



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:47 AM UTC

PDB ID : 3LAL / pdb_00003lal
Title : Crystal structure of HIV-1 reverse transcriptase in complex with N1-ethyl pyrimidinedione non-nucleoside inhibitor
Authors : Lansdon, E.B.; Mitchell, M.L.
Deposited on : 2010-01-06
Resolution : 2.51 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

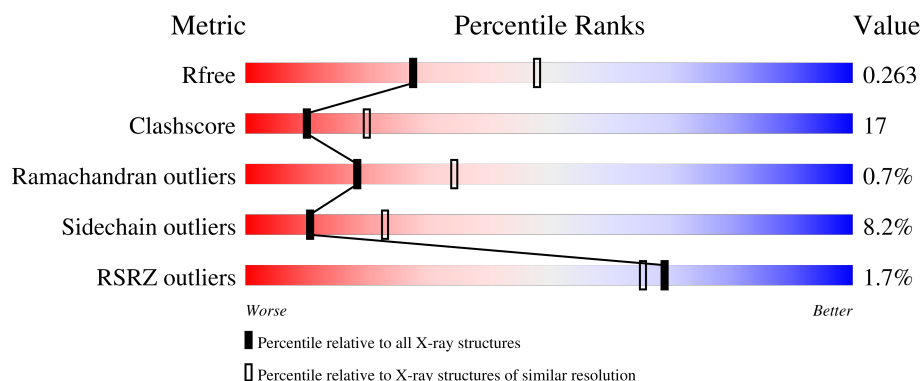
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	560	
1	B	560	

2 Entry composition [i](#)

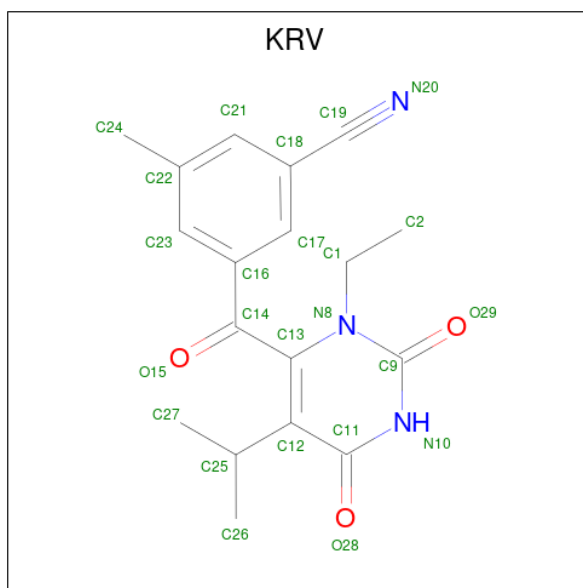
There are 4 unique types of molecules in this entry. The entry contains 8013 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Reverse transcriptase.

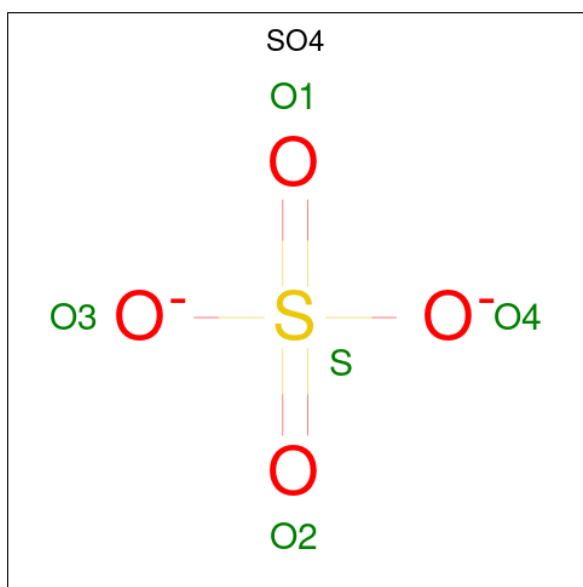
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	553	Total	C	N	O	S	0	0	0
			4508	2914	752	834	8			
1	B	403	Total	C	N	O	S	0	0	0
			3334	2173	550	605	6			

