



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 02:23 AM GMT

PDB ID : 4I2Q
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 validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

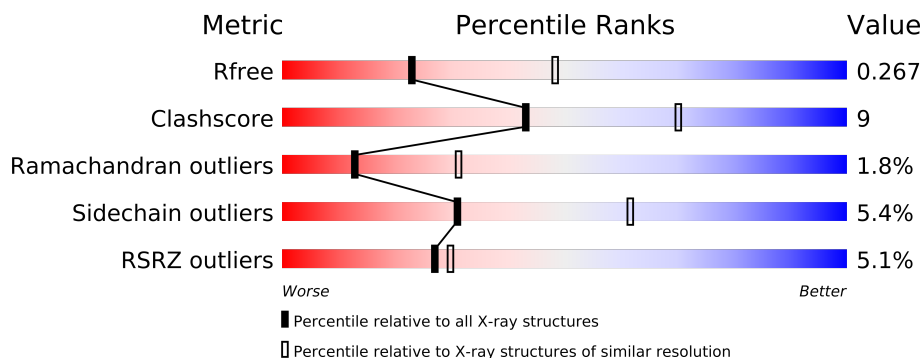
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	557	
2	B	428	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	EDO	B	501	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8150 atoms, of which 0 are hydrogen and 0 are deuterium.

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

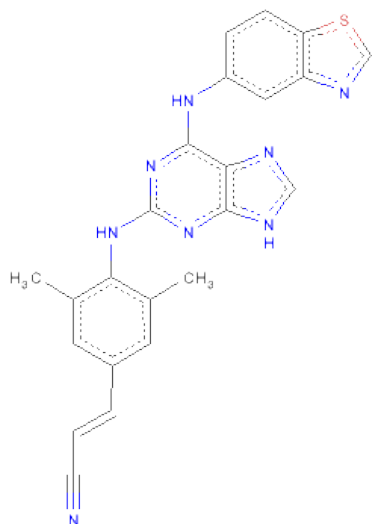
- Molecule 2 is a protein called Gag-Pol polyprotein.

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-BENZOTHAZOL-5-YLAMINO)-9H-PURIN-2-YL]AMINO}-3,5-DIMETHYLPHENYL)PROP-2-ENENITRILE (three-letter code: 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 (three-letter code: EDO) (formula: C₂H₆O₂).



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

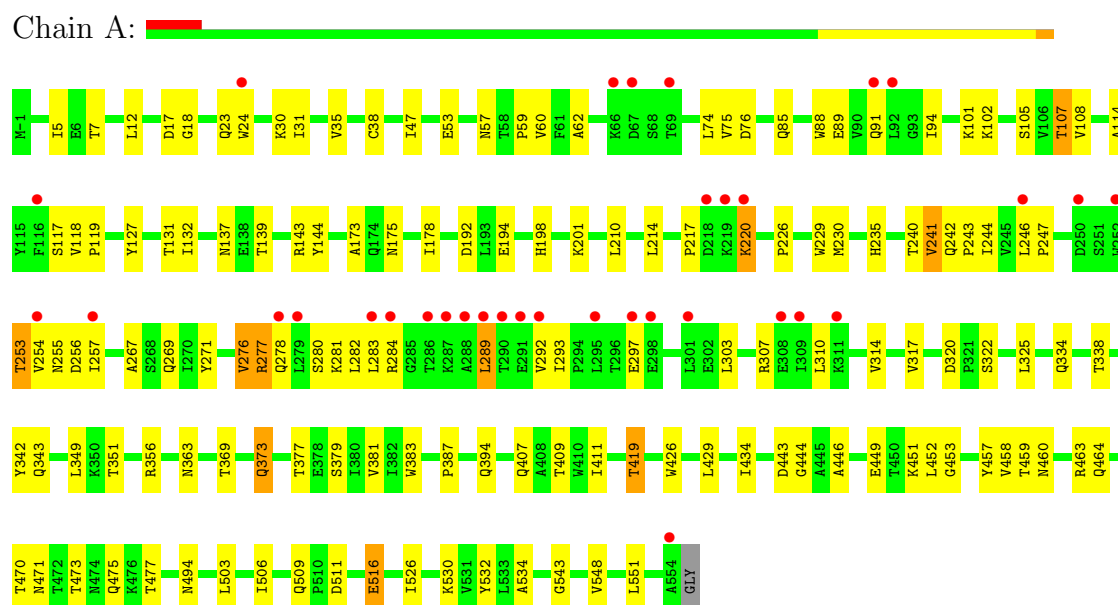
- Molecule 5 is water.

