



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:13 AM GMT

PDB ID : 3LAK
Title : Crystal structure of HIV-1 reverse transcriptase in complex with N1-heterocycle pyrimidinedione non-nucleoside inhibitor
Authors : Lansdon, E.B.; Mitchell, M.L.
Deposited on : 2010-01-06
Resolution : 2.30 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

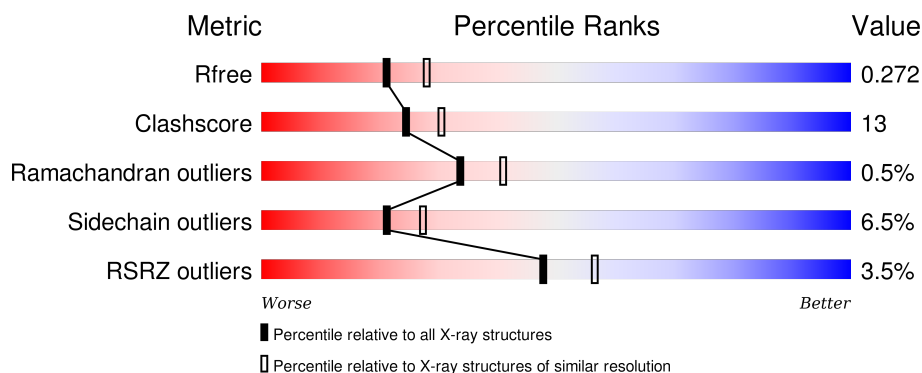
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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>3%</div> <div>67%</div> <div>28%</div> <div>..</div> </div>
1	B	560	<div> <div>3%</div> <div>53%</div> <div>18%</div> <div>28%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

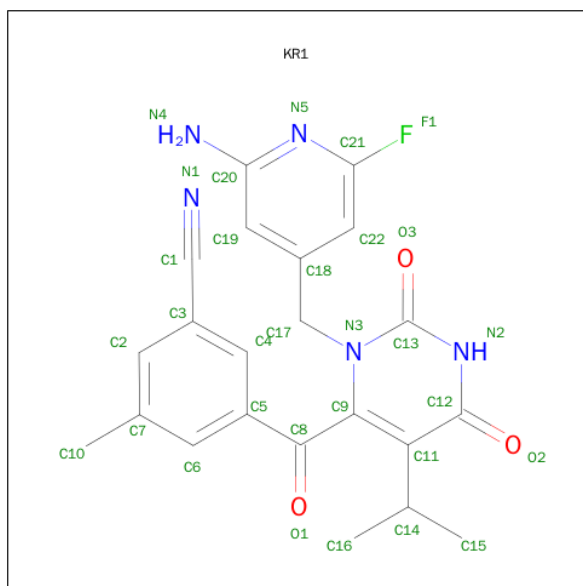
Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	563	-	-	-	X

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 Reverse transcriptase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	553	Total 4508	C 2914	N 752	O 834	S 8	0	0	0
1	B	403	Total 3334	C 2173	N 550	O 605	S 6	0	0	0

- Molecule 2 is 3-({3-[(2-AMINO-6-FLUOROPYRIDIN-4-YL)METHYL]-5-(1-METHYLETHYL)-2,6-DIOXO-1,2,3,6-TETRAHYDOPYRIMIDIN-4-YL}CARBONYL)-5-METHYLETHYL ENZONITRILE (three-letter code: KR1) (formula: $C_{22}H_{20}FN_5O_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			31	22	1	5	3		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		

