



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:42 AM UTC

PDB ID : 8U6R / pdb_00008u6r
Title : Crystal Structure of HIV-1 Reverse Transcriptase in Complex with 3-(2-((2-cyanoindolizin-8-yl)oxy)phenoxy)-N-(2,2-difluoroethyl)propanamide (JLJ756), a non-nucleoside inhibitor
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Deposited on : 2023-09-13
Resolution : 2.87 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

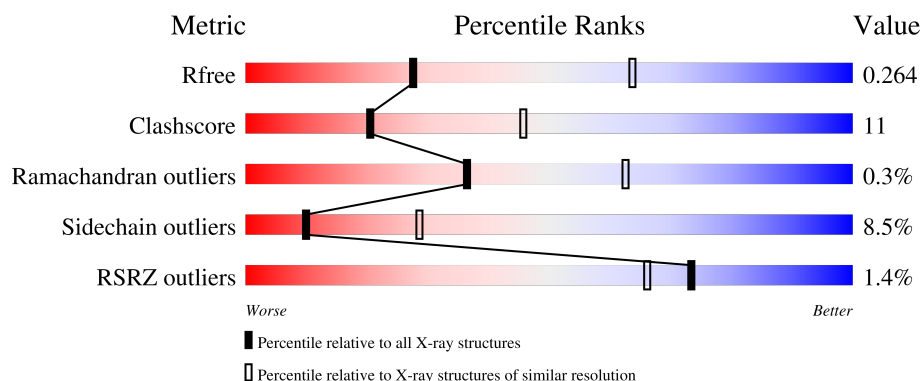
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.87 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	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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	554	2% 73% 23% ..
2	B	428	67% 25% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7455 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	546	Total	C	N	O	S	0	0	0
			4204	2723	688	786	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

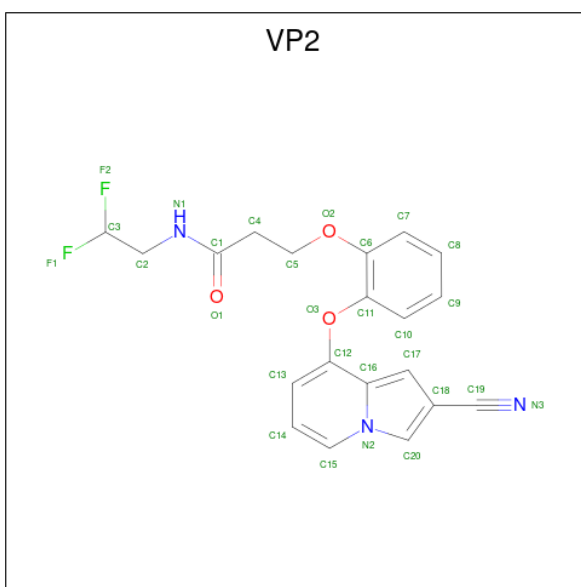
- Molecule 2 is a protein called p51 RT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	407	Total	C	N	O	S	0	0	0
			3203	2081	527	589	6			

There is a discrepancy between the modelled and reference sequences:

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

- Molecule 3 is 3-(2-{[(4R)-2-cyanoindolizin-8-yl]oxy}phenoxy)-N-(2,2-difluoroethyl)propanamide (CCD ID: VP2) (formula: C₂₀H₁₇F₂N₃O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			28	20	2	3	3		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	O P	0	0
			5	4 1		