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

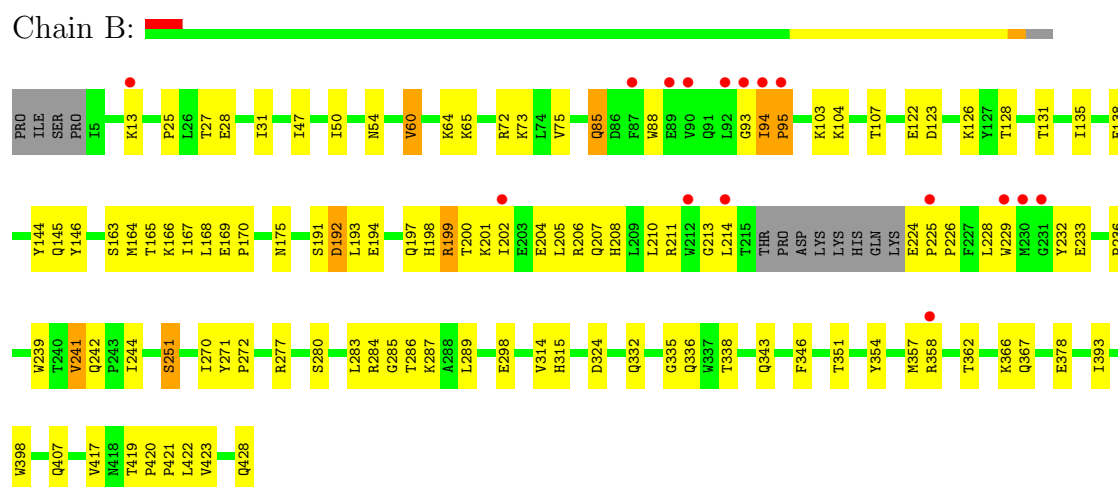
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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) 94.0 (43.02-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.211 , 0.277 0.201 , 0.267	Depositor DCC
R_{free} test set	1888 reflections (5.82%)	DCC
Wilson B-factor (Å ²)	49.5	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 31.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 34229 reflections	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 5.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.44	0/4646	0.60	0/6315
2	B	0.47	0/3565	0.62	0/4844
All	All	0.46	0/8211	0.61	0/11159