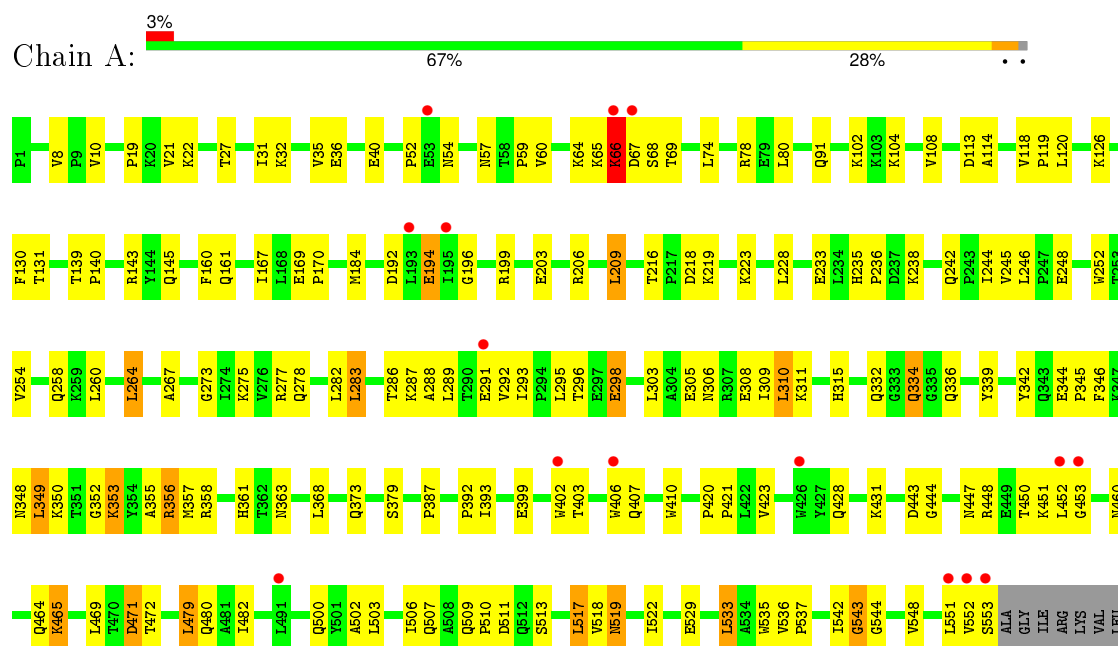
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	110	Total	O	0	0
			110	110		
5	B	97	Total	O	0	0
			97	97		

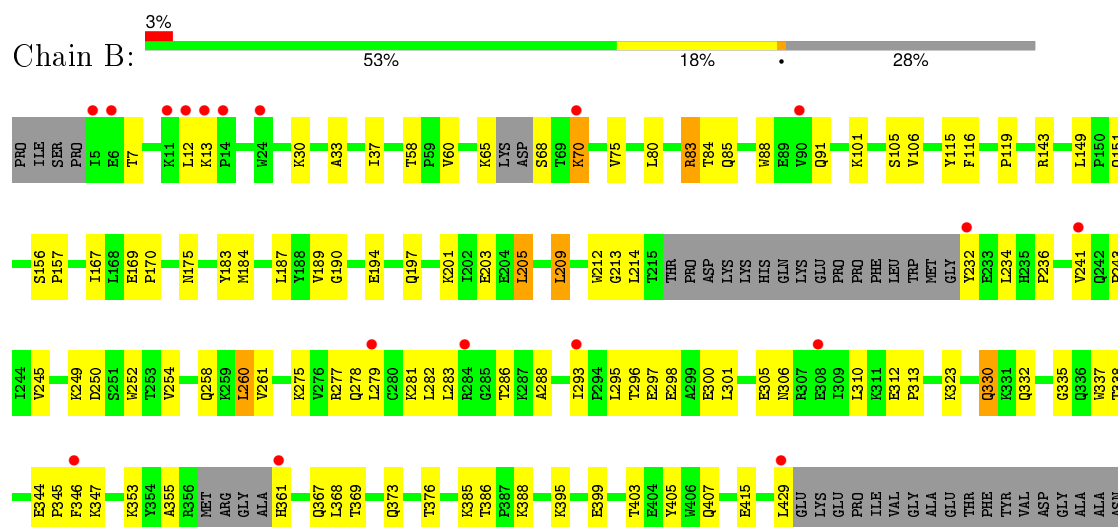
3 Residue-property plots [i](#)

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 Reverse transcriptase



• Molecule 1: HIV Reverse transcriptase



ARG	GLU	THR	LYS	LEU	GLY	LYS	ALA	GLY	TYR	VAL	THR	ASN	ARG	GLY	ARG	GLN	LYS	VAL	VAL	THR	LEU	THR	THR	ASP	THR	THR	ASN	GLN	LYS	THR	GLU	LEU	GLN	ALA	ALA	TYR	LEU	ALA	GLN	GLN	ASP	SER	GLY	LEU	GLU	VAL	ASN	ILE	VAL	THR	ASP	SER	SER	GLN	TYR	ALA	LEU	GLY	ILE	ILE	GLN
ALA	GLN	PRO	ASP	GLN	SER	GLU	SER	LEU	VAL	ASN	GLN	ILE	ILE	GLU	GLN	LEU	ILE	LYS	LYS	GLU	LYS	LYS	VAL	TYR	LEU	ALA	TRP	VAL	PRO	ALA	HIS	LYS	GLY	ILE	GLY	GLY	ASN	GLU	GLN	VAL	ASP	LYS	LEU	VAL	SER	ALA	GLY	ILE	ARG	LYS	VAL	LEU									

