



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:19 AM GMT

PDB ID : 2WOM
Title : Crystal Structure of UK-453061 bound to HIV-1 Reverse Transcriptase (K103N).
Authors : Phillips, C.; Irving, S.L.; Knoechel, T.; Ringrose, H.
Deposited on : 2009-07-27
Resolution : 3.20 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

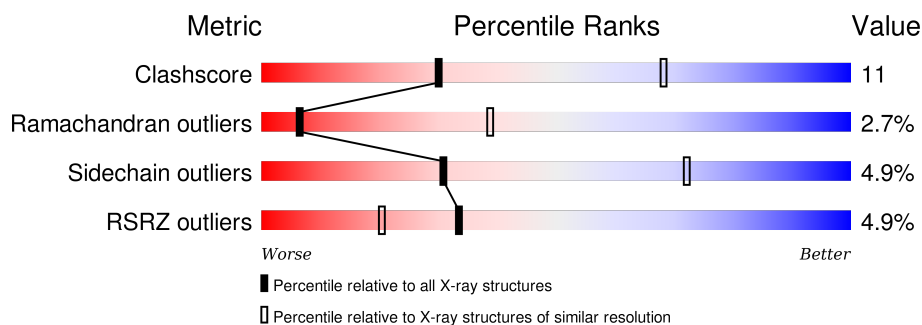
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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))
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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. 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	560	<div> <div>5%</div> <div> <div></div> <div>73%</div> <div>24%</div> <div>.</div> </div> </div>
2	B	440	<div> <div>5%</div> <div> <div></div> <div>70%</div> <div>20%</div> <div>.</div> <div>6%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7942 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 HIV-1 REVERSE TRANSCRIPTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	558	Total	C	N	O	S	0	0	1
			4536	2929	760	839	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	103	ASN	LYS	ENGINEERED MUTATION	UNP P04585

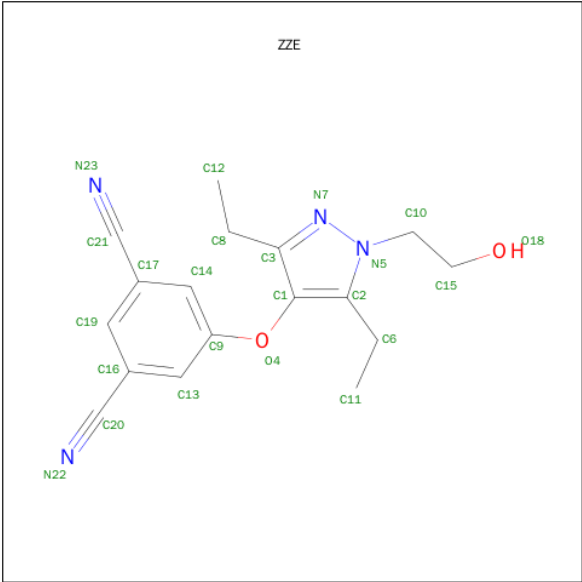
- Molecule 2 is a protein called HIV-1 REVERSE TRANSCRIPTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	413	Total	C	N	O	S	0	0	1
			3383	2195	565	616	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	103	ASN	LYS	ENGINEERED MUTATION	UNP P04585

- Molecule 3 is 5-{{[3,5-DIETHYL-1-(2-HYDROXYETHYL)-1H-PYRAZOL-4-YL]OXY}BENZENE-1,3-DICARBONITRILE (three-letter code: ZZE) (formula: C₁₇H₁₈N₄O₂).