- Molecule 2 is 3-{{[3-ethyl-5-(1-methylethyl)-2,6-dioxo-1,2,3,6-tetrahydropyrimidin-4-yl]carbonyl}-5-methylbenzonitrile (CCD ID: KRV) (formula: C₁₈H₁₉N₃O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			24	18	3	3		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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		
3	B	1	Total	O	S	0	0
			5	4	1		

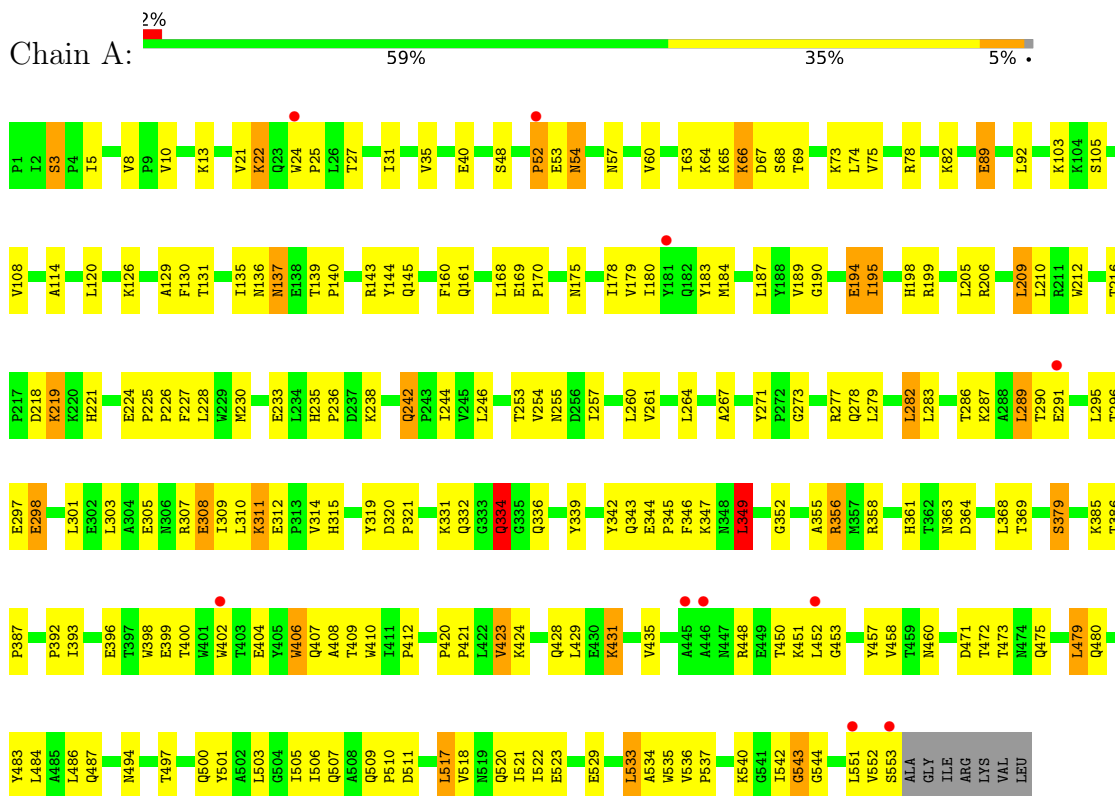
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	69	Total	O	0	0
			69	69		
4	B	53	Total	O	0	0
			53	53		

