



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 12:53 PM UTC

PDB ID : 4I2Q / pdb_00004i2q
Title : Crystal structure of K103N/Y181C mutant of HIV-1 reverse transcriptase in complex with rilpivirine (TMC278) analogue
Authors : Patel, D.; Bauman, J.D.; Das, K.; Arnold, E.
Deposited on : 2012-11-22
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

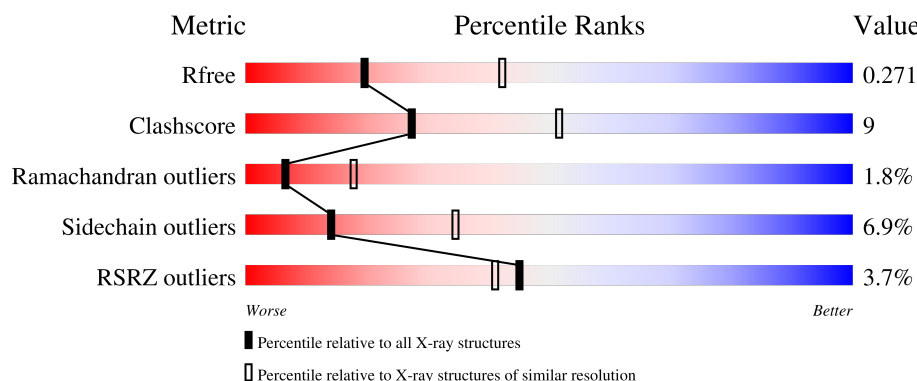
1 Overall quality at a glance

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	557	3% 72% 24% .
2	B	428	5% 69% 25% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8150 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 Gag-Pol polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	556	Total	C	N	O	S	0	2	0
			4522	2923	754	836	9			

There are 7 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	103	ASN	LYS	engineered mutation	UNP P03366
A	172	ALA	LYS	engineered mutation	UNP P03366
A	173	ALA	LYS	engineered mutation	UNP P03366
A	181	CYS	TYR	engineered mutation	UNP P03366
A	280	SER	CYS	engineered mutation	UNP P03366

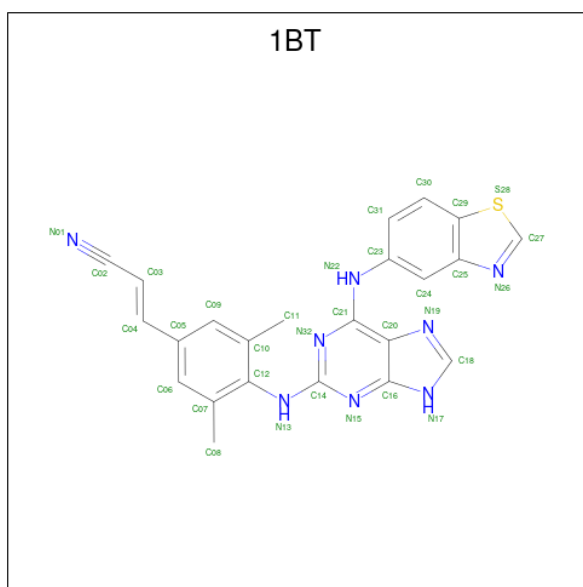
- Molecule 2 is a protein called Gag-Pol polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	416	Total	C	N	O	S	0	3	0
			3458	2254	571	626	7			

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 (2E)-3-(4-{[6-(1,3-benzothiazol-5-ylamino)-9H-purin-2-yl]amino}-3,5-dimethyl phenyl)prop-2-enitrile (CCD ID: 1BT) (formula: C₂₃H₁₈N₈S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	S	0	0
			32	23	8	1		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		