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	118.03Å 154.60Å 154.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.91 – 2.30 47.23 – 2.28	Depositor EDS
% Data completeness (in resolution range)	90.4 (29.91-2.30) 89.2 (47.23-2.28)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.27Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, R_{free}	0.223 , 0.274 0.221 , 0.272	Depositor DCC
R_{free} test set	2875 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	44.7	Xtriage
Anisotropy	0.355	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 62533 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8101	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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: KR1, SO4, CL

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	0.33	0/4625	0.49	0/6284
1	B	0.34	0/3426	0.49	0/4655
All	All	0.33	0/8051	0.49	0/10939

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	361	HIS	Peptide

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	4508	0	4558	133	0
1	B	3334	0	3365	75	0
2	A	31	0	20	0	0
3	A	20	0	0	0	0
4	B	1	0	0	0	0
5	A	110	0	0	3	0
5	B	97	0	0	4	0
All	All	8101	0	7943	204	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 13.

All (204) 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:131:THR:HG22	1:A:143:ARG:HG2	1.23	1.08
1:A:66:LYS:HG3	1:A:68:SER:H	1.40	0.87
1:B:296:THR:HG22	1:B:298:GLU:H	1.39	0.86
1:B:209:LEU:HD12	1:B:214:LEU:HD22	1.56	0.85
1:B:194:GLU:HG3	1:B:197:GLN:HG3	1.60	0.84
1:A:393:ILE:HB	1:A:423:VAL:HG13	1.59	0.84
1:B:58:THR:HG23	5:B:563:HOH:O	1.77	0.83
1:B:260:LEU:HD13	1:B:279:LEU:HD21	1.61	0.82
1:A:510:PRO:HG3	5:A:567:HOH:O	1.80	0.80
1:A:64:LYS:HE3	1:A:69:THR:HA	1.64	0.79
1:A:393:ILE:HB	1:A:423:VAL:CG1	2.13	0.79
1:B:297:GLU:H	1:B:297:GLU:CD	1.85	0.78
1:B:373:GLN:HE22	1:B:407:GLN:H	1.32	0.75
1:B:337:TRP:HE1	1:B:367:GLN:HE21	1.32	0.75
1:B:143:ARG:HD3	5:B:627:HOH:O	1.87	0.73
1:A:199:ARG:NH1	1:A:219:LYS:HE2	2.04	0.72
1:B:373:GLN:NE2	1:B:407:GLN:H	1.88	0.72
1:B:12:LEU:HD23	1:B:84:THR:HG22	1.73	0.71
1:A:219:LYS:HB3	1:A:219:LYS:NZ	2.06	0.70
1:A:131:THR:CG2	1:A:143:ARG:HH11	2.05	0.70
1:B:395:LYS:O	1:B:399:GLU:HG2	1.92	0.69
1:A:503:LEU:HD12	1:A:533:LEU:HD13	1.75	0.69
1:A:66:LYS:HD2	1:A:67:ASP:H	1.58	0.68
1:B:312:GLU:HB3	1:B:313:PRO:HD2	1.76	0.67
1:A:479:LEU:HB3	1:A:517:LEU:HD13	1.77	0.67
1:A:363:ASN:HA	1:A:511:ASP:OD1	1.94	0.67
1:A:194:GLU:H	1:A:194:GLU:CD	1.98	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:GLU:H	1:A:298:GLU:CD	1.98	0.66
1:B:376:THR:HG23	1:B:386:THR:HG23	1.78	0.66
1:A:131:THR:HG22	1:A:143:ARG:CG	2.13	0.65