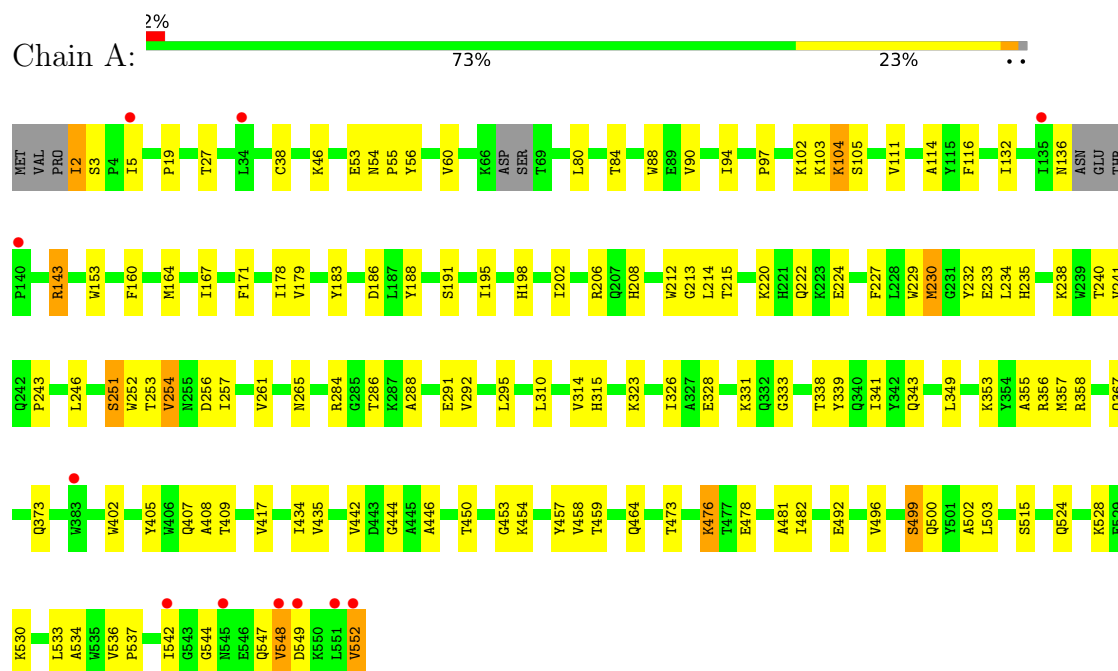
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	9	Total 9	O 9	0	0
5	B	6	Total 6	O 6	0	0

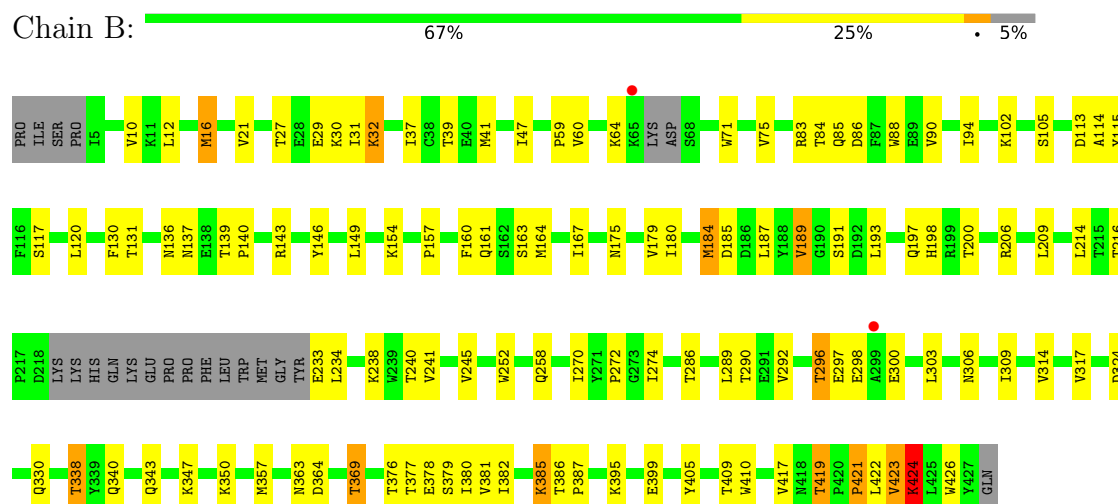
3 Residue-property plots [i](#)

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 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	224.92Å 69.98Å 106.31Å 90.00° 105.53° 90.00°	Depositor
Resolution (Å)	43.48 – 2.87 43.48 – 2.87	Depositor EDS
% Data completeness (in resolution range)	98.7 (43.48-2.87) 98.7 (43.48-2.87)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.17 (at 2.86Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487)	Depositor
R, R_{free}	0.215 , 0.261 0.228 , 0.264	Depositor DCC
R_{free} test set	1828 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	106.0	Xtriage
Anisotropy	0.234	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 87.3	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	0.95	EDS
Total number of atoms	7455	wwPDB-VP
Average B, all atoms (Å ²)	116.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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: PO4, VP2

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	0.45	0/4316	0.63	0/5911
2	B	0.45	0/3296	0.65	3/4507 (0.1%)
All	All	0.45	0/7612	0.64	3/10418 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	131	THR	CA-C-N	-9.90	115.46	122.59
2	B	131	THR	C-N-CA	-9.90	115.46	122.59
2	B	140	PRO	N-CA-C	-5.39	102.64	111.15

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	4204	0	4017	94	0
2	B	3203	0	3076	80	0
3	A	28	0	0	1	0
4	B	5	0	0	0	0
5	A	9	0	0	1	0
5	B	6	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7455	0	7093	163	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 11.