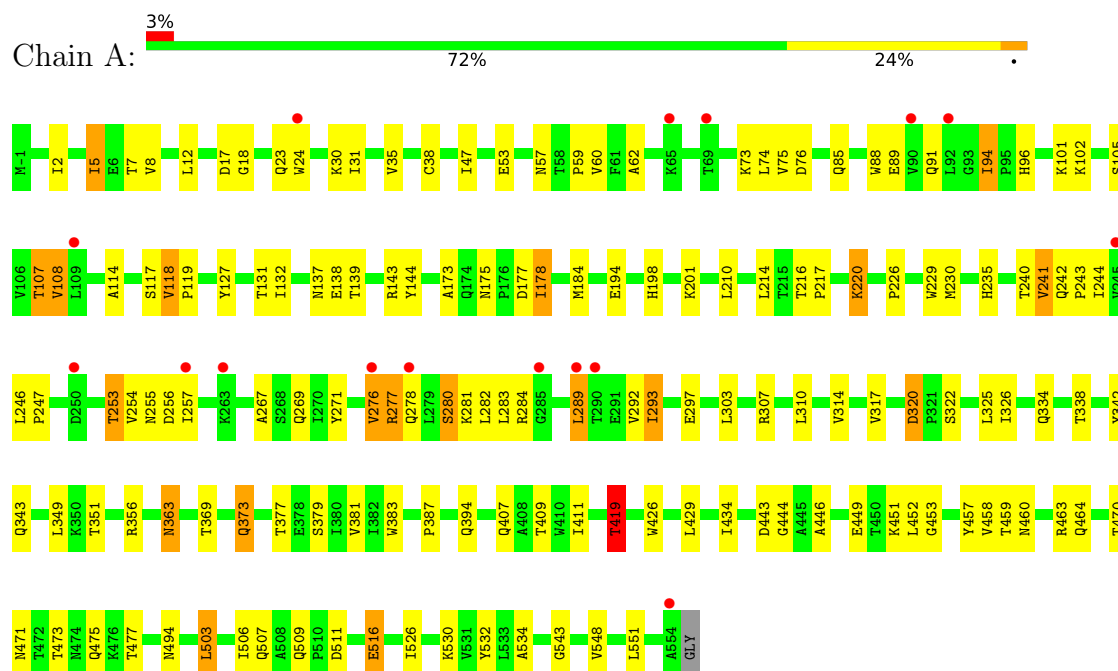
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	82	Total 82	O 82	0	0
5	B	52	Total 52	O 52	0	0

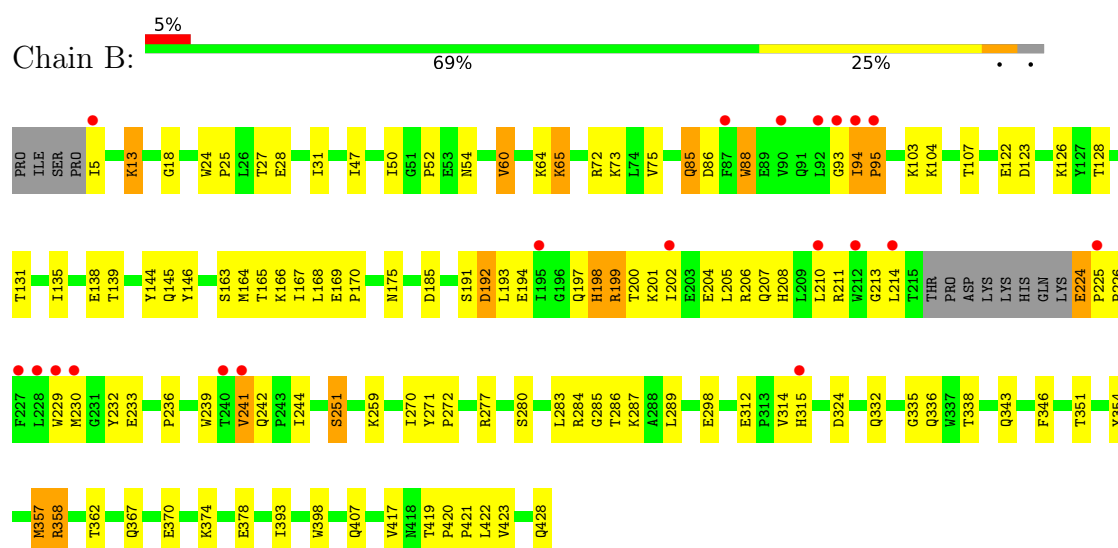
3 Residue-property plots

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: Gag-Pol polyprotein



• Molecule 2: Gag-Pol polyprotein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	162.60Å 72.95Å 108.54Å 90.00° 100.70° 90.00°	Depositor
Resolution (Å)	43.02 – 2.70 43.02 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.0 (43.02-2.70) 93.9 (43.02-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.211 , 0.277 0.203 , 0.271	Depositor DCC
R_{free} test set	2000 reflections (5.84%)	wwPDB-VP
Wilson B-factor (Å ²)	49.4	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8150	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 5.26% 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: 1BT, EDO