1:B:403:THR:HG21	5:B:596:HOH:O	1.96	0.65
1:B:30:LYS:HD3	5:B:596:HOH:O	1.97	0.64
1:B:213:GLY:C	1:B:214:LEU:HD12	2.17	0.64
1:A:356:ARG:CZ	1:A:358:ARG:HD3	2.28	0.64
1:B:252:TRP:CD1	1:B:295:LEU:HD21	2.34	0.62
1:A:334:GLN:H	1:A:334:GLN:CD	2.02	0.62
1:A:27:THR:O	1:A:31:ILE:HG13	1.99	0.62
1:A:66:LYS:HE3	1:A:68:SER:N	2.14	0.62
1:B:115:TYR:HB3	1:B:149:LEU:HB2	1.82	0.61
1:A:57:ASN:OD1	1:A:131:THR:HG23	2.00	0.61
1:A:60:VAL:HG11	1:A:130:PHE:CD2	2.35	0.61
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.83	0.61
1:A:277:ARG:HB2	1:A:336:GLN:NE2	2.16	0.61
1:A:108:VAL:HG11	1:A:223:LYS:HB2	1.82	0.61
1:B:254:VAL:O	1:B:258:GLN:HG3	2.02	0.60
1:A:368:LEU:HD22	1:A:423:VAL:HG11	1.84	0.60
1:A:244:ILE:CD1	1:A:267:ALA:HB2	2.32	0.60
1:A:469:LEU:CD1	1:A:480:GLN:HG3	2.32	0.59
1:A:513:SER:H	1:A:519:ASN:HD21	1.49	0.58
1:A:287:LYS:HD2	1:A:293:ILE:HD11	1.84	0.58
1:B:60:VAL:HG12	1:B:75:VAL:HG22	1.85	0.58
1:A:510:PRO:HB2	1:A:522:ILE:HD11	1.85	0.58
1:A:209:LEU:HD23	1:A:216:THR:HG21	1.85	0.58
1:B:282:LEU:HB3	1:B:293:ILE:CD1	2.33	0.57
1:B:65:LYS:HB2	1:B:68:SER:HB2	1.85	0.57
1:A:451:LYS:HB3	1:A:471:ASP:HA	1.86	0.57
1:A:273:GLY:HA2	1:A:332:GLN:NE2	2.20	0.57
1:A:406:TRP:CE3	1:A:407:GLN:HG2	2.40	0.57
1:A:542:ILE:HD11	1:B:261:VAL:HG11	1.87	0.57
1:A:206:ARG:CZ	1:A:218:ASP:HA	2.34	0.57
1:B:296:THR:O	1:B:300:GLU:HG3	2.06	0.56
1:B:282:LEU:HB3	1:B:293:ILE:HD13	1.86	0.56
1:B:429:LEU:HD12	1:B:429:LEU:N	2.21	0.56
1:A:431:LYS:NZ	1:A:431:LYS:HB2	2.21	0.56
1:A:233:GLU:HG2	1:A:235:HIS:NE2	2.21	0.56
1:A:161:GLN:NE2	1:A:184:MET:O	2.38	0.56
1:A:393:ILE:CB	1:A:423:VAL:HG13	2.34	0.56
1:B:194:GLU:CG	1:B:197:GLN:HG3	2.34	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:509:GLN:N	1:A:510:PRO:HD3	2.22	0.55
1:A:469:LEU:HD13	1:A:480:GLN:HG3	1.89	0.55
1:A:254:VAL:O	1:A:258:GLN:HG3	2.08	0.54
1:A:450:THR:O	1:A:451:LYS:HB2	2.07	0.54
1:B:388:LYS:HE2	1:B:415:GLU:OE1	2.08	0.54
1:A:406:TRP:HE3	1:A:407:GLN:HE21	1.56	0.54
1:A:254:VAL:HB	1:A:289:LEU:HA	1.90	0.54
1:A:273:GLY:HA2	1:A:332:GLN:HE22	1.73	0.54
1:B:345:PRO:O	1:B:346:PHE:HB2	2.08	0.53
1:B:205:LEU:HD22	1:B:209:LEU:HD22	1.90	0.53
1:A:278:GLN:HG3	1:A:298:GLU:HB2	1.91	0.53
1:A:238:LYS:HD2	1:A:315:HIS:CG	2.44	0.53
1:A:277:ARG:NH1	1:A:334:GLN:HG3	2.24	0.53
1:A:66:LYS:HD2	1:A:67:ASP:N	2.24	0.52
1:B:373:GLN:HE22	1:B:407:GLN:N	2.05	0.52
1:B:85:GLN:HA	1:B:88:TRP:NE1	2.24	0.52
1:B:214:LEU:HD12	1:B:214:LEU:N	2.24	0.52
1:B:7:THR:HG22	1:B:119:PRO:HB2	1.92	0.52
1:A:460:ASN:ND2	1:B:288:ALA:HB2	2.25	0.52
1:A:275:LYS:HG3	5:A:639:HOH:O	2.10	0.52
1:A:233:GLU:HG2	1:A:235:HIS:HE2	1.74	0.52
1:A:542:ILE:HG23	1:B:283:LEU:HD13	1.91	0.51