All (163) 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:5:ILE:HD11	1:A:167:ILE:HD11	1.50	0.93
1:A:458:VAL:HG22	1:A:464:GLN:HG2	1.58	0.84
1:A:84:THR:HG21	1:A:153:TRP:HE1	1.45	0.81
1:A:19:PRO:HG3	1:A:80:LEU:HB2	1.63	0.80
1:A:355:ALA:HB1	1:A:357:MET:HE3	1.64	0.80
2:B:423:VAL:HG23	2:B:426:TRP:HB2	1.63	0.79
2:B:240:THR:HG22	2:B:241:VAL:H	1.47	0.78
2:B:240:THR:HG22	2:B:241:VAL:N	1.99	0.77
2:B:85:GLN:HG3	2:B:154:LYS:HB3	1.66	0.76
1:A:3:SER:HB2	1:A:5:ILE:HD13	1.67	0.75
1:A:503:LEU:HD23	2:B:422:LEU:HD11	1.66	0.75
1:A:88:TRP:CE2	2:B:143:ARG:HD2	2.23	0.73
2:B:206:ARG:NH1	2:B:216:THR:OG1	2.22	0.72
2:B:143:ARG:HG2	2:B:143:ARG:HH11	1.55	0.71
1:A:343:GLN:HG3	1:A:349:LEU:HD11	1.72	0.70
2:B:180:ILE:HG12	2:B:189:VAL:HG13	1.72	0.69
2:B:47:ILE:HG22	2:B:146:TYR:HA	1.73	0.69
2:B:317:VAL:HG12	2:B:347:LYS:HB3	1.74	0.69
1:A:227:PHE:HB2	1:A:234:LEU:HB2	1.75	0.69
1:A:435:VAL:HG22	2:B:290:THR:HG21	1.74	0.69
2:B:357:MET:HG3	2:B:357:MET:O	1.93	0.67
1:A:503:LEU:CD2	2:B:422:LEU:HD11	2.24	0.66
2:B:64:LYS:HE2	2:B:71:TRP:CE2	2.29	0.66
2:B:85:GLN:HA	2:B:88:TRP:HB2	1.79	0.64
1:A:356:ARG:HB3	1:A:367:GLN:HE21	1.63	0.63
1:A:46:LYS:HD3	1:A:116:PHE:HB3	1.79	0.63
1:A:537:PRO:HG2	1:A:542:ILE:HD11	1.80	0.63
1:A:536:VAL:HB	1:A:542:ILE:HD13	1.81	0.63
1:A:191:SER:HB3	1:A:198:HIS:ND1	2.14	0.62
2:B:330:GLN:HE22	2:B:340:GLN:HE22	1.48	0.61
2:B:27:THR:HG22	2:B:29:GLU:H	1.65	0.61
1:A:356:ARG:HB3	1:A:367:GLN:NE2	2.16	0.60
1:A:53:GLU:H	1:A:53:GLU:CD	2.07	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:434:ILE:HD13	1:A:530:LYS:HB3	1.84	0.59
1:A:496:VAL:HG22	1:A:534:ALA:HB3	1.85	0.59
1:A:356:ARG:NE	1:A:358:ARG:O	2.35	0.59
1:A:3:SER:CB	1:A:5:ILE:HD13	2.32	0.58
1:A:478:GLU:HG2	1:A:499:SER:CB	2.34	0.58
2:B:240:THR:CG2	2:B:241:VAL:H	2.15	0.58
1:A:88:TRP:CZ2	2:B:143:ARG:HD2	2.40	0.57
1:A:202:ILE:O	1:A:206:ARG:HG3	2.04	0.57
1:A:53:GLU:O	1:A:54:ASN:C	2.47	0.57
1:A:454:LYS:HG3	1:A:552:VAL:HG22	1.87	0.57
2:B:379:SER:OG	2:B:387:PRO:HD3	2.05	0.56
1:A:103:LYS:O	1:A:105:SER:N	2.38	0.56
1:A:478:GLU:HG2	1:A:499:SER:OG	2.05	0.56
1:A:171:PHE:HB2	1:A:208:HIS:HD1	1.71	0.56
1:A:5:ILE:CD1	1:A:167:ILE:HD11	2.31	0.55
1:A:186:ASP:HB3	1:A:188:TYR:CE1	2.42	0.55
2:B:94:ILE:HG12	2:B:161:GLN:NE2	2.22	0.54
1:A:164:MET:HE1	1:A:214:LEU:HD13	1.89	0.54
1:A:55:PRO:HB2	1:A:143:ARG:NH1	2.23	0.54
1:A:2:ILE:HG23	1:A:213:GLY:HA3	1.90	0.54
1:A:252:TRP:CD1	1:A:295:LEU:HD11	2.43	0.53
2:B:16:MET:HE3	2:B:83:ARG:HA	1.90	0.53