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	0.55	0/4646	0.95	14/6315 (0.2%)
2	B	0.58	0/3565	0.94	15/4844 (0.3%)
All	All	0.56	0/8211	0.94	29/11159 (0.3%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	118	VAL	CA-C-N	8.33	128.26	119.85
1	A	118	VAL	C-N-CA	8.33	128.26	119.85
2	B	198	HIS	N-CA-C	-8.08	102.16	110.97
1	A	320	ASP	CA-C-N	7.15	126.82	119.82
1	A	320	ASP	C-N-CA	7.15	126.82	119.82
1	A	242	GLN	CA-C-N	6.77	128.31	119.84
1	A	242	GLN	C-N-CA	6.77	128.31	119.84
2	B	139	THR	CA-C-N	-6.65	113.79	120.31
2	B	139	THR	C-N-CA	-6.65	113.79	120.31
2	B	86	ASP	N-CA-C	-6.32	101.17	110.52
2	B	13	LYS	CA-C-N	6.10	126.46	119.93
2	B	13	LYS	C-N-CA	6.10	126.46	119.93
2	B	18	GLY	CA-C-N	5.93	125.69	119.76
2	B	18	GLY	C-N-CA	5.93	125.69	119.76
2	B	312	GLU	CA-C-N	-5.73	114.66	120.85
2	B	312	GLU	C-N-CA	-5.73	114.66	120.85
1	A	419	THR	CA-C-N	5.58	126.12	120.38
1	A	419	THR	C-N-CA	5.58	126.12	120.38
1	A	216	THR	CA-C-N	5.55	126.77	119.84
1	A	216	THR	C-N-CA	5.55	126.77	119.84
1	A	96	HIS	CA-C-N	5.40	126.59	119.84
1	A	96	HIS	C-N-CA	5.40	126.59	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	224	GLU	CA-C-N	5.31	125.85	120.38
2	B	224	GLU	C-N-CA	5.31	125.85	120.38
2	B	24	TRP	CA-C-N	5.31	125.31	119.90
2	B	24	TRP	C-N-CA	5.31	125.31	119.90
2	B	88	TRP	CB-CA-C	-5.20	103.06	111.18
1	A	8	VAL	N-CA-C	5.19	112.69	107.55
1	A	363	ASN	N-CA-C	5.19	116.95	109.07

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	4522	0	4578	87	0
2	B	3458	0	3483	66	0
3	A	32	0	18	1	0
4	B	4	0	6	0	0
5	A	82	0	0	3	0
5	B	52	0	0	3	0
All	All	8150	0	8085	146	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 9.