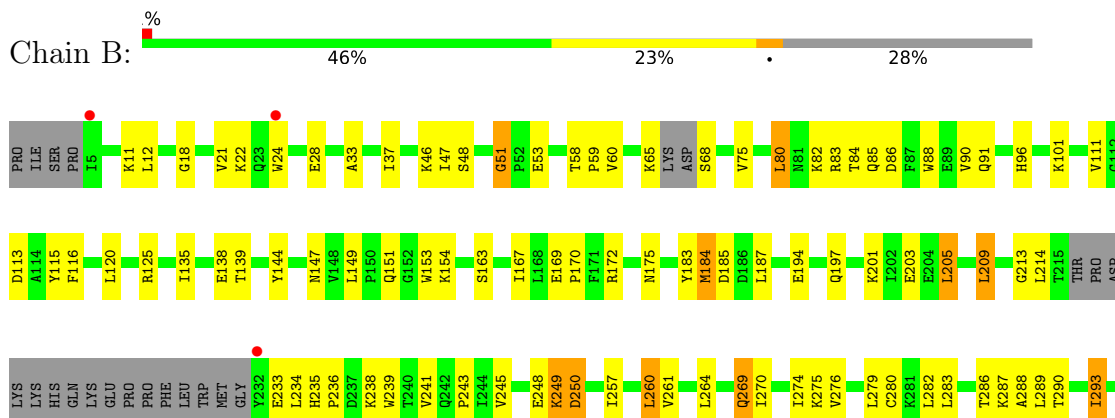
3 Residue-property plots

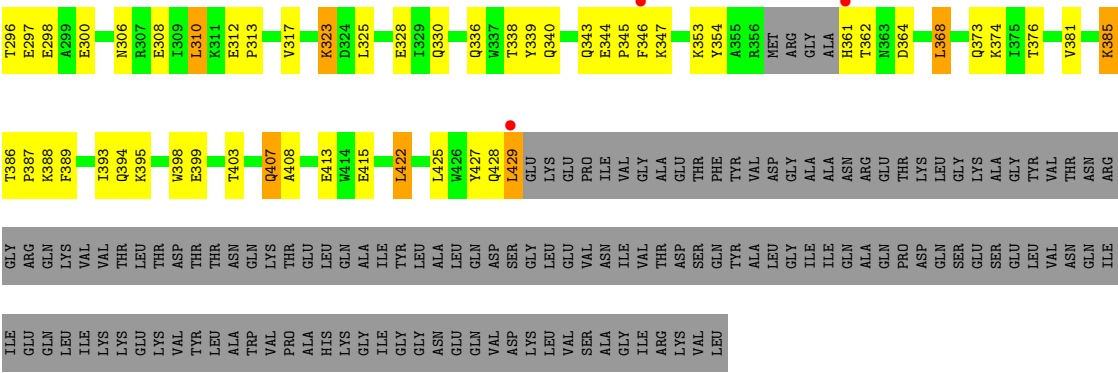
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: HIV Reverse transcriptase



• Molecule 1: HIV 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	118.15Å 154.12Å 153.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.83 – 2.51 29.83 – 2.51	Depositor EDS
% Data completeness (in resolution range)	87.8 (29.83-2.51) 87.9 (29.83-2.51)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.19 (at 2.51Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, R_{free}	0.218 , 0.267 0.209 , 0.263	Depositor DCC
R_{free} test set	2112 reflections (4.62%)	wwPDB-VP
Wilson B-factor (Å ²)	58.8	Xtriage
Anisotropy	0.359	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8013	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 3.74% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, KRV

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	0.45	0/4625	0.80	9/6284 (0.1%)
1	B	0.46	0/3426	0.81	8/4655 (0.2%)
All	All	0.45	0/8051	0.81	17/10939 (0.2%)

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	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	54	ASN	CA-C-N	7.03	127.63	120.04
1	A	54	ASN	C-N-CA	7.03	127.63	120.04
1	A	3	SER	CA-C-N	6.23	127.04	120.12
1	A	3	SER	C-N-CA	6.23	127.04	120.12
1	B	18	GLY	CA-C-N	6.22	125.98	119.76
1	B	18	GLY	C-N-CA	6.22	125.98	119.76
1	B	51	GLY	CA-C-N	6.09	126.20	119.87
1	B	51	GLY	C-N-CA	6.09	126.20	119.87
1	A	224	GLU	CA-C-N	5.88	123.95	119.66
1	A	224	GLU	C-N-CA	5.88	123.95	119.66
1	B	233	GLU	N-CA-C	5.43	117.58	108.73
1	A	386	THR	CA-C-N	5.42	125.40	120.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	386	THR	C-N-CA	5.42	125.40	120.03
1	B	235	HIS	CA-C-N	5.28	125.08	119.28
1	B	235	HIS	C-N-CA	5.28	125.08	119.28
1	B	362	THR	N-CA-C	-5.21	105.17	112.25
1	A	349	LEU	N-CA-C	-5.04	105.70	111.14

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	52	PRO	Peptide
1	B	361	HIS	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4508	0	4558	174	0
1	B	3334	0	3365	112	0
2	A	24	0	19	0	0
3	A	20	0	0	0	0
3	B	5	0	0	0	0
4	A	69	0	0	3	0
4	B	53	0	0	4	0
All	All	8013	0	7942	274	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 17.