1:A:228:LEU:HB3	1:A:242:GLN:OE1	2.11	0.51
1:B:65:LYS:CB	1:B:68:SER:HB2	2.40	0.51
1:A:342:TYR:HA	1:A:349:LEU:HD23	1.93	0.51
1:A:513:SER:N	1:A:519:ASN:HD21	2.08	0.51
1:A:503:LEU:CD1	1:A:533:LEU:HD13	2.41	0.50
1:A:356:ARG:NH1	1:A:358:ARG:HD3	2.26	0.50
1:A:392:PRO:O	1:A:423:VAL:HG12	2.11	0.50
1:B:175:ASN:HD21	1:B:201:LYS:NZ	2.10	0.50
1:A:518:VAL:O	1:A:522:ILE:HG13	2.12	0.49
1:B:330:GLN:NE2	1:B:338:THR:OG1	2.42	0.49
1:A:289:LEU:N	1:A:289:LEU:HD23	2.27	0.49
1:B:33:ALA:O	1:B:37:ILE:HG13	2.11	0.49
1:B:205:LEU:HD22	1:B:209:LEU:CD2	2.42	0.49
1:A:219:LYS:HB3	1:A:219:LYS:HZ2	1.74	0.49
1:A:464:GLN:O	1:A:465:LYS:HB2	2.12	0.49
1:A:548:VAL:O	1:A:552:VAL:HG22	2.12	0.49
1:B:183:TYR:CE2	1:B:184:MET:HG3	2.48	0.48
1:B:70:LYS:N	1:B:70:LYS:HD2	2.29	0.48
1:B:13:LYS:HG3	1:B:83:ARG:O	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:GLU:HB3	1:A:170:PRO:HD3	1.95	0.48
1:A:206:ARG:NH1	1:A:218:ASP:HA	2.28	0.48
1:A:194:GLU:N	1:A:194:GLU:CD	2.65	0.48
1:B:169:GLU:HB3	1:B:170:PRO:HD3	1.95	0.48
1:B:296:THR:HG22	1:B:298:GLU:N	2.18	0.48
1:B:277:ARG:O	1:B:281:LYS:HG3	2.13	0.48
1:B:116:PHE:CZ	1:B:151:GLN:NE2	2.82	0.48
1:B:344:GLU:HB2	1:B:347:LYS:HD3	1.96	0.48
1:B:209:LEU:HB3	1:B:214:LEU:HB2	1.96	0.47
1:A:199:ARG:O	1:A:203:GLU:HG2	2.14	0.47
1:A:167:ILE:O	1:A:170:PRO:HD2	2.13	0.47
1:A:8:VAL:O	1:A:10:VAL:HG23	2.14	0.47
1:A:52:PRO:HA	1:A:54:ASN:N	2.29	0.47
1:A:552:VAL:O	1:A:552:VAL:HG23	2.14	0.47
1:A:32:LYS:O	1:A:36:GLU:HG3	2.14	0.47
1:A:40:GLU:HA	1:A:40:GLU:OE2	2.15	0.47
1:B:353:LYS:HD3	1:B:429:LEU:CD2	2.44	0.47
1:B:156:SER:N	1:B:157:PRO:HD2	2.30	0.47
1:A:131:THR:CG2	1:A:143:ARG:NH1	2.75	0.46
1:A:536:VAL:HG13	1:A:537:PRO:HD2	1.98	0.46
1:A:66:LYS:HE3	1:A:68:SER:HB2	1.97	0.46
1:B:241:VAL:O	1:B:243:PRO:HD3	2.16	0.46
1:A:450:THR:CG2	1:A:452:LEU:HB2	2.46	0.46
1:A:289:LEU:HD23	1:A:289:LEU:H	1.79	0.46
1:B:346:PHE:CD1	1:B:346:PHE:N	2.84	0.46
1:B:105:SER:O	1:B:190:GLY:HA2	2.16	0.46
1:B:429:LEU:H	1:B:429:LEU:HD12	1.79	0.46
1:A:306:ASN:O	1:A:310:LEU:HD22	2.16	0.46
1:B:346:PHE:HD1	1:B:346:PHE:N	2.14	0.46
1:A:342:TYR:HB3	1:A:348:ASN:HD22	1.81	0.46
1:A:245:VAL:HG13	1:A:245:VAL:O	2.15	0.46
1:B:369:THR:HG21	1:B:405:TYR:HB2	1.97	0.45
1:B:101:LYS:O	1:B:236:PRO:HB2	2.17	0.45
1:A:286:THR:O	1:A:286:THR:HG23	2.17	0.45
1:A:139:THR:HB	1:A:140:PRO:HD2	1.99	0.45
1:A:126:LYS:HA	1:A:145:GLN:NE2	2.32	0.45
1:A:65:LYS:O	1:A:66:LYS:HG3	2.17	0.45
1:A:502:ALA:O	1:A:506:ILE:HG12	2.17	0.45
1:A:31:ILE:O	1:A:35:VAL:HG23	2.17	0.45
1:A:21:VAL:HB	1:A:59:PRO:HD3	1.98	0.45
1:A:252:TRP:O	1:A:292:VAL:HA	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:TRP:CE3	1:A:403:THR:HG23	2.53	0.44