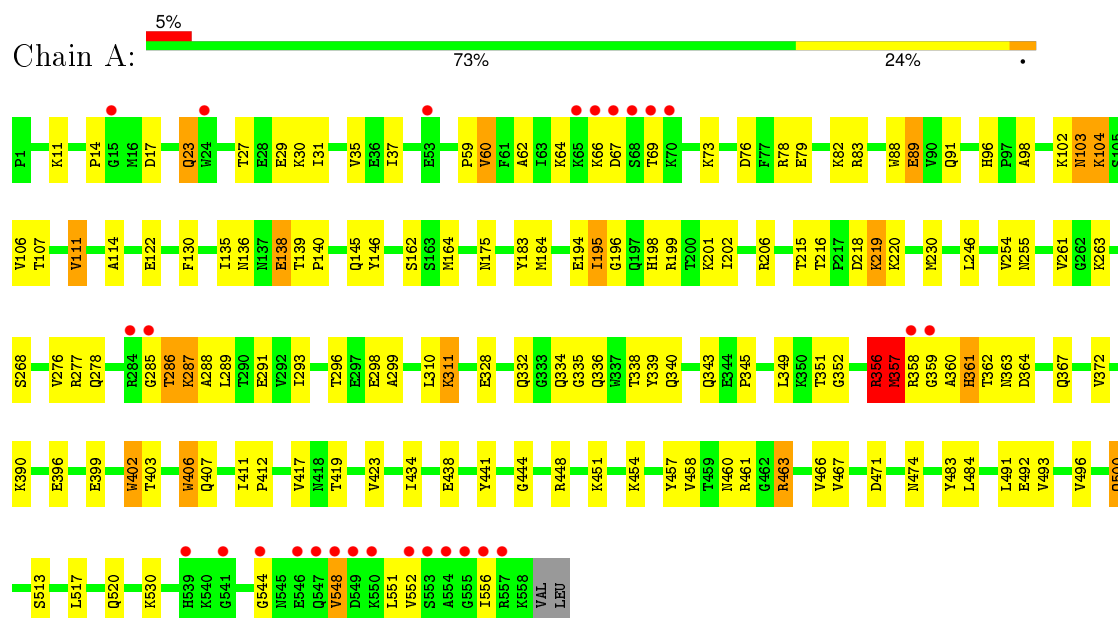


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			23	17	4	2		

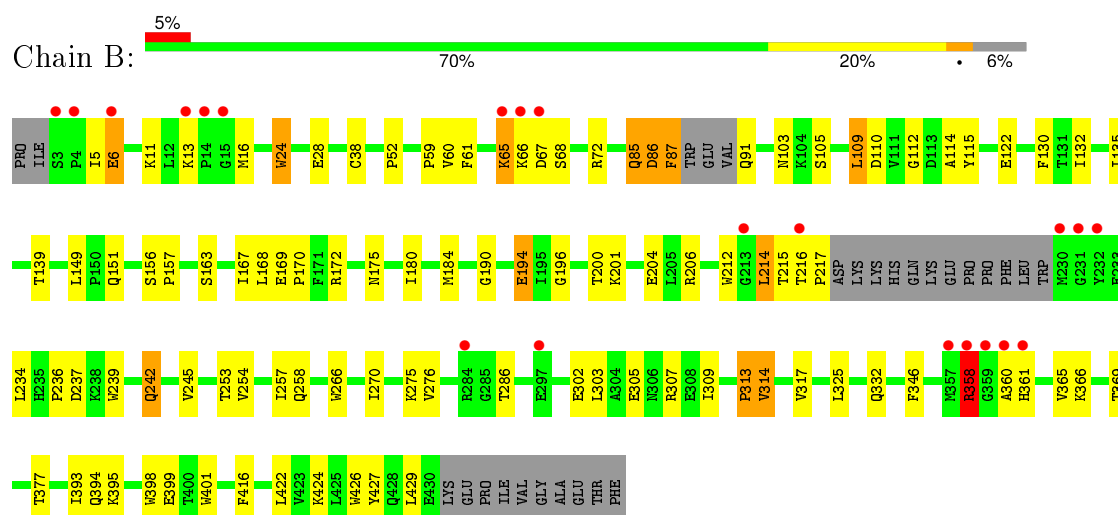
3 Residue-property plots

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: HIV-1 REVERSE TRANSCRIPTASE



• Molecule 2: HIV-1 REVERSE TRANSCRIPTASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	116.60Å 154.30Å 155.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.95 – 3.20 24.95 – 2.79	Depositor EDS
% Data completeness (in resolution range)	(Not available) (24.95-3.20) 87.5 (24.95-2.79)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.62 (at 2.80Å)	Xtriage
Refinement program	BUSTER-TNT 2.9.2	Depositor
R, R_{free}	0.228 , 0.257 0.267 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	57.3	Xtriage
Anisotropy	0.370	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 32428 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	7942	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: ZZE

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	0.46	1/4653 (0.0%)	0.60	0/6323
2	B	0.36	0/3476	0.57	0/4724
All	All	0.42	1/8129 (0.0%)	0.59	0/11047

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	103	ASN	C-N	18.72	1.77	1.34

There are no bond angle outliers.