All (274) 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	1.24	1.10
1:B:296:THR:HG22	1:B:298:GLU:H	1.28	0.95
1:A:510:PRO:HG3	4:A:568:HOH:O	1.68	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:373:GLN:HE22	1:B:407:GLN:H	1.17	0.90
1:A:393:ILE:HB	1:A:423:VAL:HG13	1.54	0.89
1:B:58:THR:HG23	4:B:563:HOH:O	1.75	0.86
1:A:131:THR:HG22	1:A:143:ARG:CG	2.05	0.85
1:A:65:LYS:O	1:A:66:LYS:HG3	1.76	0.85
1:A:131:THR:CG2	1:A:143:ARG:HH11	1.91	0.84
1:A:194:GLU:CD	1:A:194:GLU:H	1.87	0.83
1:A:537:PRO:HB2	1:A:540:LYS:HG3	1.64	0.80
1:A:60:VAL:HG12	1:A:75:VAL:HG22	1.64	0.80
1:B:260:LEU:HD13	1:B:279:LEU:HD21	1.64	0.79
1:B:376:THR:HG23	1:B:386:THR:HG23	1.68	0.76
1:B:65:LYS:HG3	1:B:68:SER:HB2	1.67	0.75
1:A:228:LEU:HB3	1:A:242:GLN:NE2	2.03	0.73
1:B:138:GLU:HG2	1:B:139:THR:HG23	1.69	0.73
1:A:297:GLU:O	1:A:301:LEU:HG	1.89	0.73
1:A:393:ILE:HB	1:A:423:VAL:CG1	2.21	0.70
1:A:277:ARG:HB2	1:A:336:GLN:NE2	2.07	0.70
1:B:373:GLN:NE2	1:B:407:GLN:H	1.88	0.70
1:B:297:GLU:H	1:B:297:GLU:CD	2.00	0.69
1:B:395:LYS:O	1:B:399:GLU:HG2	1.92	0.69
1:A:114:ALA:HB1	1:A:160:PHE:CE1	2.28	0.69
1:A:199:ARG:HH11	1:A:219:LYS:HE2	1.58	0.68
1:A:298:GLU:H	1:A:298:GLU:CD	1.99	0.68
1:B:241:VAL:HG23	1:B:243:PRO:HD3	1.75	0.68
1:A:66:LYS:HD3	1:A:67:ASP:H	1.59	0.68
1:A:278:GLN:HG3	1:A:298:GLU:HB2	1.75	0.68
1:A:520:GLN:O	1:A:523:GLU:HG2	1.94	0.68
1:A:277:ARG:NH1	1:A:334:GLN:HG3	2.10	0.67
1:B:393:ILE:HD13	1:B:398:TRP:HB2	1.77	0.66
1:B:194:GLU:HG3	1:B:197:GLN:HG3	1.76	0.66
1:B:345:PRO:O	1:B:346:PHE:HB2	1.94	0.66
1:B:175:ASN:HD21	1:B:201:LYS:HZ3	1.41	0.66
1:A:542:ILE:HD11	1:B:261:VAL:HG11	1.77	0.66
1:B:24:TRP:HZ3	1:B:59:PRO:CG	2.07	0.66
1:A:27:THR:O	1:A:31:ILE:HG13	1.95	0.66
1:A:273:GLY:HA2	1:A:332:GLN:NE2	2.12	0.65
1:A:343:GLN:HG3	1:A:349:LEU:HD21	1.78	0.65
1:A:368:LEU:HD22	1:A:423:VAL:HG11	1.79	0.65
1:B:24:TRP:HZ3	1:B:59:PRO:HG2	1.60	0.65
1:A:334:GLN:CD	1:A:334:GLN:H	2.03	0.65
1:B:60:VAL:HG12	1:B:75:VAL:HG22	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:480:GLN:HG2	1:A:517:LEU:HD11	1.79	0.64
