



Full wwPDB X-ray Structure Validation Report

Oct 28, 2014 – 07:46 PM EDT

PDB ID : 4O4G
Title : Crystal Structure of HIV-1 Reverse Transcriptase in complex with 4-((4-(methylamino)-1,3,5-triazin-2-yl)amino)benzonitrile(JLJ527), a non-nucleoside inhibitor
Authors : Mislak, A.C.; Frey, K.M.
Deposited on : 2013-12-18
Resolution : 2.71 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>

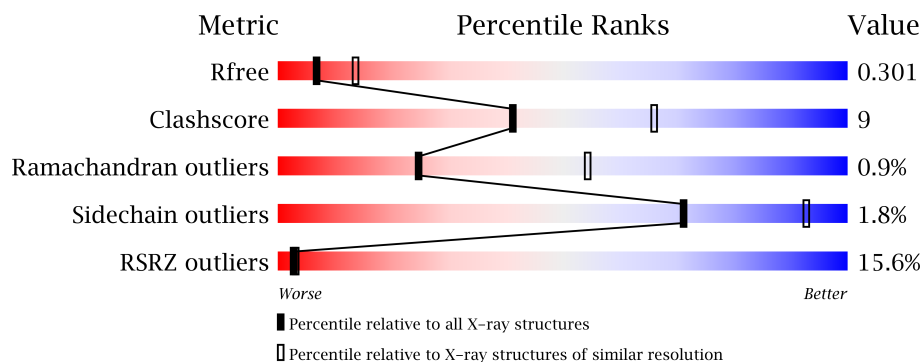
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable24103
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.1.3
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable24103

1 Overall quality at a glance



The reported resolution of this entry is 2.71 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	1770 (2.74-2.70)
Clashscore	79885	2183 (2.74-2.70)
Ramachandran outliers	78287	2147 (2.74-2.70)
Sidechain outliers	78261	2148 (2.74-2.70)
RSRZ outliers	66119	1772 (2.74-2.70)

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	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8104 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 HIV-1 reverse transcriptase, p66 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	554	Total	C	N	O	S	0	0	0
			4505	2917	748	832	8			

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

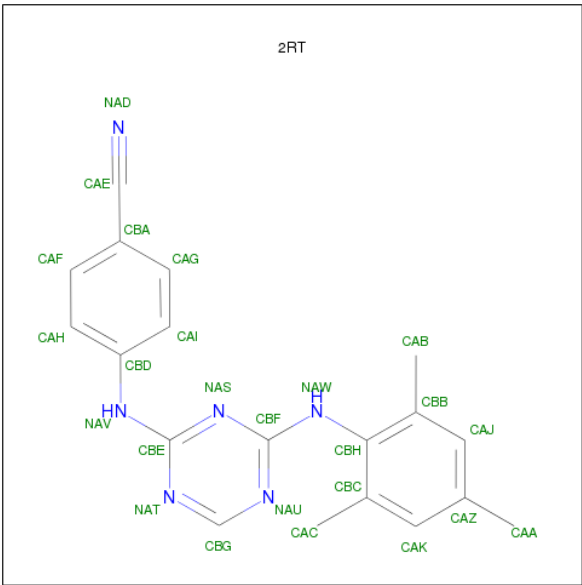
- Molecule 2 is a protein called HIV-1 reverse transcriptase, p51 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	428	Total	C	N	O	S	0	0	0
			3539	2305	586	641	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 4-({4-[(2,4,6-TRIMETHYLPHENYL)AMINO]-1,3,5-TRIAZIN-2-YL}AMINO)BENZONITRILE (three-letter code: 2RT) (formula: C₁₉H₁₈N₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			25	19	6		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	22	Total	O	0	0
			22	22		
4	B	13	Total	O	0	0
			13	13		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	163.04Å 73.40Å 109.73Å 90.00° 100.06° 90.00°	Depositor
Resolution (Å)	43.24 – 2.71 43.24 – 2.71	Depositor EDS
% Data completeness (in resolution range)	98.5 (43.24-2.71) 95.8 (43.24-2.71)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.46 (at 2.73Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1069)	Depositor
R, R_{free}	0.236 , 0.299 0.235 , 0.301	Depositor DCC
R_{free} test set	1948 reflections (5.84%)	DCC
Wilson B-factor (Å ²)	67.3	Xtriage
Anisotropy	0.364	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 22.9	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 34315 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8104	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.14% 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: 2RT