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	4536	0	4582	106	0
2	B	3383	0	3410	77	0
3	A	23	0	18	2	0
All	All	7942	0	8010	177	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 (177) 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:103:ASN:C	1:A:104:LYS:N	1.77	1.36
1:A:139:THR:HB	1:A:140:PRO:HD2	1.43	1.00
1:A:23:GLN:HE22	1:A:60:VAL:H	1.06	0.93
2:B:13:LYS:HB2	2:B:16:MET:HG3	1.62	0.82
1:A:23:GLN:NE2	1:A:60:VAL:H	1.79	0.81
1:A:278:GLN:HG2	1:A:298:GLU:HB3	1.63	0.80
1:A:296:THR:HG23	1:A:299:ALA:H	1.47	0.79
2:B:85:GLN:HA	2:B:87:PHE:N	1.98	0.79
2:B:85:GLN:HA	2:B:87:PHE:H	1.53	0.74
1:A:111:VAL:HG12	1:A:111:VAL:O	1.90	0.71
1:A:461:ARG:HH11	1:A:461:ARG:HG3	1.55	0.71
2:B:360:ALA:HB2	2:B:366:LYS:HD2	1.72	0.70
1:A:372:VAL:HG11	1:A:411:ILE:HG23	1.73	0.70
1:A:328:GLU:HG2	1:A:390:LYS:HB2	1.72	0.70
1:A:500:GLN:NE2	2:B:422:LEU:HG	2.07	0.69
2:B:86:ASP:HA	2:B:91:GLN:HB2	1.75	0.69
1:A:175:ASN:OD1	1:A:201:LYS:HE3	1.92	0.69
2:B:13:LYS:HZ1	2:B:85:GLN:HG2	1.58	0.68
1:A:412:PRO:HG3	2:B:401:TRP:HZ2	1.61	0.65
1:A:492:GLU:HG2	1:A:530:LYS:HB2	1.79	0.65
1:A:138:GLU:HG2	1:A:139:THR:N	2.12	0.64
1:A:111:VAL:CG1	1:A:111:VAL:O	2.46	0.64
2:B:13:LYS:HD2	2:B:16:MET:HE3	1.80	0.64
2:B:5:ILE:HG22	2:B:6:GLU:H	1.63	0.62
2:B:114:ALA:HB2	2:B:214:LEU:HD13	1.80	0.62
1:A:458:VAL:HG12	1:A:548:VAL:HG22	1.79	0.62
1:A:362:THR:HG22	1:A:363:ASN:N	2.16	0.61
2:B:24:TRP:HE1	2:B:59:PRO:HB3	1.65	0.61
1:A:500:GLN:HE21	1:A:500:GLN:HA	1.65	0.60
1:A:399:GLU:HA	1:A:402:TRP:CD1	2.37	0.60
1:A:27:THR:CG2	1:A:29:GLU:HG2	2.31	0.60
1:A:106:VAL:HG21	3:A:1558:ZZE:H152	1.83	0.60
1:A:254:VAL:HG22	1:A:293:ILE:HD11	1.84	0.60
1:A:500:GLN:HE22	2:B:422:LEU:HG	1.66	0.59
2:B:206:ARG:HD2	2:B:216:THR:O	2.02	0.59
1:A:277:ARG:HB3	1:A:336:GLN:OE1	2.02	0.59
1:A:73:LYS:NZ	1:A:146:TYR:OH	2.36	0.59
2:B:24:TRP:NE1	2:B:59:PRO:HB3	2.18	0.59
1:A:311:LYS:HE2	1:A:311:LYS:HA	1.83	0.59
1:A:103:ASN:C	1:A:104:LYS:CA	2.72	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:THR:HB	1:A:140:PRO:CD	2.26	0.57