All (146) 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)
2:B:357:MET:SD	5:B:601:HOH:O	2.26	0.92
1:A:178:ILE:HD11	1:A:201:LYS:HD3	1.62	0.81
1:A:276:VAL:HG12	1:A:277:ARG:H	1.49	0.78
1:A:23:GLN:HE22	1:A:60:VAL:HG12	1.50	0.77
1:A:282:LEU:HB3	1:A:293:ILE:HG21	1.71	0.73
1:A:343:GLN:HG3	1:A:349:LEU:HD11	1.74	0.69
1:A:253:THR:H	1:A:256:ASP:HB2	1.56	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:ASN:HB3	1:A:178:ILE:HD13	1.78	0.66
1:A:394:GLN:NE2	5:A:750:HOH:O	2.29	0.65
2:B:103:LYS:HB3	2:B:191:SER:O	2.02	0.60
1:A:255:ASN:HB2	1:A:289:LEU:HG	1.85	0.59
1:A:246:LEU:HD12	1:A:307:ARG:HG2	1.86	0.57
1:A:60:VAL:HG23	1:A:75:VAL:HG22	1.85	0.57
1:A:253:THR:HG23	1:A:255:ASN:H	1.70	0.57
1:A:226:PRO:HB3	1:A:235:HIS:NE2	2.19	0.56
1:A:247:PRO:O	1:A:307:ARG:NH2	2.35	0.56
2:B:242:GLN:HB2	2:B:351:THR:OG1	2.05	0.56
1:A:377:THR:O	1:A:381:VAL:HG23	2.06	0.56
1:A:23:GLN:NE2	1:A:131:THR:O	2.33	0.56
1:A:108:VAL:O	1:A:220:LYS:NZ	2.39	0.55
2:B:73:LYS:NZ	2:B:146:TYR:OH	2.39	0.55
2:B:50:ILE:HG21	2:B:145:GLN:HB3	1.88	0.55
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.42	0.55
1:A:31:ILE:O	1:A:35:VAL:HG23	2.07	0.55
1:A:107:THR:HG1	1:A:198:HIS:HE2	1.51	0.55
2:B:393:ILE:HD13	2:B:398:TRP:HB2	1.88	0.55
1:A:543:GLY:HA3	2:B:285:GLY:H	1.71	0.55
1:A:253:THR:HG22	1:A:256:ASP:H	1.72	0.55
1:A:107:THR:OG1	1:A:198:HIS:NE2	2.25	0.54
1:A:277:ARG:O	1:A:281:LYS:N	2.40	0.54
3:A:601:1BT:N32	3:A:601:1BT:H16	2.23	0.53
2:B:198:HIS:C	2:B:200:THR:H	2.17	0.53
1:A:516:GLU:N	1:A:516:GLU:OE1	2.42	0.53
2:B:13:LYS:HG3	2:B:85:GLN:HA	1.90	0.52
2:B:354:TYR:OH	2:B:378:GLU:OE2	2.25	0.52
1:A:449:GLU:OE1	1:A:449:GLU:N	2.41	0.52
1:A:379:SER:CB	1:A:387:PRO:HD3	2.39	0.52
2:B:175:ASN:HD21	2:B:201:LYS:HE3	1.75	0.52
1:A:278:GLN:NE2	1:A:334:GLN:HG2	2.24	0.52
2:B:200:THR:O	2:B:204:GLU:HG2	2.09	0.52
1:A:276:VAL:HG12	1:A:280:SER:OG	2.09	0.52
2:B:65:LYS:HD2	2:B:72:ARG:HD2	1.90	0.52
1:A:443:ASP:OD1	1:A:444:GLY:N	2.41	0.51
1:A:117:SER:O	1:A:119:PRO:HD3	2.11	0.51
1:A:363:ASN:ND2	1:A:509:GLN:O	2.33	0.51
1:A:419:THR:O	1:A:419:THR:OG1	2.25	0.50
1:A:381:VAL:HG22	2:B:25:PRO:HG3	1.92	0.50
1:A:244:ILE:HG13	1:A:267:ALA:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:336:GLN:HA	2:B:354:TYR:O	2.11	0.50
1:A:434:ILE:HD13	1:A:530:LYS:HB3	1.94	0.49
2:B:324:ASP:O	2:B:343:GLN:HG2	2.12	0.49
2:B:270:ILE:HG12	2:B:346:PHE:HB3	1.95	0.48
1:A:240:THR:HG23	1:A:241:VAL:O	2.13	0.48
1:A:460:ASN:HA	2:B:286:THR:O	2.14	0.48
2:B:298:GLU:OE1	2:B:298:GLU:N	2.36	0.48
2:B:206:ARG:HE	2:B:229:TRP:HA	1.77	0.48
1:A:426:TRP:HB3	1:A:526:ILE:HG12	1.96	0.48
1:A:257:ILE:CD1	1:A:283:LEU:HG	2.44	0.47
2:B:104:LYS:HD3	2:B:192:ASP:OD1	2.13	0.47
2:B:128:THR:OG1	2:B:146:TYR:HB2	2.13	0.47
1:A:446:ALA:HA	1:A:453:GLY:HA3	1.96	0.47
2:B:199:ARG:HA	2:B:202:ILE:CG1	2.45	0.47
1:A:342:TYR:HA	1:A:349:LEU:HD13	1.96	0.47
2:B:210:LEU:O	2:B:213:GLY:N	2.48	0.47
1:A:320:ASP:OD1	1:A:322:SER:OG	2.25	0.47
1:A:325:LEU:HD11	1:A:383:TRP:CD2	2.50	0.47
2:B:244:ILE:HD12	2:B:271:TYR:HE2	1.80	0.47
1:A:548:VAL:HA	1:A:551:LEU:HB2	1.97	0.46
1:A:57:ASN:OD1	1:A:143:ARG:NH1	2.48	0.46
1:A:532:TYR:CE2	1:A:534:ALA:HB2	2.50	0.46
1:A:107:THR:HG22	1:A:220:LYS:NZ	2.31	0.46
1:A:23:GLN:OE1	1:A:24:TRP:N	2.49	0.46
1:A:89:GLU:HB3	1:A:91:GLN:OE1	2.15	0.46
1:A:303:LEU:O	1:A:307:ARG:HG3	2.16	0.46