1:A:235:HIS:HB2	1:A:238:LYS:O	2.08	0.53
2:B:120:LEU:HD13	2:B:149:LEU:HD23	1.90	0.53
1:A:5:ILE:HD11	1:A:167:ILE:CD1	2.31	0.53
2:B:60:VAL:HG12	2:B:75:VAL:HG22	1.92	0.52
2:B:143:ARG:HG2	2:B:143:ARG:NH1	2.22	0.52
2:B:330:GLN:NE2	2:B:338:THR:CB	2.72	0.52
2:B:423:VAL:HG22	2:B:424:LYS:N	2.24	0.52
2:B:350:LYS:NZ	2:B:378:GLU:OE1	2.44	0.51
2:B:114:ALA:HB2	2:B:214:LEU:HD13	1.92	0.51
1:A:5:ILE:HD12	1:A:212:TRP:HE3	1.77	0.50
1:A:252:TRP:HD1	1:A:295:LEU:HD11	1.75	0.50
1:A:254:VAL:HG11	1:A:288:ALA:O	2.11	0.50
1:A:284:ARG:O	1:A:286:THR:HG23	2.12	0.50
1:A:446:ALA:HA	1:A:453:GLY:HA2	1.94	0.50
2:B:369:THR:HG21	2:B:405:TYR:HB2	1.94	0.50
1:A:195:ILE:HD12	1:A:195:ILE:H	1.77	0.49
1:A:251:SER:HB2	1:A:292:VAL:HG11	1.94	0.49
2:B:191:SER:OG	2:B:198:HIS:ND1	2.29	0.49
1:A:235:HIS:ND1	1:A:238:LYS:HE3	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:12:LEU:HG	2:B:84:THR:HG22	1.94	0.49
1:A:80:LEU:O	1:A:84:THR:HG23	2.13	0.49
1:A:254:VAL:HG12	1:A:291:GLU:O	2.13	0.49
1:A:253:THR:HB	1:A:256:ASP:H	1.76	0.49
1:A:500:GLN:HG2	2:B:422:LEU:HD21	1.95	0.49
2:B:27:THR:O	2:B:31:ILE:HG13	2.12	0.48
2:B:330:GLN:NE2	2:B:338:THR:HB	2.28	0.48
2:B:160:PHE:HE2	2:B:164:MET:HE2	1.77	0.48
1:A:405:TYR:CE2	1:A:407:GLN:HB2	2.48	0.48
1:A:524:GLN:O	1:A:528:LYS:HG2	2.13	0.48
1:A:19:PRO:O	1:A:56:TYR:HB3	2.13	0.48
2:B:385:LYS:HE2	2:B:385:LYS:HB3	1.45	0.48
1:A:229:TRP:CE2	1:A:230:MET:HG2	2.49	0.47
2:B:330:GLN:NE2	2:B:338:THR:OG1	2.48	0.47
2:B:27:THR:HG22	2:B:29:GLU:N	2.29	0.47
1:A:492:GLU:HA	1:A:530:LYS:O	2.15	0.47
2:B:37:ILE:HD11	2:B:71:TRP:O	2.15	0.46
2:B:300:GLU:OE1	2:B:303:LEU:HD23	2.15	0.46
1:A:446:ALA:HA	1:A:453:GLY:CA	2.46	0.46
1:A:97:PRO:HG2	1:A:232:TYR:CD2	2.51	0.46
2:B:270:ILE:O	2:B:272:PRO:HD3	2.16	0.46
2:B:395:LYS:HE3	2:B:399:GLU:OE2	2.15	0.46
2:B:330:GLN:N	2:B:330:GLN:OE1	2.49	0.46
2:B:378:GLU:O	2:B:382:ILE:HD12	2.16	0.46
2:B:379:SER:CB	2:B:387:PRO:HD3	2.46	0.46
1:A:94:ILE:HG22	1:A:183:TYR:CE1	2.51	0.46
1:A:240:THR:HG22	1:A:241:VAL:N	2.31	0.46
2:B:21:VAL:HB	2:B:59:PRO:HD3	1.97	0.46
1:A:328:GLU:O	1:A:339:TYR:HA	2.16	0.45
2:B:240:THR:CG2	2:B:241:VAL:N	2.67	0.45
1:A:478:GLU:HG3	1:A:502:ALA:HB2	1.99	0.45
2:B:115:TYR:OH	2:B:157:PRO:HB3	2.17	0.45
2:B:324:ASP:O	2:B:343:GLN:HG2	2.17	0.45
1:A:444:GLY:HA2	1:A:552:VAL:HG21	1.97	0.45
1:A:496:VAL:HG21	2:B:289:LEU:HD21	1.99	0.45
1:A:442:VAL:HB	1:A:481:ALA:HB1	1.97	0.45
1:A:533:LEU:HD12	1:A:534:ALA:H	1.82	0.45