1:A:551:LEU:HD23	1:A:551:LEU:H	1.69	0.57
2:B:215:THR:O	2:B:217:PRO:HD3	2.03	0.57
1:A:263:LYS:HA	1:A:263:LYS:HE2	1.86	0.57
1:A:206:ARG:HG3	1:A:216:THR:OG1	2.05	0.57
1:A:451:LYS:HE2	1:A:471:ASP:HA	1.87	0.57
1:A:261:VAL:HG13	1:A:276:VAL:HG11	1.85	0.57
1:A:412:PRO:HG3	2:B:401:TRP:CZ2	2.40	0.57
2:B:365:VAL:O	2:B:369:THR:HG23	2.04	0.56
1:A:23:GLN:HE22	1:A:60:VAL:N	1.90	0.55
1:A:548:VAL:HA	1:A:551:LEU:HD21	1.89	0.55
2:B:242:GLN:HE22	2:B:429:LEU:CD1	2.19	0.55
2:B:110:ASP:O	2:B:216:THR:HG23	2.06	0.55
1:A:122:GLU:CD	1:A:122:GLU:H	2.11	0.54
1:A:444:GLY:HA2	1:A:552:VAL:HG11	1.89	0.54
2:B:66:LYS:HG3	2:B:67:ASP:OD1	2.07	0.54
1:A:268:SER:O	1:A:351:THR:HG22	2.07	0.54
2:B:360:ALA:CB	2:B:366:LYS:HD2	2.38	0.54
1:A:194:GLU:H	1:A:194:GLU:CD	2.10	0.53
1:A:343:GLN:HG3	1:A:349:LEU:HD11	1.90	0.53
1:A:288:ALA:HB3	1:A:291:GLU:HG2	1.90	0.53
1:A:111:VAL:HG21	1:A:164:MET:CE	2.39	0.52
1:A:466:VAL:HB	1:A:551:LEU:HD12	1.91	0.52
1:A:27:THR:HG21	1:A:29:GLU:HG2	1.90	0.52
2:B:172:ARG:HH21	2:B:180:ILE:HB	1.75	0.52
2:B:175:ASN:HD21	2:B:201:LYS:NZ	2.08	0.52
2:B:266:TRP:CH2	2:B:427:TYR:CZ	2.98	0.51
1:A:286:THR:O	1:A:287:LYS:HG2	2.10	0.51
1:A:59:PRO:HG2	1:A:76:ASP:HB3	1.93	0.51
2:B:266:TRP:HH2	2:B:427:TYR:CZ	2.28	0.51
2:B:168:LEU:HD13	2:B:180:ILE:HG21	1.93	0.51
1:A:335:GLY:HA2	1:A:367:GLN:OE1	2.11	0.51
1:A:23:GLN:NE2	1:A:59:PRO:HA	2.25	0.51
2:B:314:VAL:HG13	2:B:317:VAL:HG22	1.93	0.51
1:A:27:THR:O	1:A:31:ILE:HG13	2.12	0.50
1:A:96:HIS:HD1	1:A:98:ALA:H	1.59	0.50
1:A:64:LYS:NZ	1:A:69:THR:HG23	2.26	0.50
1:A:79:GLU:HG3	1:A:83:ARG:NH1	2.25	0.50
1:A:195:ILE:HB	1:A:199:ARG:NH2	2.26	0.50
2:B:13:LYS:NZ	2:B:85:GLN:HG2	2.26	0.50
2:B:266:TRP:HH2	2:B:427:TYR:CE1	2.29	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:156:SER:HB2	2:B:157:PRO:HD3	1.93	0.50
2:B:157:PRO:HG2	2:B:184:MET:HA	1.93	0.50
2:B:109:LEU:HG	2:B:216:THR:HG22	1.93	0.49
2:B:60:VAL:HG21	2:B:130:PHE:CD2	2.47	0.49
1:A:441:TYR:CD2	1:A:544:GLY:HA3	2.47	0.49
1:A:556:ILE:O	1:A:556:ILE:HG22	2.13	0.49
2:B:303:LEU:HD22	2:B:307:ARG:HH21	1.76	0.49