1:A:199:ARG:NH1	1:A:219:LYS:HE2	2.12	0.64
1:A:305:GLU:O	1:A:309:ILE:HG13	1.98	0.64
1:B:205:LEU:HD22	1:B:209:LEU:HD22	1.80	0.64
1:B:33:ALA:O	1:B:37:ILE:HG13	1.98	0.63
1:A:230:MET:HE2	1:A:230:MET:H	1.63	0.63
1:A:356:ARG:CZ	1:A:358:ARG:HD3	2.29	0.63
1:B:86:ASP:O	1:B:90:VAL:HG22	1.99	0.62
1:A:408:ALA:HB1	1:B:364:ASP:HB3	1.80	0.62
1:A:257:ILE:O	1:A:261:VAL:HG23	2.00	0.62
1:A:544:GLY:HA2	1:B:286:THR:HG22	1.81	0.61
1:A:175:ASN:HB3	1:A:178:ILE:HD12	1.82	0.61
1:B:175:ASN:HD21	1:B:201:LYS:NZ	1.98	0.61
1:A:409:THR:O	1:B:364:ASP:HB2	2.01	0.61
1:A:406:TRP:HE3	1:A:407:GLN:HE21	1.49	0.60
1:B:425:LEU:O	1:B:428:GLN:HG2	2.01	0.60
1:A:503:LEU:HD12	1:A:533:LEU:HD13	1.83	0.60
1:B:209:LEU:HD12	1:B:214:LEU:HD22	1.85	0.59
1:A:206:ARG:NH1	1:A:218:ASP:HA	2.17	0.59
1:A:379:SER:OG	1:A:387:PRO:HD3	2.02	0.58
1:B:312:GLU:HB3	1:B:313:PRO:HD2	1.85	0.58
1:A:168:LEU:HD11	1:A:187:LEU:HD21	1.86	0.58
1:A:238:LYS:HE2	1:A:315:HIS:CG	2.38	0.58
1:A:479:LEU:HB3	1:A:517:LEU:HD13	1.86	0.58
1:B:249:LYS:HG2	1:B:250:ASP:H	1.66	0.58
1:A:342:TYR:HA	1:A:349:LEU:HD23	1.86	0.58
1:A:254:VAL:HB	1:A:289:LEU:HA	1.86	0.58
1:B:296:THR:HG22	1:B:298:GLU:N	2.10	0.57
1:A:60:VAL:HG11	1:A:130:PHE:CD2	2.40	0.57
1:A:287:LYS:HB3	1:A:291:GLU:HG3	1.86	0.57
1:A:233:GLU:HG2	1:A:235:HIS:NE2	2.20	0.56
1:B:163:SER:O	1:B:167:ILE:HG13	2.04	0.56
1:A:65:LYS:HE2	1:A:66:LYS:HG2	1.88	0.56
1:A:358:ARG:NH2	1:B:394:GLN:HG3	2.20	0.56
1:B:169:GLU:HB3	1:B:170:PRO:HD3	1.87	0.56
1:B:317:VAL:HG22	1:B:347:LYS:HB3	1.87	0.56
1:A:228:LEU:HB3	1:A:242:GLN:HE22	1.69	0.55
1:A:244:ILE:CD1	1:A:267:ALA:HB2	2.37	0.55
1:A:89:GLU:HG2	1:A:92:LEU:HD21	1.88	0.55
1:B:213:GLY:C	1:B:214:LEU:HD12	2.31	0.55
1:A:64:LYS:NZ	1:A:69:THR:HA	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:483:TYR:O	1:A:487:GLN:HG2	2.07	0.55
1:B:51:GLY:HA3	1:B:53:GLU:OE2	2.07	0.55
1:A:40:GLU:OE2	1:A:40:GLU:HA	2.06	0.55
1:A:406:TRP:CE3	1:A:407:GLN:HG2	2.42	0.54
1:A:57:ASN:OD1	1:A:131:THR:HG23	2.08	0.54