1:A:447:ASN:HB3	1:A:450:THR:HB	1.99	0.44
1:A:355:ALA:O	1:A:356:ARG:C	2.56	0.44
1:A:289:LEU:CD2	1:A:289:LEU:H	2.31	0.44
1:A:552:VAL:O	1:A:553:SER:HB3	2.18	0.44
1:A:296:THR:HB	1:A:298:GLU:OE1	2.18	0.44
1:A:406:TRP:CE3	1:A:407:GLN:NE2	2.86	0.44
1:B:116:PHE:HZ	1:B:151:GLN:NE2	2.16	0.44
1:A:420:PRO:HA	1:A:421:PRO:C	2.37	0.44
1:A:469:LEU:HD11	1:A:480:GLN:HG3	1.99	0.44
1:A:66:LYS:CG	1:A:68:SER:H	2.22	0.43
1:A:482:ILE:HD12	1:A:502:ALA:HB1	2.00	0.43
1:A:21:VAL:HG12	1:A:22:LYS:N	2.32	0.43
1:A:353:LYS:HB3	1:A:353:LYS:HE3	1.59	0.43
1:A:196:GLY:HA3	5:A:618:HOH:O	2.18	0.43
1:A:379:SER:OG	1:A:387:PRO:HD3	2.17	0.43
1:B:115:TYR:HB3	1:B:149:LEU:CB	2.49	0.43
1:A:66:LYS:CG	1:A:67:ASP:N	2.82	0.43
1:A:345:PRO:HA	1:A:346:PHE:HA	1.59	0.42
1:A:443:ASP:OD2	1:A:444:GLY:N	2.51	0.42
1:A:288:ALA:HB3	1:A:291:GLU:HG3	2.01	0.42
1:A:235:HIS:HB3	1:A:236:PRO:CD	2.50	0.42
1:A:544:GLY:HA2	1:B:286:THR:HG22	2.02	0.42
1:A:104:LYS:HB3	1:A:192:ASP:HA	2.01	0.42
1:B:295:LEU:CD2	1:B:295:LEU:N	2.82	0.42
1:B:243:PRO:O	1:B:245:VAL:HG13	2.19	0.42
1:A:108:VAL:HG13	1:A:108:VAL:O	2.20	0.42
1:A:450:THR:O	1:A:451:LYS:CB	2.68	0.42
1:B:278:GLN:HG3	1:B:298:GLU:HB3	2.00	0.41
1:A:339:TYR:CZ	1:A:352:GLY:HA3	2.55	0.41
1:A:260:LEU:O	1:A:264:LEU:HD22	2.20	0.41
1:A:431:LYS:HZ3	1:A:431:LYS:HB2	1.85	0.41
1:A:283:LEU:HA	1:A:283:LEU:HD12	1.82	0.41
1:A:244:ILE:HD12	1:A:267:ALA:HB2	2.02	0.41
1:B:332:GLN:HG3	1:B:338:THR:HG23	2.02	0.41
1:A:118:VAL:HG13	1:A:119:PRO:HD2	2.03	0.41
1:A:102:LYS:HB3	1:A:102:LYS:HE2	1.92	0.41
1:B:295:LEU:HD22	1:B:295:LEU:N	2.36	0.41
1:A:542:ILE:O	1:A:543:GLY:C	2.58	0.41
1:B:330:GLN:HB2	1:B:330:GLN:HE21	1.75	0.41
1:A:453:GLY:HA3	1:A:472:THR:HG21	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:VAL:HA	1:B:189:VAL:O	2.21	0.41
1:A:244:ILE:HD13	1:A:267:ALA:HB2	2.00	0.41
1:B:282:LEU:HB3	1:B:293:ILE:HD11	2.03	0.41
1:B:275:LYS:H	1:B:306:ASN:HD21	1.69	0.40
1:B:335:GLY:O	1:B:355:ALA:HA	2.22	0.40
1:B:301:LEU:O	1:B:305:GLU:HG3	2.22	0.40
1:A:305:GLU:O	1:A:309:ILE:HG13	2.21	0.40
1:A:19:PRO:HG3	1:A:80:LEU:HB2	2.03	0.40
1:A:64:LYS:CE	1:A:69:THR:HA	2.44	0.40
1:A:235:HIS:HB3	1:A:236:PRO:HD2	2.03	0.40
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.56	0.40
1:B:167:ILE:HG12	1:B:212:TRP:CD2	2.56	0.40
1:A:246:LEU:HD22	1:A:260:LEU:HD12	2.02	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	551/560 (98%)	527 (96%)	19 (3%)	5 (1%)	21	24
1	B	395/560 (70%)	381 (96%)	14 (4%)	0	100	100
All	All	946/1120 (84%)	908 (96%)	33 (4%)	5 (0%)	34	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	356	ARG
1	A	543	GLY
1	A	66	LYS
1	A	465	LYS