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	0.27	0/4623	0.45	0/6284
2	B	0.27	0/3644	0.48	0/4952
All	All	0.27	0/8267	0.46	0/11236

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	4505	0	4562	79	0
2	B	3539	0	3576	80	0
3	A	25	0	0	0	0
4	A	22	0	0	1	0
4	B	13	0	0	3	0
All	All	8104	0	8138	149	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 (149) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:544:GLY:HA2	2:B:286:THR:HG22	1.57	0.86
1:A:281:LYS:HB2	1:A:284:ARG:HE	1.52	0.74
1:A:50:ILE:HD13	1:A:145:GLN:HG3	1.71	0.71
2:B:116:PHE:HB2	2:B:224:GLU:HG3	1.73	0.69
2:B:114:ALA:HB3	2:B:224:GLU:HB2	1.76	0.68
1:A:28:GLU:HG3	1:A:135:ILE:HD12	1.79	0.65
1:A:491:LEU:HB3	1:A:529:GLU:HG3	1.79	0.64
1:A:259:LYS:HZ2	1:A:265:ASN:HB2	1.65	0.62
2:B:380:ILE:HG12	4:B:509:HOH:O	2.01	0.61
1:A:543:GLY:H	2:B:283:LEU:HB3	1.66	0.61
1:A:249:LYS:NZ	4:A:712:HOH:O	2.33	0.61
2:B:241:VAL:HG22	2:B:350:LYS:HA	1.83	0.60
2:B:214:LEU:HD22	2:B:226:PRO:HB3	1.84	0.60
1:A:246:LEU:HD11	1:A:264:LEU:HD21	1.84	0.60
2:B:191:SER:OG	2:B:198:HIS:ND1	2.27	0.60
1:A:542:ILE:HG23	2:B:283:LEU:HD13	1.85	0.59
1:A:343:GLN:HG3	1:A:349:LEU:HD11	1.84	0.59
2:B:424:LYS:HD3	2:B:425:LEU:HG	1.84	0.59
2:B:218:ASP:HB2	2:B:221:HIS:HB2	1.85	0.59
2:B:180:ILE:HG12	2:B:189:VAL:HG22	1.85	0.59
2:B:193:LEU:HD13	2:B:201:LYS:HE3	1.83	0.59
1:A:544:GLY:N	2:B:285:GLY:O	2.35	0.59
1:A:303:LEU:HD23	1:A:306:ASN:HD22	1.68	0.58
1:A:259:LYS:HA	1:A:260:LEU:HB2	1.86	0.57
1:A:417:VAL:HG22	1:A:419:THR:HG23	1.86	0.56
1:A:225:PRO:HB3	1:A:236:PRO:HD3	1.87	0.56
1:A:80:LEU:HD22	1:A:127:TYR:CZ	2.41	0.56
2:B:80:LEU:O	2:B:84:THR:N	2.39	0.56
1:A:536:VAL:HB	1:A:542:ILE:HD13	1.88	0.56
2:B:210:LEU:HD12	2:B:214:LEU:HD12	1.87	0.55
2:B:104:LYS:HA	2:B:237:ASP:HB2	1.88	0.55
1:A:88:TRP:CD1	2:B:143:ARG:HD2	2.43	0.54
2:B:79:GLU:OE2	2:B:83:ARG:NE	2.41	0.54
2:B:86:ASP:O	2:B:88:TRP:N	2.40	0.54
1:A:209:LEU:HB3	1:A:214:LEU:HB2	1.89	0.53
1:A:405:TYR:CE2	1:A:407:GLN:HB2	2.43	0.53
1:A:195:ILE:O	1:A:199:ARG:HB2	2.08	0.53
2:B:113:ASP:HB3	2:B:214:LEU:HD21	1.92	0.52
1:A:195:ILE:HG13	1:A:199:ARG:HD2	1.90	0.52
1:A:457:TYR:HE1	1:A:463:ARG:HG2	1.74	0.52
2:B:116:PHE:HB3	2:B:223:LYS:HA	1.91	0.52
2:B:241:VAL:HG13	2:B:351:THR:HG23	1.90	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:111:VAL:HB	1:A:114:ALA:HB3	1.91	0.51
2:B:91:GLN:HG3	2:B:161:GLN:HG3	1.93	0.51