2:B:193:LEU:HB2	2:B:197:GLN:NE2	2.31	0.46
1:A:178:ILE:CD1	1:A:201:LYS:HD3	2.41	0.46
1:A:457:TYR:HE1	1:A:463:ARG:HG2	1.81	0.46
2:B:357:MET:HE1	2:B:374:LYS:HE2	1.96	0.46
1:A:173:ALA:O	5:A:753:HOH:O	2.21	0.46
2:B:166:LYS:HE3	2:B:166:LYS:HB2	1.68	0.46
1:A:5:ILE:HG13	1:A:117:SER:O	2.16	0.45
2:B:239:TRP:CZ2	2:B:378:GLU:HG2	2.51	0.45
1:A:543:GLY:N	2:B:283:LEU:O	2.49	0.45
2:B:27:THR:O	2:B:31:ILE:HG13	2.15	0.45
1:A:494:ASN:HB3	2:B:289:LEU:HD12	1.99	0.45
2:B:191:SER:HG	2:B:198:HIS:CE1	2.35	0.45
1:A:12:LEU:HG	1:A:127:TYR:CE2	2.51	0.45
2:B:47:ILE:HD12	2:B:144:TYR:CD1	2.52	0.45
2:B:207:GLN:O	2:B:211:ARG:N	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:GLN:HA	1:A:351:THR:O	2.18	0.44
1:A:473:THR:O	1:A:477:THR:HG23	2.17	0.44
2:B:107:THR:OG1	2:B:198:HIS:NE2	2.50	0.44
1:A:267:ALA:HB1	1:A:271:TYR:HD2	1.82	0.44
1:A:369:THR:O	1:A:373[B]:GLN:HG2	2.17	0.44
1:A:114:ALA:HA	1:A:214:LEU:HD22	1.99	0.44
1:A:503:LEU:HD12	1:A:507:GLN:CD	2.42	0.44
1:A:30:LYS:HD3	1:A:62:ALA:HB3	1.99	0.44
1:A:276:VAL:HG12	1:A:277:ARG:N	2.27	0.43
2:B:191:SER:HB2	2:B:193:LEU:HG	2.00	0.43
2:B:126:LYS:HE2	2:B:126:LYS:HB3	1.91	0.43
1:A:17:ASP:OD1	1:A:18:GLY:N	2.50	0.43
1:A:101:LYS:HD2	1:A:102:LYS:H	1.83	0.43
2:B:88:TRP:HZ3	2:B:93:GLY:HA2	1.84	0.43
1:A:47:ILE:HD12	1:A:144:TYR:CD1	2.54	0.43
2:B:60:VAL:HG23	2:B:75:VAL:HG22	2.01	0.43
1:A:38:CYS:SG	1:A:132:ILE:HD11	2.58	0.43
1:A:94:ILE:O	1:A:94:ILE:HG13	2.19	0.43
1:A:254:VAL:HG23	1:A:293:ILE:HD13	2.01	0.43
2:B:52:PRO:C	2:B:54:ASN:H	2.27	0.42
2:B:164:MET:O	2:B:168:LEU:HG	2.19	0.42
2:B:280:SER:O	2:B:283:LEU:N	2.51	0.42
1:A:407:GLN:NE2	2:B:417:VAL:O	2.51	0.42
1:A:278:GLN:HE22	1:A:334:GLN:HG2	1.84	0.42
1:A:88:TRP:HB2	2:B:54:ASN:O	2.20	0.42
2:B:107:THR:HA	2:B:232:TYR:O	2.20	0.42
1:A:429:LEU:HD11	1:A:506:ILE:HG22	2.02	0.42
2:B:259:LYS:NZ	5:B:611:HOH:O	2.40	0.42
2:B:335:GLY:HA2	2:B:367:GLN:OE1	2.20	0.42
2:B:208:HIS:HA	2:B:211:ARG:HB2	2.02	0.42
1:A:118:VAL:HA	1:A:119:PRO:HD3	1.76	0.42
2:B:239:TRP:CH2	2:B:378:GLU:HG2	2.55	0.42
2:B:358:ARG:HH22	2:B:370[B]:GLU:CD	2.28	0.42
2:B:332:GLN:HG3	2:B:338:THR:HG23	2.02	0.41
1:A:246:LEU:HG	1:A:310:LEU:HD12	2.03	0.41
2:B:94:ILE:HA	2:B:95:PRO:HD2	1.91	0.41
1:A:451:LYS:HB3	1:A:471:ASN:HA	2.03	0.41
1:A:59:PRO:HG2	1:A:76:ASP:HB3	2.01	0.41
2:B:419:THR:HA	2:B:420:PRO:HD2	1.88	0.41
2:B:64:LYS:O	2:B:407:GLN:HG2	2.20	0.41
1:A:543:GLY:HA3	2:B:284:ARG:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:205:LEU:HD23	2:B:205:LEU:O	2.21	0.41
1:A:452:LEU:HD23	1:A:470:THR:HG22	2.01	0.41
2:B:28:GLU:HB2	2:B:135:ILE:HD11	2.03	0.41
2:B:224:GLU:HA	2:B:225:PRO:HD3	1.79	0.41
2:B:420:PRO:HA	2:B:421:PRO:HD3	1.86	0.41
2:B:251:SER:HA	5:B:617:HOH:O	2.20	0.41
2:B:422:LEU:HD23	2:B:422:LEU:HA	1.93	0.41
1:A:229:TRP:CE2	1:A:230:MET:HG2	2.56	0.40
2:B:198:HIS:O	2:B:200:THR:N	2.52	0.40
1:A:281:LYS:HE3	1:A:284:ARG:CZ	2.51	0.40
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.21	0.40
2:B:13:LYS:HE3	2:B:85:GLN:HA	2.03	0.40
2:B:194:GLU:HG2	2:B:197:GLN:HE22	1.86	0.40
1:A:177:ASP:C	1:A:178:ILE:HD12	2.47	0.40
1:A:475:GLN:HG3	5:A:771:HOH:O	2.21	0.40
1:A:458:VAL:HG12	1:A:464:GLN:HG2	2.04	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	556/557 (100%)	519 (93%)	30 (5%)	7 (1%)	9	25
2	B	415/428 (97%)	370 (89%)	35 (8%)	10 (2%)	4	12
All	All	971/985 (99%)	889 (92%)	65 (7%)	17 (2%)	6	18