2:B:13:LYS:HD2	2:B:16:MET:CE	2.41	0.49
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.48	0.49
1:A:60:VAL:HG22	1:A:130:PHE:HB2	1.95	0.49
2:B:5:ILE:HG22	2:B:6:GLU:N	2.25	0.49
1:A:27:THR:HG22	1:A:29:GLU:HG2	1.95	0.49
2:B:369:THR:HG22	2:B:398:TRP:CH2	2.48	0.48
1:A:17:ASP:O	1:A:83:ARG:HD3	2.13	0.48
1:A:106:VAL:HG11	3:A:1558:ZZE:C1	2.43	0.48
1:A:434:ILE:HD13	1:A:530:LYS:HB3	1.96	0.48
1:A:417:VAL:HG13	1:A:419:THR:HG22	1.96	0.48
1:A:246:LEU:HD11	1:A:310:LEU:HD12	1.96	0.48
1:A:361:HIS:HD2	1:A:513:SER:OG	1.96	0.48
1:A:285:GLY:O	1:A:287:LYS:N	2.47	0.47
2:B:24:TRP:NE1	2:B:61:PHE:CZ	2.82	0.47
2:B:112:GLY:HA3	2:B:151:GLN:HE21	1.79	0.47
1:A:362:THR:CG2	1:A:363:ASN:N	2.78	0.47
1:A:139:THR:CB	1:A:140:PRO:HD2	2.31	0.47
1:A:461:ARG:HH11	1:A:461:ARG:CG	2.26	0.47
2:B:163:SER:O	2:B:167:ILE:HG13	2.14	0.46
1:A:332:GLN:HG3	1:A:338:THR:HG23	1.97	0.46
1:A:206:ARG:HD3	1:A:218:ASP:OD1	2.15	0.46
2:B:275:LYS:HB2	2:B:302:GLU:HG3	1.98	0.46
2:B:200:THR:O	2:B:204:GLU:HG3	2.15	0.46
2:B:234:LEU:HD21	2:B:377:THR:CG2	2.45	0.46
1:A:23:GLN:NE2	1:A:60:VAL:N	2.54	0.46
2:B:358:ARG:HD3	2:B:358:ARG:H	1.81	0.46
1:A:362:THR:HG22	1:A:363:ASN:H	1.81	0.46
2:B:266:TRP:CZ3	2:B:426:TRP:HB3	2.50	0.46
1:A:406:TRP:HE3	1:A:407:GLN:NE2	2.14	0.46
1:A:454:LYS:HG3	1:A:556:ILE:CD1	2.46	0.45
2:B:65:LYS:O	2:B:68:SER:HB3	2.17	0.45
2:B:122:GLU:CD	2:B:122:GLU:H	2.20	0.45
1:A:458:VAL:CG1	1:A:548:VAL:HG22	2.47	0.45
1:A:454:LYS:O	1:A:552:VAL:HG13	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:ASN:HA	2:B:286:THR:O	2.17	0.45
2:B:270:ILE:HG12	2:B:346:PHE:HB3	1.99	0.44
1:A:198:HIS:O	1:A:202:ILE:HG12	2.17	0.44
2:B:242:GLN:HE22	2:B:429:LEU:HD13	1.82	0.44
1:A:255:ASN:HB2	1:A:289:LEU:HG	2.00	0.44
2:B:395:LYS:HG3	2:B:416:PHE:CE2	2.51	0.44
1:A:30:LYS:HE2	1:A:62:ALA:O	2.18	0.44
1:A:467:VAL:HG23	1:A:484:LEU:HD11	1.99	0.44
2:B:194:GLU:HG3	2:B:196:GLY:H	1.81	0.44
2:B:266:TRP:HE1	2:B:346:PHE:HE1	1.65	0.44
2:B:242:GLN:O	2:B:242:GLN:HG3	2.17	0.43
2:B:169:GLU:HB3	2:B:170:PRO:HD3	2.00	0.43
