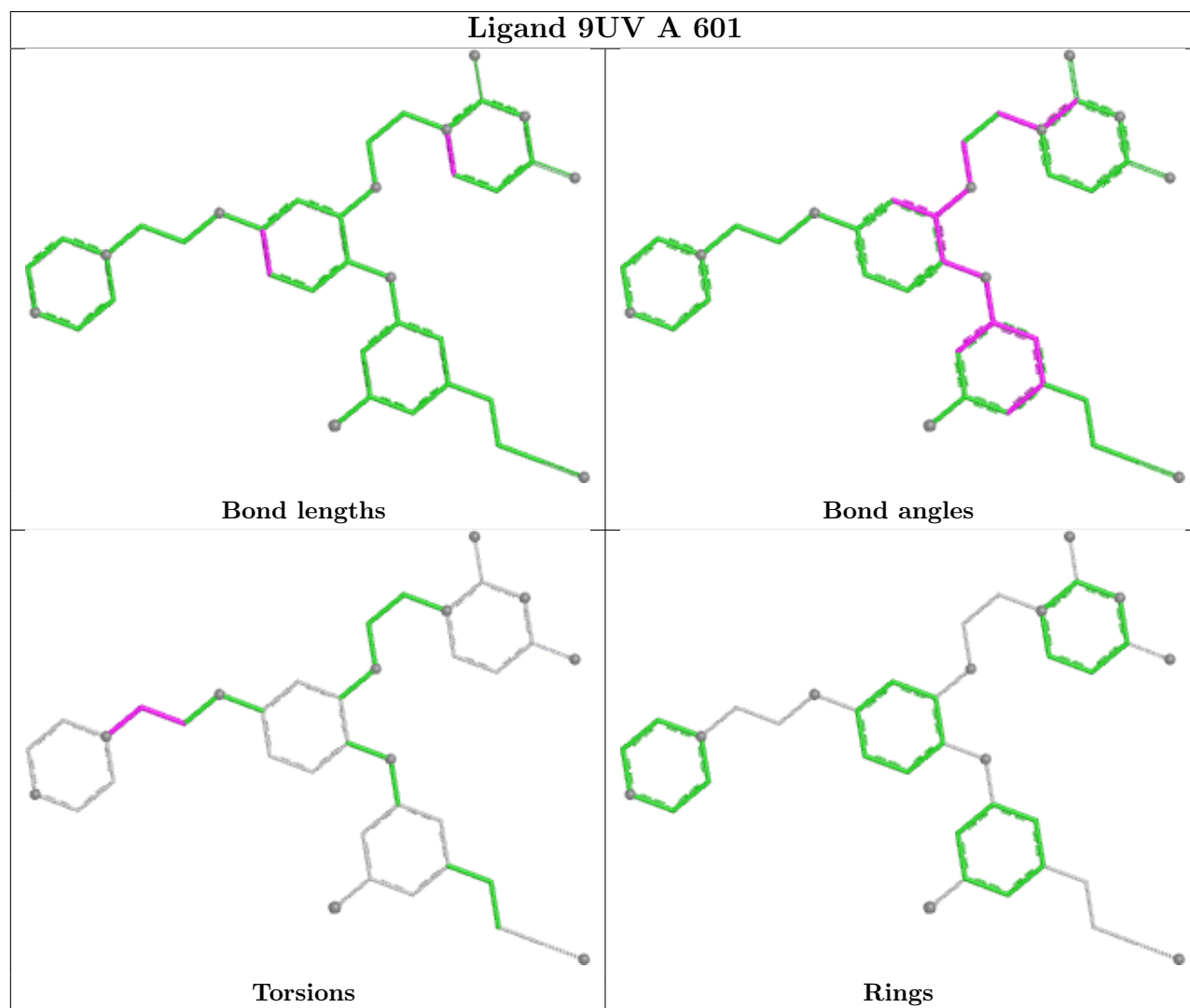
Mol	Chain	Res	Type	Atoms
3	A	601	9UV	C1A-C1B-N1E-C1I
3	A	601	9UV	O07-C1A-C1B-N1E