1:A:191:SER:OG	1:A:198:HIS:ND1	2.38	0.51
1:A:56:TYR:O	1:A:143:ARG:NH2	2.26	0.50
1:A:73:LYS:NZ	1:A:146:TYR:OH	2.43	0.50
2:B:112:GLY:C	2:B:114:ALA:HB2	2.31	0.50
1:A:266:TRP:O	1:A:269:GLN:NE2	2.44	0.49
1:A:246:LEU:O	1:A:307:ARG:NH1	2.40	0.49
2:B:120:LEU:HD23	2:B:125:ARG:HG2	1.93	0.49
2:B:116:PHE:O	2:B:117:SER:OG	2.22	0.48
1:A:203:GLU:CD	1:A:206:ARG:HH12	2.15	0.48
1:A:209:LEU:HD12	1:A:216:THR:HG21	1.95	0.48
1:A:279:LEU:HD23	1:A:299:ALA:HB1	1.94	0.48
2:B:21:VAL:HB	2:B:59:PRO:HD3	1.95	0.48
2:B:356:ARG:HD2	2:B:374:LYS:HZ2	1.78	0.48
1:A:373:GLN:OE1	2:B:401:TRP:NE1	2.43	0.48
1:A:80:LEU:HD13	1:A:127:TYR:CE2	2.49	0.48
2:B:376:THR:O	2:B:380:ILE:HG13	2.14	0.48
1:A:104:LYS:HB2	1:A:192:ASP:HA	1.96	0.47
2:B:232:TYR:OH	2:B:356:ARG:NH1	2.47	0.47
1:A:490:GLY:O	1:A:528:LYS:NZ	2.42	0.47
1:A:60:VAL:HG11	1:A:130:PHE:HD1	1.80	0.47
1:A:543:GLY:H	2:B:283:LEU:CB	2.26	0.47
2:B:374:LYS:HE3	2:B:374:LYS:HB2	1.70	0.46
1:A:76:ASP:OD1	1:A:78:ARG:HG3	2.15	0.46
2:B:116:PHE:HD1	2:B:148:VAL:HG21	1.79	0.46
2:B:2:ILE:HG22	2:B:4:PRO:HD3	1.98	0.46
2:B:197:GLN:O	2:B:200:THR:OG1	2.31	0.46
2:B:212:TRP:HA	2:B:220:LYS:HG3	1.98	0.46
2:B:163:SER:O	2:B:167:ILE:HG13	2.16	0.46
1:A:58:THR:N	1:A:129:ALA:O	2.50	0.45
1:A:200:THR:HA	1:A:203:GLU:HG2	1.98	0.45
1:A:271:TYR:CD1	1:A:274:ILE:HG13	2.51	0.45
2:B:115:TYR:HB3	2:B:149:LEU:HB2	1.99	0.45
1:A:24:TRP:HB2	1:A:61:PHE:CZ	2.52	0.45
1:A:257:ILE:HG21	1:A:282:LEU:HD12	1.99	0.45
2:B:49:LYS:HE3	2:B:142:ILE:HG23	1.99	0.45
2:B:47:ILE:HD12	2:B:144:TYR:CG	2.52	0.44
2:B:224:GLU:HB3	2:B:225:PRO:HD2	1.98	0.44
1:A:80:LEU:O	1:A:84:THR:OG1	2.26	0.44
1:A:369:THR:O	1:A:373:GLN:HG2	2.17	0.44
1:A:281:LYS:HD2	1:A:284:ARG:HH21	1.82	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:236:PRO:HA	2:B:239:TRP:HD1	1.83	0.44
1:A:379:SER:CB	1:A:387:PRO:HD3	2.47	0.44
2:B:100:LEU:HG	2:B:381:VAL:HG13	2.00	0.44
1:A:441:TYR:CG	1:A:544:GLY:HA3	2.53	0.43
1:A:107:THR:OG1	1:A:198:HIS:NE2	2.41	0.43
1:A:407:GLN:OE1	2:B:394:GLN:HG2	2.18	0.43
1:A:411:ILE:HG22	1:A:412:PRO:O	2.18	0.43
1:A:543:GLY:HA2	2:B:284:ARG:HA	1.98	0.43
1:A:172:ALA:HB2	1:A:180:ILE:HD12	2.00	0.43
2:B:172:LYS:N	4:B:505:HOH:O	2.51	0.43
2:B:369:THR:HG22	2:B:398:TRP:CZ3	2.53	0.43
2:B:11:LYS:HB2	2:B:11:LYS:HE3	1.85	0.43
2:B:308:GLU:HA	2:B:311:LYS:HE2	2.00	0.43
1:A:126:LYS:HA	1:A:145:GLN:NE2	2.33	0.43
1:A:59:PRO:HB2	1:A:61:PHE:CE2	2.53	0.43