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4522	0	4578	85	0
2	B	3458	0	3483	63	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	141	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (141) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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
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
2:B:73:LYS:NZ	2:B:146:TYR:OH	2.39	0.55
1:A:108:VAL:O	1:A:220:LYS:NZ	2.39	0.55
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.42	0.55
2:B:50:ILE:HG21	2:B:145:GLN:HB3	1.88	0.55
1:A:31:ILE:O	1:A:35:VAL:HG23	2.07	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:H16	3:A:601:1BT:N32	2.23	0.53
1:A:107:THR:HG1	1:A:198:HIS:HE2	1.51	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
2:B:200:THR:O	2:B:204:GLU:HG2	2.09	0.52
1:A:278:GLN:NE2	1:A:334:GLN:HG2	2.24	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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:363:ASN:ND2	1:A:509:GLN:O	2.33	0.51
1:A:117:SER:O	1:A:119:PRO:HD3	2.11	0.51
1:A:419:THR:OG1	1:A:419:THR:O	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
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
2:B:198:HIS:C	2:B:200:THR:H	2.17	0.48
1:A:460:ASN:HA	2:B:286:THR:O	2.14	0.48
2:B:298:GLU:N	2:B:298:GLU:OE1	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
2:B:104:LYS:HD3	2:B:192:ASP:OD1	2.13	0.47
1:A:257:ILE:CD1	1:A:283:LEU:HG	2.44	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:532:TYR:CE2	1:A:534:ALA:HB2	2.50	0.46
1:A:57:ASN:OD1	1:A:143:ARG:NH1	2.48	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
2:B:193:LEU:HB2	2:B:197:GLN:NE2	2.31	0.46
1:A:303:LEU:O	1:A:307:ARG:HG3	2.16	0.46
1:A:89:GLU:HB3	1:A:91:GLN:OE1	2.15	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
1:A:173:ALA:O	5:A:753:HOH:O	2.21	0.46
2:B:166:LYS:HB2	2:B:166:LYS:HE3	1.68	0.46
2:B:239:TRP:CZ2	2:B:378:GLU:HG2	2.51	0.45
1:A:5:ILE:HG13	1:A:117:SER:O	2.16	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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:207:GLN:O	2:B:211:ARG:N	2.47	0.45
2:B:47:ILE:HD12	2:B:144:TYR:CD1	2.52	0.45
1:A:269:GLN:HA	1:A:351:THR:O	2.18	0.44
2:B:366:LYS:NZ	5:B:612:HOH:O	2.49	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: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:HB3	2:B:126:LYS:HE2	1.91	0.43
2:B:88:TRP:HZ3	2:B:93:GLY:HA2	1.84	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:60:VAL:HG23	2:B:75:VAL:HG22	2.01	0.43
1:A:47:ILE:HD12	1:A:144:TYR:CD1	2.54	0.43
1:A:38:CYS:SG	1:A:132:ILE:HD11	2.58	0.43
1:A:254:VAL:HG23	1:A:293:ILE:HD13	2.01	0.43
1:A:94:ILE:O	1:A:94:ILE:HG13	2.19	0.43
2:B:280:SER:O	2:B:283:LEU:N	2.51	0.42
2:B:164:MET:O	2:B:168:LEU:HG	2.19	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
2:B:107:THR:HA	2:B:232:TYR:O	2.20	0.42
1:A:88:TRP:HB2	2:B:54:ASN:O	2.20	0.42
1:A:429:LEU:HD11	1:A:506:ILE:HG22	2.02	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
2:B:239:TRP:CH2	2:B:378:GLU:HG2	2.55	0.42
1:A:118:VAL:HA	1:A:119:PRO:HD3	1.76	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	Distance(Å)	Clash(Å)
2:B:205:LEU:HD23	2:B:205:LEU:O	2.21	0.41
2:B:420:PRO:HA	2:B:421:PRO:HD3	1.86	0.41
2:B:224:GLU:HA	2:B:225:PRO:HD3	1.79	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:251:SER:HA	5:B:617:HOH:O	2.20	0.41
2:B:422:LEU:HA	2:B:422:LEU:HD23	1.93	0.41
2:B:198:HIS:O	2:B:200:THR:N	2.52	0.40
1:A:229:TRP:CE2	1:A:230:MET:HG2	2.56	0.40
1:A:281:LYS:HE3	1:A:284:ARG:CZ	2.51	0.40
2:B:194:GLU:HG2	2:B:197:GLN:HE22	1.86	0.40
2:B:13:LYS:HE3	2:B:85:GLN:HA	2.03	0.40
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.21	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

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	556/557 (100%)	519 (93%)	30 (5%)	7 (1%)	18	43
2	B	415/428 (97%)	370 (89%)	35 (8%)	10 (2%)	9	22
All	All	971/985 (99%)	889 (92%)	65 (7%)	17 (2%)	13	31

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	ARG
1	A	243	PRO
2	B	85	GLN
1	A	85	GLN
2	B	122	GLU

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Mol	Chain	Res	Type
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 ⓘ

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%)	470 (95%)	27 (5%)	31	61
2	B	380/390 (97%)	358 (94%)	22 (6%)	28	57
All	All	877/885 (99%)	828 (94%)	49 (6%)	31	59

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	THR
1	A	53	GLU
1	A	74	LEU
1	A	105	SER
1	A	107	THR
1	A	137	ASN
1	A	139	THR
1	A	192	ASP
1	A	194	GLU
1	A	210	LEU
1	A	220	LYS
1	A	241	VAL
1	A	242	GLN