2:B:417:VAL:HG23	2:B:419:THR:HG23	1.98	0.45
2:B:193:LEU:HD13	2:B:197:GLN:HB3	1.98	0.44
2:B:274:ILE:HG23	2:B:306:ASN:OD1	2.18	0.44
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:160:PHE:CE2	2:B:164:MET:HE2	2.51	0.44
1:A:476:LYS:NZ	5:A:701:HOH:O	2.43	0.44
1:A:84:THR:HG21	1:A:153:TRP:NE1	2.22	0.44
1:A:103:LYS:O	1:A:104:LYS:C	2.61	0.44
1:A:233:GLU:OE2	1:A:243:PRO:HD3	2.18	0.44
1:A:240:THR:OG1	1:A:315:HIS:HA	2.18	0.44
1:A:341:ILE:O	1:A:349:LEU:N	2.45	0.43
2:B:41:MET:HB3	2:B:47:ILE:HG12	2.01	0.43
1:A:548:VAL:O	1:A:549:ASP:C	2.61	0.43
2:B:289:LEU:HD23	2:B:289:LEU:HA	1.88	0.43
1:A:408:ALA:HB1	2:B:364:ASP:HB3	2.01	0.43
2:B:184:MET:HE3	2:B:410:TRP:HB3	2.01	0.43
2:B:330:GLN:NE2	2:B:340:GLN:HE22	2.13	0.43
2:B:85:GLN:HG3	2:B:154:LYS:CB	2.43	0.43
1:A:224:GLU:N	1:A:224:GLU:OE1	2.52	0.43
2:B:209:LEU:HD22	2:B:214:LEU:HD23	2.01	0.43
1:A:252:TRP:HD1	1:A:295:LEU:CD1	2.31	0.42
1:A:323:LYS:HB2	1:A:343:GLN:NE2	2.34	0.42
2:B:113:ASP:O	2:B:114:ALA:C	2.61	0.42
1:A:257:ILE:O	1:A:261:VAL:HG22	2.18	0.42
1:A:38:CYS:SG	1:A:132:ILE:HD11	2.60	0.42
1:A:331:LYS:HG2	1:A:333:GLY:H	1.85	0.42
1:A:94:ILE:HG22	1:A:183:TYR:HE1	1.84	0.42
2:B:136:ASN:O	2:B:137:ASN:C	2.62	0.42
2:B:234:LEU:HD11	2:B:377:THR:HG21	2.01	0.42
1:A:102:LYS:HA	3:A:601:VP2:C4	2.50	0.42
2:B:252:TRP:O	2:B:292:VAL:HG13	2.19	0.42
1:A:88:TRP:CD2	2:B:143:ARG:HD2	2.54	0.42
1:A:195:ILE:HD12	1:A:195:ILE:N	2.34	0.42
1:A:246:LEU:HD11	1:A:310:LEU:HD12	2.01	0.42
2:B:32:LYS:HE2	2:B:32:LYS:HB3	1.90	0.42
1:A:478:GLU:HG2	1:A:499:SER:HB3	2.02	0.42
1:A:5:ILE:HD12	1:A:212:TRP:CE3	2.53	0.41
1:A:442:VAL:HG12	1:A:457:TYR:HB3	2.02	0.41
2:B:10:VAL:HG13	2:B:88:TRP:CZ2	2.55	0.41
1:A:547:GLN:O	1:A:548:VAL:C	2.63	0.41
2:B:417:VAL:O	2:B:417:VAL:HG13	2.21	0.41
1:A:238:LYS:HD2	1:A:315:HIS:CG	2.56	0.41
2:B:306:ASN:HA	2:B:309:ILE:HG22	2.03	0.41
2:B:163:SER:O	2:B:167:ILE:HG13	2.21	0.41
2:B:296:THR:OG1	2:B:297:GLU:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:376:THR:HG23	2:B:386:THR:HB	2.03	0.41
1:A:536:VAL:HG12	2:B:258:GLN:HB3	2.03	0.41
2:B:60:VAL:HG11	2:B:130:PHE:CD2	2.56	0.41
2:B:421:PRO:O	2:B:422:LEU:HB3	2.21	0.40
1:A:295:LEU:N	1:A:295:LEU:HD12	2.36	0.40
1:A:544:GLY:HA2	2:B:286:THR:HG22	2.03	0.40
2:B:30:LYS:HG2	2:B:71:TRP:CZ3	2.56	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	540/554 (98%)	500 (93%)	39 (7%)	1 (0%)	43	70
2	B	401/428 (94%)	378 (94%)	21 (5%)	2 (0%)	24	51
All	All	941/982 (96%)	878 (93%)	60 (6%)	3 (0%)	36	62