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	ARG
1	A	243	PRO

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Mol	Chain	Res	Type
2	B	85	GLN
1	A	85	GLN
2	B	122	GLU
2	B	138	GLU
2	B	170	PRO
2	B	199	ARG
1	A	516	GLU
2	B	272	PRO
2	B	167	ILE
2	B	95	PRO
2	B	226	PRO
1	A	217	PRO
1	A	276	VAL
2	B	241	VAL
1	A	317	VAL

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	497/495 (100%)	461 (93%)	36 (7%)	13	32
2	B	380/390 (97%)	354 (93%)	26 (7%)	14	35
All	All	877/885 (99%)	815 (93%)	62 (7%)	14	33

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

Mol	Chain	Res	Type
1	A	2	ILE
1	A	5	ILE
1	A	7	THR
1	A	53	GLU
1	A	73	LYS
1	A	74	LEU
1	A	94	ILE
1	A	105	SER

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Mol	Chain	Res	Type
1	A	107	THR
1	A	108	VAL
1	A	137	ASN
1	A	138	GLU
1	A	139	THR
1	A	178	ILE
1	A	184	MET
1	A	194	GLU
1	A	210	LEU
1	A	220	LYS
1	A	241	VAL
1	A	253	THR
1	A	280	SER
1	A	289	LEU
1	A	292	VAL
1	A	293	ILE
1	A	297	GLU
1	A	314	VAL
1	A	326	ILE
1	A	338	THR
1	A	356	ARG
1	A	373[A]	GLN
1	A	373[B]	GLN
1	A	409	THR
1	A	411	ILE
1	A	419	THR
1	A	459	THR
1	A	503	LEU
2	B	5	ILE
2	B	60	VAL
2	B	65	LYS
2	B	94	ILE
2	B	123	ASP
2	B	131	THR
2	B	163	SER
2	B	165	THR
2	B	169	GLU
2	B	185	ASP
2	B	192	ASP
2	B	214	LEU
2	B	230	MET
2	B	233	GLU