2:B:276:VAL:HA	2:B:302:GLU:OE2	2.18	0.43
1:A:78:ARG:O	1:A:82:LYS:HG3	2.18	0.43
1:A:27:THR:HG22	1:A:29:GLU:H	1.81	0.43
2:B:24:TRP:HE1	2:B:59:PRO:CB	2.28	0.43
2:B:72:ARG:HH22	2:B:151:GLN:NE2	2.16	0.43
1:A:364:ASP:HB3	1:A:423:VAL:HG13	2.00	0.43
2:B:313:PRO:HB2	2:B:314:VAL:H	1.65	0.43
1:A:359:GLY:C	1:A:361:HIS:H	2.21	0.43
1:A:457:TYR:HA	1:A:548:VAL:HG11	2.00	0.42
1:A:339:TYR:CZ	1:A:352:GLY:HA3	2.54	0.42
1:A:23:GLN:NE2	1:A:60:VAL:HG23	2.34	0.42
1:A:402:TRP:CZ2	1:A:403:THR:HG22	2.55	0.42
1:A:162:SER:OG	2:B:52:PRO:HD3	2.19	0.42
1:A:31:ILE:O	1:A:35:VAL:HG23	2.20	0.42
2:B:303:LEU:HD22	2:B:307:ARG:NH2	2.35	0.42
2:B:254:VAL:O	2:B:258:GLN:HG3	2.19	0.42
1:A:438:GLU:OE1	1:A:463:ARG:NH2	2.49	0.42
1:A:218:ASP:O	1:A:220:LYS:N	2.53	0.42
2:B:167:ILE:HG12	2:B:212:TRP:CD2	2.54	0.42
2:B:115:TYR:HB3	2:B:149:LEU:HB2	2.01	0.42
2:B:332:GLN:HA	2:B:424:LYS:HE3	2.02	0.42
1:A:438:GLU:CG	1:A:461:ARG:HD2	2.50	0.42
2:B:103:ASN:O	2:B:236:PRO:HG2	2.20	0.41
2:B:393:ILE:HG12	2:B:394:GLN:N	2.36	0.41
2:B:105:SER:O	2:B:190:GLY:HA2	2.21	0.41
1:A:111:VAL:HG21	1:A:164:MET:HE1	2.01	0.41
2:B:266:TRP:CE3	2:B:426:TRP:HB3	2.56	0.41
2:B:13:LYS:HB2	2:B:16:MET:CG	2.43	0.41
1:A:183:TYR:CE1	1:A:184:MET:HG3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:LYS:CE	1:A:311:LYS:HA	2.50	0.41
1:A:278:GLN:HG2	1:A:298:GLU:CB	2.42	0.41
1:A:357:MET:O	1:A:359:GLY:N	2.36	0.41
1:A:66:LYS:NZ	1:A:67:ASP:H	2.19	0.41
1:A:454:LYS:HG3	1:A:556:ILE:HD11	2.03	0.41
1:A:356:ARG:HD2	1:A:357:MET:H	1.86	0.41
2:B:305:GLU:O	2:B:309:ILE:HG13	2.21	0.41
2:B:28:GLU:HB2	2:B:135:ILE:HD11	2.03	0.41
2:B:253:THR:O	2:B:257:ILE:HG13	2.21	0.40
1:A:483:TYR:CE1	1:A:520:GLN:HB3	2.57	0.40
2:B:38:CYS:SG	2:B:132:ILE:HD11	2.61	0.40
2:B:314:VAL:HG13	2:B:317:VAL:CG2	2.50	0.40
1:A:548:VAL:HA	1:A:551:LEU:CD2	2.51	0.40
1:A:396:GLU:H	1:A:396:GLU:CD	2.25	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	556/560 (99%)	506 (91%)	31 (6%)	19 (3%)	5	31
2	B	407/440 (92%)	383 (94%)	17 (4%)	7 (2%)	11	52
All	All	963/1000 (96%)	889 (92%)	48 (5%)	26 (3%)	6	39