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	549/557 (98%)	-0.04	13 (2%) 59 50	8, 55, 130, 232	1 (0%)
2	B	428/428 (100%)	-0.14	14 (3%) 49 40	8, 47, 158, 291	1 (0%)
All	All	977/985 (99%)	-0.09	27 (2%) 55 46	8, 52, 144, 291	2 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	423	VAL	9.3
1	A	2	ILE	5.1
2	B	420	PRO	4.9
1	A	543	GLY	4.7
2	B	419	THR	4.0
2	B	67	ASP	4.0
1	A	221	HIS	3.9
1	A	1	PRO	3.8
2	B	299	ALA	3.6
2	B	5	ILE	3.3
2	B	2	ILE	3.1
1	A	474	ASN	2.9
1	A	55	PRO	2.8
2	B	241	VAL	2.8
1	A	24	TRP	2.7
1	A	133	PRO	2.7
1	A	252	TRP	2.6
1	A	545	ASN	2.5
2	B	14	PRO	2.4
2	B	356	ARG	2.4
2	B	354	TYR	2.3
1	A	227	PHE	2.3
2	B	1	PRO	2.1
2	B	24	TRP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	23	GLN	2.0
1	A	411	ILE	2.0
2	B	232	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](#)

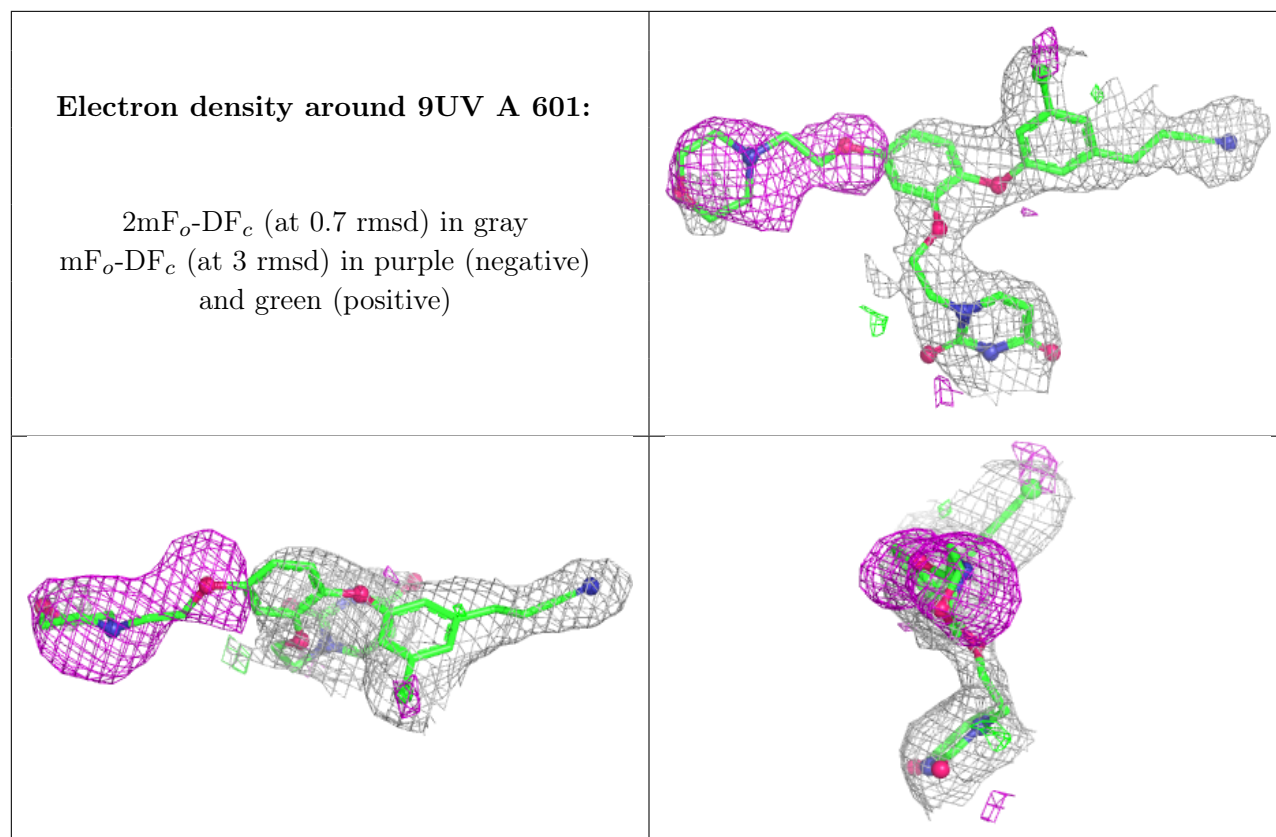
There are no oligosaccharides in this entry.

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	9UV	A	601	38/38	0.80	0.13	53,53,53,53	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.