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	495/500 (99%)	457 (92%)	38 (8%)	16	20
1	B	367/500 (73%)	349 (95%)	18 (5%)	31	41
All	All	862/1000 (86%)	806 (94%)	56 (6%)	21	27

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	LYS
1	A	74	LEU
1	A	78	ARG
1	A	113	ASP
1	A	120	LEU
1	A	194	GLU
1	A	209	LEU
1	A	248	GLU
1	A	264	LEU
1	A	282	LEU
1	A	283	LEU
1	A	295	LEU
1	A	298	GLU
1	A	303	LEU
1	A	308	GLU
1	A	310	LEU
1	A	311	LYS
1	A	334	GLN
1	A	344	GLU
1	A	349	LEU
1	A	350	LYS
1	A	353	LYS
1	A	357	MET
1	A	361	HIS
1	A	373	GLN
1	A	399	GLU
1	A	410	TRP

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Mol	Chain	Res	Type
1	A	428	GLN
1	A	448	ARG
1	A	471	ASP
1	A	479	LEU
1	A	500	GLN
1	A	507	GLN
1	A	517	LEU
1	A	519	ASN
1	A	529	GLU
1	A	533	LEU
1	A	551	LEU
1	B	70	LYS
1	B	80	LEU
1	B	83	ARG
1	B	91	GLN
1	B	187	LEU
1	B	203	GLU
1	B	205	LEU
1	B	209	LEU
1	B	232	TYR
1	B	234	LEU
1	B	249	LYS
1	B	250	ASP
1	B	260	LEU
1	B	310	LEU
1	B	323	LYS
1	B	330	GLN
1	B	368	LEU
1	B	385	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	GLN
1	A	258	GLN
1	A	332	GLN
1	A	336	GLN
1	A	348	ASN
1	A	407	GLN
1	A	428	GLN
1	A	487	GLN
1	A	519	ASN

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Mol	Chain	Res	Type
1	A	524	GLN
1	B	85	GLN
1	B	91	GLN
1	B	151	GLN
1	B	175	ASN
1	B	182	GLN
1	B	197	GLN
1	B	258	GLN
1	B	306	ASN
1	B	330	GLN
1	B	336	GLN
1	B	340	GLN
1	B	348	ASN
1	B	367	GLN
1	B	373	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 for Mogul analysis.