1:A:376:THR:HG21	2:B:401:TRP:CZ2	2.54	0.43
2:B:393:ILE:HD13	2:B:398:TRP:HB2	1.99	0.43
1:A:366:LYS:HE2	1:A:405:TYR:OH	2.18	0.43
2:B:116:PHE:CD1	2:B:148:VAL:HG11	2.52	0.43
1:A:114:ALA:HB2	1:A:214:LEU:HB3	2.01	0.43
2:B:123:ASP:O	2:B:126:LYS:NZ	2.51	0.43
2:B:332:GLN:HG3	2:B:338:THR:HG23	2.01	0.43
1:A:199:ARG:O	1:A:203:GLU:HG2	2.20	0.42
2:B:311:LYS:HE3	2:B:311:LYS:HB2	1.84	0.42
2:B:65:LYS:HG3	2:B:72:ARG:HH11	1.85	0.42
2:B:225:PRO:HA	2:B:226:PRO:HD3	1.65	0.42
2:B:320:ASP:HA	2:B:321:PRO:HD3	1.79	0.42
2:B:74:LEU:HD12	4:B:502:HOH:O	2.19	0.42
1:A:342:TYR:HA	1:A:349:LEU:HD13	2.01	0.42
2:B:38:CYS:SG	2:B:132:ILE:HD11	2.60	0.42
1:A:0:VAL:HA	1:A:1:PRO:HD3	1.85	0.42
2:B:287:LYS:HD3	2:B:293:ILE:HD11	2.01	0.41
1:A:80:LEU:O	1:A:127:TYR:OH	2.34	0.41
2:B:19:PRO:HG3	2:B:80:LEU:HB2	2.02	0.41
2:B:208:HIS:HD1	2:B:209:LEU:HD23	1.85	0.41
2:B:365:VAL:O	2:B:369:THR:HG23	2.20	0.41
1:A:118:VAL:HA	1:A:119:PRO:HD3	1.91	0.41
1:A:122:GLU:HA	1:A:125:ARG:HG3	2.03	0.41
2:B:224:GLU:CB	2:B:225:PRO:HD2	2.51	0.41
2:B:72:ARG:HD2	2:B:227:PHE:CE2	2.55	0.41
2:B:335:GLY:O	2:B:355:ALA:HA	2.20	0.41
1:A:171:PHE:CD2	1:A:205:LEU:HD13	2.56	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:114:ALA:HA	1:A:214:LEU:HD22	2.01	0.41
1:A:271:TYR:HA	1:A:272:PRO:HD3	1.75	0.41
1:A:276:VAL:O	1:A:280:SER:OG	2.23	0.41
2:B:113:ASP:N	2:B:114:ALA:HA	2.35	0.41
2:B:218:ASP:HA	2:B:220:LYS:N	2.35	0.41
1:A:7:THR:HG22	1:A:119:PRO:HB2	2.03	0.41
2:B:64:LYS:HE3	2:B:71:TRP:CE2	2.56	0.41
1:A:116:PHE:HD2	1:A:148:VAL:HG21	1.84	0.41
2:B:46:LYS:HZ1	2:B:224:GLU:HA	1.86	0.41
1:A:365:VAL:HG11	1:A:401:TRP:CG	2.56	0.41
1:A:253:THR:HG22	1:A:292:VAL:HG22	2.03	0.41
1:A:303:LEU:HA	1:A:306:ASN:HB2	2.02	0.41
1:A:31:ILE:O	1:A:35:VAL:HG23	2.21	0.41
1:A:441:TYR:CD2	1:A:544:GLY:HA3	2.56	0.41
2:B:149:LEU:HD21	2:B:159:ILE:HD12	2.02	0.41
1:A:97:PRO:HG2	1:A:232:TYR:CD2	2.55	0.41
2:B:171:PHE:C	2:B:173:LYS:H	2.25	0.41
2:B:153:TRP:HB3	2:B:156:SER:OG	2.20	0.40
1:A:401:TRP:O	1:A:405:TYR:HB2	2.22	0.40
2:B:240:THR:OG1	2:B:241:VAL:N	2.55	0.40
2:B:368:LEU:O	2:B:372:VAL:HG23	2.21	0.40
2:B:423:VAL:HG12	2:B:426:TRP:CE3	2.56	0.40
2:B:205:LEU:O	2:B:209:LEU:HG	2.21	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	552/557 (99%)	516 (94%)	33 (6%)	3 (0%)	38	70
2	B	426/428 (100%)	381 (89%)	39 (9%)	6 (1%)	16	39
All	All	978/985 (99%)	897 (92%)	72 (7%)	9 (1%)	25	54