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Mol	Chain	Res	Type
1	A	253	THR
1	A	289	LEU
1	A	292	VAL
1	A	297	GLU
1	A	314	VAL
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	60	VAL
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	192	ASP
2	B	214	LEU
2	B	228	LEU
2	B	233	GLU
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	358	ARG
2	B	362	THR
2	B	423	VAL
2	B	428	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates 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	1.83	5 (13%)	50,51,51	2.25	16 (32%)
4	EDO	B	501	-	3,3,3	0.58	0	2,2,2	0.44	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	-	-	0/12/12/12	0/1/5/5
4	EDO	B	501	-	-	0/1/1/1	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	1BT	C29-S28	-6.64	1.67	1.74
3	A	601	1BT	C14-N13	4.14	1.42	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	1BT	C21-N22	4.07	1.44	1.36
3	A	601	1BT	C16-N15	3.43	1.38	1.34
3	A	601	1BT	C25-N26	2.63	1.47	1.38

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	1BT	N17-C16-N15	8.36	131.32	124.37
3	A	601	1BT	C30-C29-S28	5.30	134.92	125.67
3	A	601	1BT	C25-C29-S28	-4.02	106.00	112.09
3	A	601	1BT	C31-C30-C29	3.69	121.63	118.81
3	A	601	1BT	C20-C16-N15	-3.62	120.69	125.94
3	A	601	1BT	C29-C25-N26	3.52	116.03	108.11
3	A	601	1BT	C16-C20-N19	-3.15	106.87	109.93
3	A	601	1BT	C05-C06-C07	-2.98	119.25	122.28
3	A	601	1BT	C14-N15-C16	2.80	118.66	115.15
3	A	601	1BT	C27-S28-C29	2.56	98.92	92.69
3	A	601	1BT	C20-C21-N32	-2.44	117.79	120.45
3	A	601	1BT	C10-C12-N13	2.32	121.88	119.31
3	A	601	1BT	N15-C14-N32	-2.29	122.66	126.19
3	A	601	1BT	C14-N32-C21	2.22	121.59	116.97
3	A	601	1BT	C11-C10-C12	2.20	123.94	121.44
3	A	601	1BT	C24-C25-N26	-2.14	124.70	130.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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%)	0.13	34 (6%) 21 22	25, 59, 119, 152	0
2	B	416/428 (97%)	0.14	16 (3%) 38 43	26, 53, 108, 140	0
All	All	972/985 (98%)	0.13	50 (5%) 27 30	25, 57, 114, 152	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	67	ASP	8.1
2	B	92	LEU	7.3
1	A	290	THR	6.2
1	A	286	THR	6.1
1	A	219	LYS	5.8
1	A	250	ASP	5.5
2	B	90	VAL	5.0
2	B	89	GLU	4.9
2	B	95	PRO	4.7
2	B	93	GLY	4.5
1	A	298	GLU	4.4
2	B	231	GLY	4.3
1	A	554	ALA	4.1
1	A	92	LEU	4.0
2	B	358	ARG	4.0
1	A	220	LYS	4.0
1	A	283	LEU	4.0
1	A	291	GLU	3.7
1	A	297	GLU	3.4
2	B	214	LEU	3.3
1	A	287	LYS	3.3
2	B	212	TRP	3.2
1	A	218	ASP	3.1
1	A	295	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
2	B	225	PRO	3.0
1	A	278	GLN	2.9
1	A	257	ILE	2.9
1	A	252	TRP	2.8
1	A	254	VAL	2.8
2	B	229	TRP	2.7
1	A	69	THR	2.6
1	A	308	GLU	2.6
2	B	94	ILE	2.5
1	A	292	VAL	2.5
2	B	230	MET	2.5
1	A	24	TRP	2.4
1	A	279	LEU	2.4
1	A	301	LEU	2.4
1	A	284	ARG	2.4
1	A	311	LYS	2.3
2	B	87	PHE	2.3
1	A	288	ALA	2.3
1	A	289	LEU	2.2
1	A	116	PHE	2.2
1	A	66	LYS	2.2
1	A	246	LEU	2.1
1	A	91	GLN	2.1
1	A	309	ILE	2.1
2	B	202	ILE	2.1
2	B	13	LYS	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	EDO	B	501	4/4	0.24	3.06	42,45,47,54	0
3	1BT	A	601	32/32	0.20	0.49	52,62,72,76	0

6.5 Other polymers ⓘ

There are no such residues in this entry.