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	LYS
2	B	424	LYS
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	429/494 (87%)	396 (92%)	33 (8%)	12	33
2	B	332/390 (85%)	300 (90%)	32 (10%)	8	23
All	All	761/884 (86%)	696 (92%)	65 (8%)	10	28

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ILE
1	A	27	THR
1	A	60	VAL
1	A	90	VAL
1	A	111	VAL
1	A	136	ASN
1	A	143	ARG
1	A	178	ILE
1	A	179	VAL
1	A	215	THR
1	A	220	LYS
1	A	222	GLN
1	A	230	MET
1	A	251	SER
1	A	254	VAL
1	A	265	ASN
1	A	314	VAL
1	A	326	ILE
1	A	338	THR
1	A	353	LYS
1	A	373	GLN
1	A	402	TRP
1	A	409	THR
1	A	417	VAL
1	A	450	THR
1	A	459	THR
1	A	473	THR
1	A	476	LYS
1	A	482	ILE
1	A	499	SER
1	A	515	SER
1	A	548	VAL
1	A	552	VAL
2	B	16	MET

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Mol	Chain	Res	Type
2	B	32	LYS
2	B	39	THR
2	B	86	ASP
2	B	90	VAL
2	B	102	LYS
2	B	105	SER
2	B	117	SER
2	B	139	THR
2	B	175	ASN
2	B	179	VAL
2	B	184	MET
2	B	185	ASP
2	B	187	LEU
2	B	189	VAL
2	B	200	THR
2	B	233	GLU
2	B	238	LYS
2	B	245	VAL
2	B	296	THR
2	B	298	GLU
2	B	314	VAL
2	B	338	THR
2	B	363	ASN
2	B	369	THR
2	B	380	ILE
2	B	381	VAL
2	B	385	LYS
2	B	409	THR
2	B	419	THR
2	B	423	VAL
2	B	424	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	136	ASN
1	A	197	GLN
1	A	278	GLN
1	A	474	ASN
1	A	487	GLN
1	A	519	ASN
1	A	545	ASN

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Mol	Chain	Res	Type
2	B	255	ASN
2	B	269	GLN
2	B	330	GLN
2	B	332	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](#)

2 ligands are 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	VP2	A	601	-	29,30,30	2.74	7 (24%)	32,40,40	2.25	7 (21%)
4	PO4	B	501	-	4,4,4	0.87	0	6,6,6	0.75	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	VP2	A	601	-	-	3/15/17/17	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	VP2	C16-N2	-8.05	1.31	1.40
3	A	601	VP2	C19-C18	7.56	1.54	1.44
3	A	601	VP2	C20-N2	-5.57	1.28	1.37
3	A	601	VP2	C1-N1	5.37	1.46	1.33
3	A	601	VP2	C17-C18	3.23	1.45	1.40
3	A	601	VP2	O1-C1	-2.16	1.19	1.23
3	A	601	VP2	C14-C13	-2.13	1.35	1.41