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	87	PHE
1	A	243	PRO
2	B	93	GLY
2	B	214	LEU
2	B	114	ALA
1	A	9	PRO
1	A	545	ASN
2	B	111	VAL
2	B	94	ILE

5.3.2 Protein sidechains ⓘ

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	494/495 (100%)	485 (98%)	9 (2%)	71	93
2	B	390/390 (100%)	383 (98%)	7 (2%)	71	93
All	All	884/885 (100%)	868 (98%)	16 (2%)	71	93

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	LYS
1	A	80	LEU
1	A	102	LYS
1	A	194	GLU
1	A	283	LEU
1	A	291	GLU
1	A	367	GLN
1	A	449	GLU
1	A	497	THR
2	B	39	THR
2	B	65	LYS
2	B	96	HIS
2	B	171	PHE
2	B	215	THR
2	B	220	LYS

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Mol	Chain	Res	Type
2	B	424	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN

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 ⓘ

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	2RT	A	601	-	27,27,27	1.51	5 (18%)	37,37,37	2.22	5 (13%)

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	2RT	A	601	-	-	0/10/10/10	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	2RT	CBE-NAV	3.02	1.41	1.36
3	A	601	2RT	CAH-CBD	2.78	1.43	1.39
3	A	601	2RT	CBH-CBC	2.26	1.43	1.40
3	A	601	2RT	CAI-CBD	2.17	1.42	1.39
3	A	601	2RT	CBF-NAW	2.07	1.39	1.36

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	2RT	NAT-CBE-NAS	-7.74	120.08	126.70
3	A	601	2RT	NAU-CBF-NAS	-6.17	121.43	126.70
3	A	601	2RT	CBF-NAS-CBE	5.12	121.14	114.05
3	A	601	2RT	CBC-CBH-NAW	3.29	122.90	119.31
3	A	601	2RT	NAU-CBG-NAT	-2.91	123.73	128.66

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	554/557 (99%)	0.81	89 (16%) 2 3	9, 66, 126, 145	0
2	B	428/428 (100%)	0.88	64 (14%) 3 3	13, 54, 121, 162	0
All	All	982/985 (99%)	0.84	153 (15%) 3 3	9, 61, 124, 162	0

All (153) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	69	THR	9.3
2	B	216	THR	8.8
1	A	263	LYS	8.2
2	B	217	PRO	8.0
1	A	273	GLY	7.8
1	A	286	THR	7.8
1	A	271	TYR	7.7
1	A	309	ILE	7.7
1	A	257	ILE	7.5
1	A	289	LEU	7.4
1	A	74	LEU	7.2
2	B	227	PHE	6.7
2	B	95	PRO	6.7
2	B	96	HIS	6.5
1	A	290	THR	6.3
2	B	202	ILE	6.3
2	B	85	GLN	6.0
1	A	24	TRP	5.7
1	A	308	GLU	5.7
1	A	221	HIS	5.5
2	B	15	GLY	5.3
2	B	232	TYR	5.3
2	B	215	THR	5.2
2	B	225	PRO	5.1