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	GLU
1	A	135	ILE
1	A	195	ILE
1	A	219	LYS

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Mol	Chain	Res	Type
1	A	286	THR
1	A	358	ARG
1	A	491	LEU
2	B	85	GLN
2	B	313	PRO
1	A	14	PRO
1	A	111	VAL
1	A	114	ALA
1	A	357	MET
2	B	86	ASP
2	B	314	VAL
2	B	361	HIS
1	A	88	TRP
1	A	196	GLY
1	A	360	ALA
1	A	361	HIS
2	B	358	ARG
1	A	138	GLU
1	A	287	LYS
1	A	345	PRO
1	A	356	ARG
2	B	245	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	497/500 (99%)	468 (94%)	29 (6%)	25	66
2	B	372/400 (93%)	358 (96%)	14 (4%)	40	78
All	All	869/900 (97%)	826 (95%)	43 (5%)	31	72

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LYS

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Mol	Chain	Res	Type
1	A	23	GLN
1	A	37	ILE
1	A	60	VAL
1	A	89	GLU
1	A	91	GLN
1	A	102	LYS
1	A	104	LYS
1	A	107	THR
1	A	136	ASN
1	A	145	GLN
1	A	215	THR
1	A	219	LYS
1	A	230	MET
1	A	311	LYS
1	A	334	GLN
1	A	340	GLN
1	A	356	ARG
1	A	357	MET
1	A	402	TRP
1	A	406	TRP
1	A	448	ARG
1	A	463	ARG
1	A	474	ASN
1	A	493	VAL
1	A	496	VAL
1	A	500	GLN
1	A	517	LEU
1	A	548	VAL
2	B	6	GLU
2	B	11	LYS
2	B	24	TRP
2	B	65	LYS
2	B	87	PHE
2	B	109	LEU
2	B	139	THR
2	B	194	GLU
2	B	214	LEU
2	B	237	ASP
2	B	242	GLN
2	B	325	LEU
2	B	358	ARG
2	B	399	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	137	ASN
1	A	174	GLN
1	A	182	GLN
1	A	207	GLN
1	A	269	GLN
1	A	278	GLN
1	A	334	GLN
1	A	361	HIS
1	A	373	GLN
1	A	407	GLN
1	A	428	GLN
1	A	480	GLN
1	A	500	GLN
1	A	512	GLN
1	A	519	ASN
1	A	545	ASN
1	A	547	GLN
2	B	137	ASN
2	B	147	ASN
2	B	151	GLN
2	B	175	ASN
2	B	182	GLN
2	B	208	HIS
2	B	242	GLN
2	B	278	GLN
2	B	336	GLN
2	B	394	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 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	ZZE	A	1558	-	23,24,24	2.15	9 (39%)	26,32,32	3.91	7 (26%)

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	ZZE	A	1558	-	-	0/13/15/15	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1558	ZZE	C10-N5	-3.66	1.40	1.47
3	A	1558	ZZE	O4-C9	-2.00	1.35	1.39
3	A	1558	ZZE	C20-N22	2.08	1.19	1.14
3	A	1558	ZZE	C21-N23	2.51	1.20	1.14
3	A	1558	ZZE	C19-C16	2.52	1.44	1.39
3	A	1558	ZZE	C10-C15	2.56	1.58	1.50
3	A	1558	ZZE	C13-C16	3.25	1.45	1.39
3	A	1558	ZZE	C19-C17	4.22	1.47	1.39
3	A	1558	ZZE	C13-C9	4.87	1.47	1.38

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1558	ZZE	C14-C17-C21	-11.05	105.55	119.51
3	A	1558	ZZE	C15-C10-N5	-2.54	105.66	110.85
3	A	1558	ZZE	C6-C2-N5	3.78	128.87	120.20
3	A	1558	ZZE	C2-C1-C3	4.68	110.81	104.05
3	A	1558	ZZE	C3-N7-N5	7.63	111.00	104.42
3	A	1558	ZZE	C9-O4-C1	8.59	133.42	118.46
3	A	1558	ZZE	C19-C17-C21	9.14	131.07	119.51

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1558	ZZE	2	0

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	558/560 (99%)	0.05	27 (4%) 34 21	22, 49, 90, 129	0
2	B	413/440 (93%)	-0.05	21 (5%) 32 18	17, 47, 90, 116	0
All	All	971/1000 (97%)	0.01	48 (4%) 33 20	17, 49, 90, 129	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	549	ASP	8.2
1	A	359	GLY	6.9
2	B	216	THR	6.1
1	A	548	VAL	5.9
1	A	553	SER	5.2
2	B	3	SER	5.2
1	A	550	LYS	5.2
1	A	547	GLN	5.1
2	B	357	MET	5.0
2	B	4	PRO	4.9
2	B	358	ARG	4.8
2	B	230	MET	4.7
2	B	13	LYS	4.1
1	A	556	ILE	4.1
2	B	359	GLY	4.0
2	B	14	PRO	3.8
1	A	358	ARG	3.7
1	A	546	GLU	3.6
2	B	66	LYS	3.5
1	A	555	GLY	3.5
2	B	297	GLU	3.4
1	A	67	ASP	3.4
1	A	544	GLY	3.3
2	B	231	GLY	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	552	VAL	3.1
1	A	557	ARG	3.0
1	A	24	TRP	3.0
1	A	539	HIS	3.0
1	A	69	THR	2.9
1	A	541	GLY	2.9
2	B	67	ASP	2.9
1	A	53	GLU	2.8
2	B	65	LYS	2.7
2	B	361	HIS	2.7
1	A	554	ALA	2.4
1	A	66	LYS	2.4
2	B	360	ALA	2.4
1	A	284	ARG	2.3
2	B	232	TYR	2.2
2	B	213	GLY	2.2
2	B	284	ARG	2.2
1	A	15	GLY	2.2
1	A	65	LYS	2.1
1	A	285	GLY	2.1
1	A	70	LYS	2.1
1	A	68	SER	2.0
2	B	6	GLU	2.0
2	B	15	GLY	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 carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ZZE	A	1558	23/23	0.86	0.27	0.69	32,41,44,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.