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	VP2	F1-C3-C2	-7.95	98.79	110.18
3	A	601	VP2	C15-N2-C20	-4.80	121.75	130.31
3	A	601	VP2	F2-C3-C2	-4.16	104.22	110.18
3	A	601	VP2	C2-N1-C1	3.54	130.30	122.69
3	A	601	VP2	C17-C18-C19	3.20	130.86	125.61
3	A	601	VP2	O2-C5-C4	2.35	112.83	107.28
3	A	601	VP2	C17-C16-C12	2.13	137.15	131.70

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	601	VP2	O1-C1-C4-C5
3	A	601	VP2	C3-C2-N1-C1
3	A	601	VP2	C6-C11-O3-C12

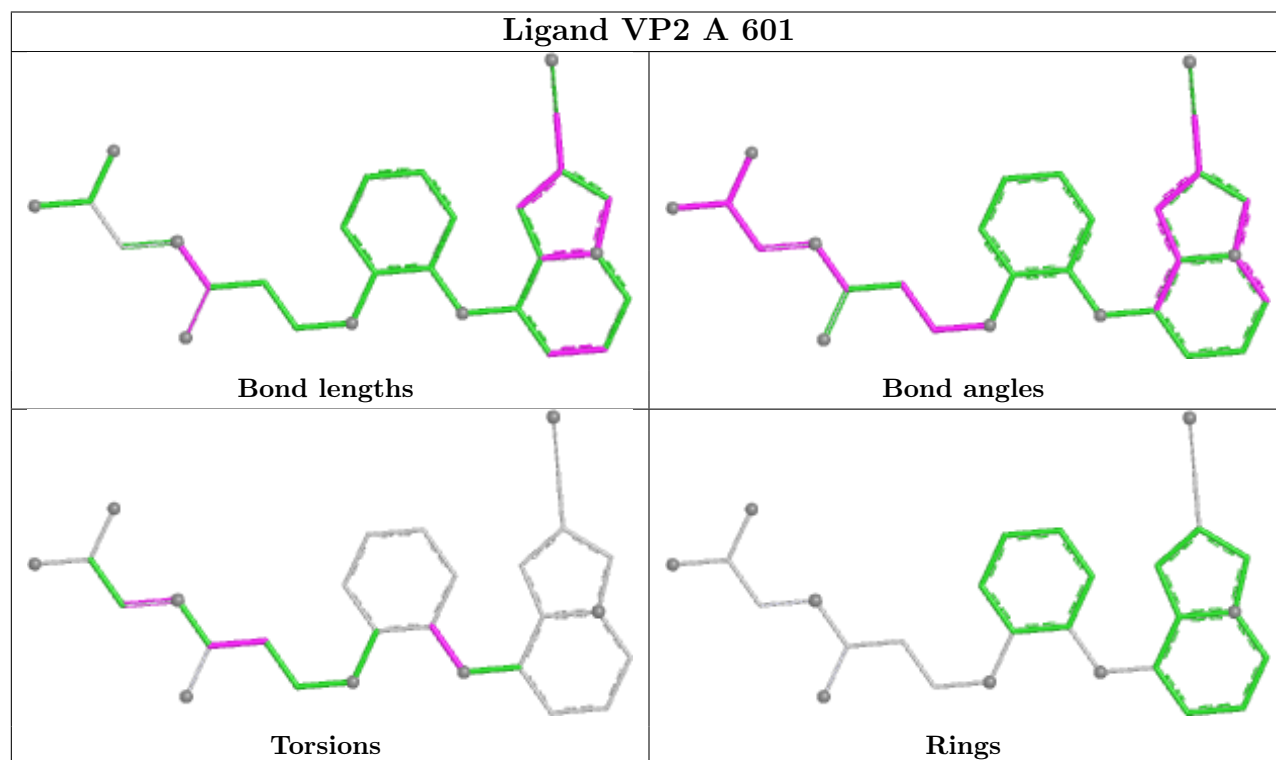
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	VP2	1	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	546/554 (98%)	-0.05	11 (2%) 65 57	75, 115, 166, 198	0
2	B	407/428 (95%)	-0.11	2 (0%) 87 83	80, 110, 163, 188	0
All	All	953/982 (97%)	-0.08	13 (1%) 73 66	75, 113, 164, 198	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	551	LEU	3.6
1	A	5	ILE	3.0
1	A	545	ASN	2.8
2	B	65	LYS	2.7
1	A	549	ASP	2.5
1	A	552	VAL	2.5
1	A	383	TRP	2.5
1	A	548	VAL	2.3
1	A	542	ILE	2.3
1	A	34	LEU	2.2
2	B	299	ALA	2.2
1	A	140	PRO	2.1
1	A	135	ILE	2.1