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Mol	Chain	Res	Type
2	B	241	VAL
2	B	251	SER
2	B	277	ARG
2	B	287	LYS
2	B	314	VAL
2	B	315[A]	HIS
2	B	315[B]	HIS
2	B	357	MET
2	B	358	ARG
2	B	362	THR
2	B	423	VAL
2	B	428	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	197	GLN
1	A	474	ASN
1	A	475	GLN
2	B	330	GLN
2	B	332	GLN
2	B	340	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	1BT	A	601	-	36,36,36	2.00	9 (25%)	47,51,51	2.65	13 (27%)
4	EDO	B	501	-	3,3,3	0.48	0	2,2,2	0.43	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	1BT	A	601	-	-	2/11/12/12	0/5/5/5
4	EDO	B	501	-	-	1/1/1/1	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	1BT	C27-N26	7.46	1.39	1.29
3	A	601	1BT	C21-N22	4.16	1.44	1.36
3	A	601	1BT	C25-N26	3.97	1.47	1.39
3	A	601	1BT	C27-S28	-3.69	1.68	1.73
3	A	601	1BT	C14-N13	3.03	1.42	1.36
3	A	601	1BT	C29-S28	-2.21	1.67	1.73
3	A	601	1BT	C16-N15	2.10	1.38	1.34
3	A	601	1BT	C24-C25	2.06	1.43	1.40
3	A	601	1BT	C18-N19	2.01	1.36	1.32

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	1BT	S28-C27-N26	-12.30	109.29	117.01
3	A	601	1BT	C20-C16-N15	-5.97	120.69	126.55
3	A	601	1BT	C20-N19-C18	4.54	108.60	103.48
3	A	601	1BT	C30-C29-S28	4.30	134.92	127.36
3	A	601	1BT	N17-C18-N19	-3.50	108.53	113.87
3	A	601	1BT	C23-C24-C25	3.48	121.48	118.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	1BT	C24-C25-N26	-3.46	124.70	130.51
3	A	601	1BT	C16-C20-C21	2.74	118.61	116.95
3	A	601	1BT	C16-C20-N19	-2.74	106.87	110.67
3	A	601	1BT	N17-C16-N15	2.41	131.32	127.45
3	A	601	1BT	N32-C14-N15	-2.20	122.66	126.26
3	A	601	1BT	C14-N32-C21	2.13	121.59	116.36
3	A	601	1BT	C11-C10-C12	2.07	123.94	121.42

There are no chirality outliers.

All (3) torsion outliers are listed below:

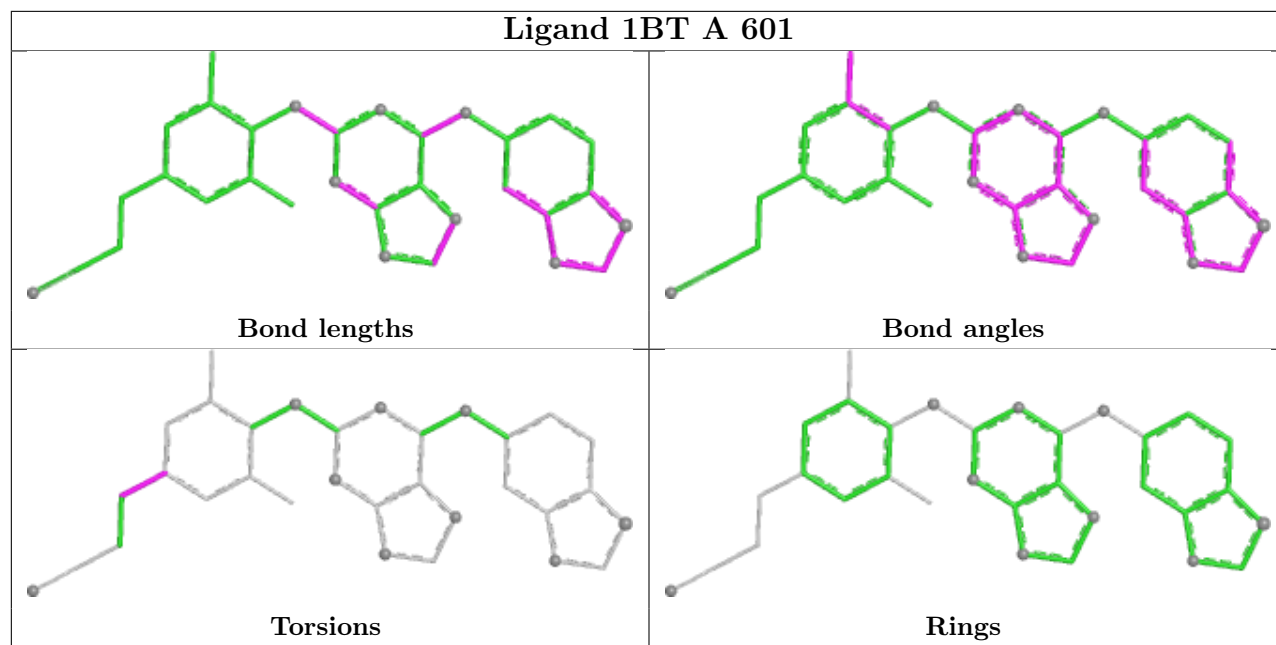
Mol	Chain	Res	Type	Atoms
3	A	601	1BT	C03-C04-C05-C09
3	A	601	1BT	C03-C04-C05-C06
4	B	501	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	1BT	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