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	SO4	A	561	-	4,4,4	0.19	0	6,6,6	0.23	0
3	SO4	A	562	-	4,4,4	0.22	0	6,6,6	0.16	0
3	SO4	A	563	-	4,4,4	0.28	0	6,6,6	0.09	0
3	SO4	A	564	-	4,4,4	0.26	0	6,6,6	0.08	0
2	KR1	A	701	-	29,33,33	3.01	15 (51%)	36,48,48	3.26	12 (33%)

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	SO4	A	561	-	-	0/0/0/0	0/0/0/0
3	SO4	A	562	-	-	0/0/0/0	0/0/0/0
3	SO4	A	563	-	-	0/0/0/0	0/0/0/0
3	SO4	A	564	-	-	0/0/0/0	0/0/0/0
2	KR1	A	701	-	-	0/18/18/18	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	KR1	C9-C8	-2.49	1.46	1.50
2	A	701	KR1	O2-C12	2.02	1.29	1.24
2	A	701	KR1	C2-C7	2.36	1.43	1.38
2	A	701	KR1	C15-C14	2.57	1.59	1.52
2	A	701	KR1	C1-N1	2.87	1.21	1.14
2	A	701	KR1	F1-C21	2.97	1.40	1.35
2	A	701	KR1	O1-C8	3.65	1.29	1.22
2	A	701	KR1	C2-C3	3.88	1.46	1.39
2	A	701	KR1	C21-N5	4.09	1.39	1.32
2	A	701	KR1	C22-C18	4.09	1.46	1.39
2	A	701	KR1	C4-C5	4.28	1.45	1.39
2	A	701	KR1	C12-N2	5.26	1.42	1.33
2	A	701	KR1	C9-N3	5.58	1.47	1.35
2	A	701	KR1	C6-C5	5.85	1.48	1.39
2	A	701	KR1	C19-C18	6.05	1.50	1.39

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	KR1	O1-C8-C9	-6.67	112.53	120.21
2	A	701	KR1	C11-C12-N2	-5.03	117.14	125.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	KR1	C22-C21-N5	-4.32	120.62	127.07
2	A	701	KR1	C17-N3-C13	-4.19	113.10	118.05
2	A	701	KR1	C21-C22-C18	-3.91	117.02	119.49
2	A	701	KR1	C5-C4-C3	-2.83	117.91	120.46
2	A	701	KR1	C9-C11-C14	3.18	125.84	122.34
2	A	701	KR1	F1-C21-N5	3.70	120.00	114.33
2	A	701	KR1	C17-N3-C9	4.00	127.52	119.48
2	A	701	KR1	C21-N5-C20	4.10	125.65	117.61
2	A	701	KR1	C5-C8-C9	5.79	126.20	119.53
2	A	701	KR1	C12-N2-C13	12.27	125.85	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	553/560 (98%)	0.14	15 (2%) 58 67	30, 49, 82, 101	0
1	B	403/560 (71%)	0.22	18 (4%) 37 46	26, 47, 82, 95	0
All	All	956/1120 (85%)	0.17	33 (3%) 48 56	26, 48, 82, 101	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	14	PRO	5.9
1	B	232	TYR	4.5
1	B	279	LEU	4.0
1	A	452	LEU	4.0
1	B	13	LYS	3.8
1	B	346	PHE	3.4
1	B	11	LYS	3.3
1	A	66	LYS	3.2
1	A	67	ASP	3.0
1	B	308	GLU	2.8
1	A	551	LEU	2.7
1	B	6	GLU	2.7
1	A	553	SER	2.6
1	B	284	ARG	2.6
1	A	193	LEU	2.6
1	A	402	TRP	2.5
1	B	361	HIS	2.5
1	B	24	TRP	2.5
1	A	53	GLU	2.4
1	B	5	ILE	2.4
1	B	90	VAL	2.3
1	A	426	TRP	2.3
1	B	293	ILE	2.2
1	B	241	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	552	VAL	2.2
1	B	429	LEU	2.2
1	B	70	LYS	2.1
1	A	195	ILE	2.1
1	A	453	GLY	2.0
1	A	491	LEU	2.0
1	A	406	TRP	2.0
1	A	291	GLU	2.0
1	B	12	LEU	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(Å ²)	Q<0.9
3	SO4	A	563	5/5	0.83	0.31	5.05	120,120,120,120	0
3	SO4	A	561	5/5	0.96	0.17	1.13	61,65,67,68	0
2	KR1	A	701	31/31	0.94	0.15	0.45	34,41,44,48	0
3	SO4	A	562	5/5	0.96	0.12	-0.39	58,60,61,64	0
4	CL	B	561	1/1	0.95	0.24	-	62,62,62,62	0
3	SO4	A	564	5/5	0.73	0.16	-	107,107,108,108	0

6.5 Other polymers [i](#)

There are no such residues in this entry.