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](#)

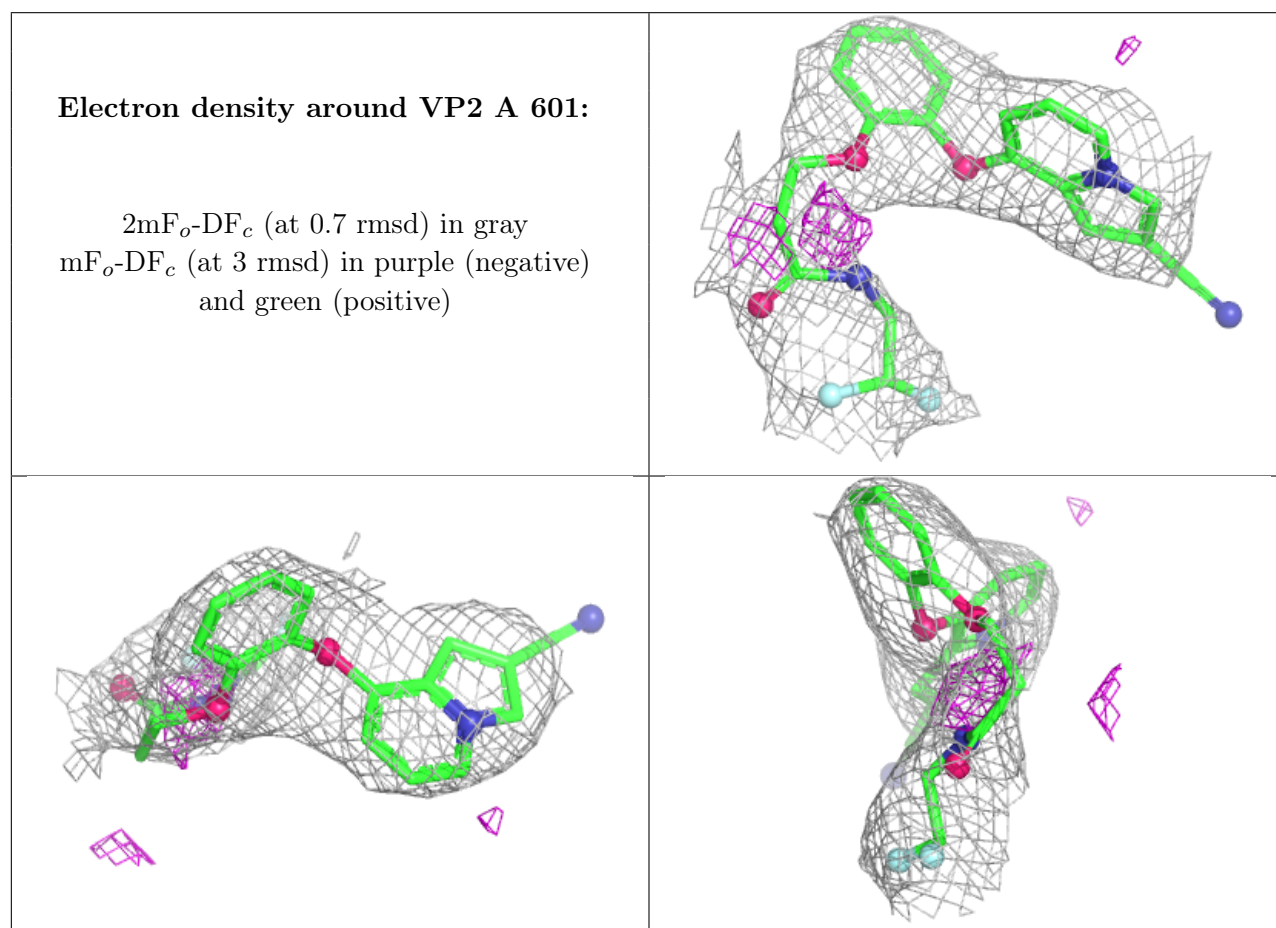
There are no oligosaccharides in this entry.

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(Å ²)	Q<0.9
4	PO4	B	501	5/5	0.70	0.08	132,132,158,186	0
3	VP2	A	601	28/28	0.93	0.10	86,100,107,111	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.