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	556/557 (99%)	0.12	16 (2%)	53 50	21, 59, 119, 152	2 (0%)
2	B	416/428 (97%)	0.10	20 (4%)	35 32	20, 53, 108, 140	3 (0%)
All	All	972/985 (98%)	0.11	36 (3%)	45 41	20, 57, 114, 152	5 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	92	LEU	5.6
2	B	94	ILE	4.3
2	B	95	PRO	4.0
1	A	92	LEU	4.0
1	A	250	ASP	3.4
1	A	263	LYS	3.4
1	A	69	THR	3.3
1	A	554	ALA	3.2
2	B	315[A]	HIS	3.1
2	B	90	VAL	3.0
2	B	5	ILE	3.0
2	B	93	GLY	2.9
2	B	230	MET	2.9
2	B	228	LEU	2.7
2	B	240	THR	2.7
2	B	202	ILE	2.7
1	A	245	VAL	2.7
2	B	225	PRO	2.5
1	A	65	LYS	2.5
1	A	109	LEU	2.4
1	A	289	LEU	2.4
2	B	229	TRP	2.3
2	B	195	ILE	2.3
1	A	90	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	210	LEU	2.2
1	A	290	THR	2.2
2	B	227	PHE	2.2
1	A	257	ILE	2.2
2	B	87	PHE	2.2
1	A	278	GLN	2.1
1	A	285	GLY	2.1
1	A	24	TRP	2.1
2	B	212	TRP	2.1
2	B	241	VAL	2.1
2	B	214	LEU	2.0
1	A	276	VAL	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](#)

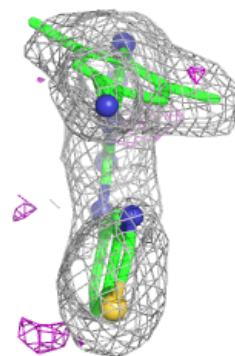
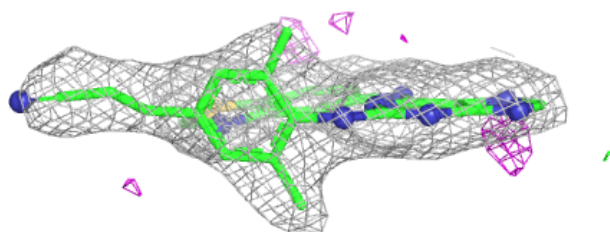
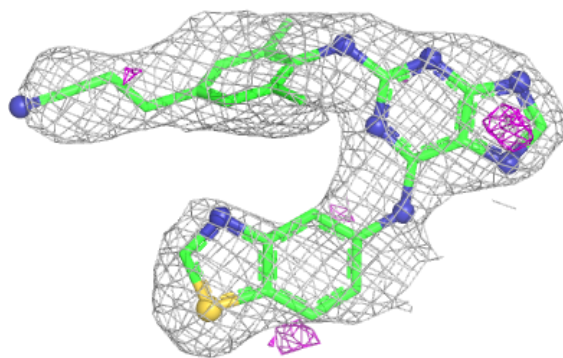
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	EDO	B	501	4/4	0.89	0.13	42,45,47,54	0
3	1BT	A	601	32/32	0.91	0.12	52,62,72,76	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 1BT A 601:

$2mF_o-DF_c$ (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.