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Mol	Chain	Res	Type	RSRZ
2	B	237	ASP	5.1
1	A	314	VAL	5.1
2	B	88	TRP	5.0
1	A	285	GLY	4.9
1	A	275	LYS	4.9
2	B	222	GLN	4.7
1	A	301	LEU	4.7
2	B	149	LEU	4.7
2	B	212	TRP	4.6
1	A	250	ASP	4.6
1	A	298	GLU	4.5
1	A	151	GLN	4.5
2	B	209	LEU	4.5
1	A	67	ASP	4.4
2	B	239	TRP	4.3
1	A	254	VAL	4.3
2	B	219	LYS	4.3
2	B	14	PRO	4.2
1	A	71	TRP	4.2
1	A	70	LYS	4.2
1	A	284	ARG	4.1
1	A	217	PRO	4.1
1	A	255	ASN	4.1
2	B	2	ILE	4.0
1	A	288	ALA	4.0
1	A	551	LEU	4.0
1	A	287	LYS	4.0
2	B	1	PRO	3.9
1	A	547	GLN	3.9
1	A	295	LEU	3.9
2	B	231	GLY	3.8
2	B	221	HIS	3.8
2	B	226	PRO	3.7
1	A	2	ILE	3.7
1	A	256	ASP	3.7
1	A	241	VAL	3.7
1	A	260	LEU	3.6
1	A	291	GLU	3.5
1	A	274	ILE	3.5
1	A	272	PRO	3.5
1	A	63	ILE	3.5
2	B	16	MET	3.4

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Mol	Chain	Res	Type	RSRZ
2	B	201	LYS	3.4
2	B	3	SER	3.4
1	A	142	ILE	3.4
2	B	90	VAL	3.4
2	B	4	PRO	3.3
2	B	223	LYS	3.3
2	B	12	LEU	3.3
2	B	189	VAL	3.3
2	B	69	THR	3.3
1	A	549	ASP	3.2
1	A	248	GLU	3.2
1	A	258	GLN	3.2
2	B	9	PRO	3.2
1	A	252	TRP	3.2
2	B	190	GLY	3.2
1	A	40	GLU	3.2
1	A	19	PRO	3.1
1	A	282	LEU	3.0
1	A	127	TYR	3.0
1	A	552	VAL	3.0
2	B	116	PHE	3.0
1	A	259	LYS	2.9
1	A	171	PHE	2.9
1	A	53	GLU	2.9
1	A	249	LYS	2.9
1	A	228	LEU	2.9
1	A	305	GLU	2.9
1	A	262	GLY	2.9
1	A	52	PRO	2.8
1	A	300	GLU	2.8
1	A	294	PRO	2.8
1	A	167	ILE	2.8
1	A	279	LEU	2.8
1	A	72	ARG	2.7
2	B	66	LYS	2.7
2	B	230	MET	2.7
2	B	150	PRO	2.7
1	A	548	VAL	2.7
1	A	148	VAL	2.6
2	B	428	GLN	2.6
2	B	284	ARG	2.6
1	A	20	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	91	GLN	2.6
2	B	124	PHE	2.6
2	B	236	PRO	2.5
2	B	171	PHE	2.5
1	A	270	ILE	2.5
1	A	546	GLU	2.5
2	B	170	PRO	2.5
2	B	167	ILE	2.5
1	A	222	GLN	2.5
1	A	278	GLN	2.5
1	A	310	LEU	2.4
1	A	132	ILE	2.4
1	A	293	ILE	2.4
2	B	93	GLY	2.4
1	A	116	PHE	2.4
1	A	14	PRO	2.4
1	A	315	HIS	2.4
1	A	357	MET	2.4
1	A	130	PHE	2.3
2	B	205	LEU	2.3
2	B	356	ARG	2.3
2	B	166	LYS	2.3
2	B	206	ARG	2.3
1	A	144	TYR	2.3
1	A	51	GLY	2.3
1	A	58	THR	2.3
1	A	264	LEU	2.3
2	B	211	ARG	2.3
2	B	172	LYS	2.2
2	B	229	TRP	2.2
2	B	11	LYS	2.2
1	A	193	LEU	2.2
2	B	358	ARG	2.2
1	A	223	LYS	2.2
2	B	238	LYS	2.2
2	B	68	SER	2.1
2	B	214	LEU	2.1
1	A	43	LYS	2.1
2	B	65	LYS	2.1
2	B	318	TYR	2.1
1	A	47	ILE	2.1
1	A	73	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	-1	MET	2.1
2	B	359	GLY	2.0
2	B	182	GLN	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
3	2RT	A	601	25/25	0.17	-0.36	37,48,63,72	0

6.5 Other polymers ⓘ

There are no such residues in this entry.