1:A:219:LYS:HB3	1:A:219:LYS:HZ2	1.72	0.54
1:B:243:PRO:O	1:B:245:VAL:HG13	2.08	0.53
1:A:356:ARG:NH2	1:A:358:ARG:HD3	2.23	0.53
1:A:358:ARG:HH21	1:B:394:GLN:HG3	1.73	0.53
1:B:274:ILE:HA	1:B:306:ASN:HD21	1.72	0.53
1:A:319:TYR:OH	1:A:385:LYS:HE2	2.08	0.53
1:B:282:LEU:HB3	1:B:293:ILE:CD1	2.38	0.53
1:A:131:THR:HG23	1:A:143:ARG:HH11	1.73	0.53
1:A:194:GLU:O	1:A:195:ILE:C	2.51	0.53
1:A:169:GLU:HB3	1:A:170:PRO:HD3	1.90	0.53
1:A:420:PRO:HA	1:A:421:PRO:C	2.34	0.53
1:A:228:LEU:HD12	1:A:228:LEU:N	2.24	0.53
1:A:369:THR:HG1	1:A:398:TRP:HZ3	1.56	0.53
1:B:151:GLN:HB3	1:B:185:ASP:OD2	2.08	0.53
1:A:31:ILE:O	1:A:35:VAL:HG23	2.07	0.52
1:B:46:LYS:HE2	1:B:116:PHE:CD1	2.43	0.52
1:B:429:LEU:N	1:B:429:LEU:HD12	2.23	0.52
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.09	0.52
1:A:228:LEU:CG	1:A:242:GLN:HE21	2.22	0.52
1:B:28:GLU:HB2	1:B:135:ILE:HD11	1.91	0.52
1:B:101:LYS:HG2	4:B:574:HOH:O	2.09	0.52
1:A:246:LEU:HD12	1:A:307:ARG:HD3	1.92	0.52
1:A:450:THR:HG22	1:A:452:LEU:HB2	1.91	0.52
1:A:65:LYS:C	1:A:66:LYS:HG3	2.35	0.52
1:A:453:GLY:HA3	1:A:472:THR:HG21	1.92	0.52
1:B:282:LEU:HB3	1:B:293:ILE:HD13	1.90	0.52
1:A:289:LEU:CD2	1:A:289:LEU:H	2.21	0.52
1:A:279:LEU:O	1:A:282:LEU:HB2	2.10	0.52
1:B:306:ASN:O	1:B:310:LEU:HD22	2.10	0.51
1:A:273:GLY:HA2	1:A:332:GLN:HE22	1.75	0.51
1:A:536:VAL:HG13	1:A:537:PRO:HD2	1.92	0.51
1:A:66:LYS:HD3	1:A:68:SER:H	1.76	0.51
1:A:225:PRO:HG3	1:A:227:PHE:CE2	2.46	0.51
1:A:355:ALA:O	1:A:356:ARG:C	2.53	0.51
1:A:331:LYS:HE3	1:A:364:ASP:OD2	2.11	0.51
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:542:ILE:O	1:A:543:GLY:O	2.28	0.51
1:B:113:ASP:O	1:B:116:PHE:HD2	1.94	0.51
1:A:392:PRO:O	1:A:423:VAL:HG12	2.11	0.51
1:B:236:PRO:O	1:B:239:TRP:HB2	2.10	0.51
1:A:225:PRO:HG3	1:A:227:PHE:CZ	2.46	0.50
1:A:450:THR:CG2	1:A:452:LEU:HB2	2.41	0.50
1:A:13:LYS:HE3	1:A:82:LYS:O	2.12	0.50
1:A:105:SER:O	1:A:190:GLY:HA2	2.12	0.50
1:A:542:ILE:HD13	1:B:280:CYS:HB2	1.92	0.50
1:B:115:TYR:HB3	1:B:149:LEU:HB2	1.93	0.50
1:B:323:LYS:HB3	1:B:343:GLN:NE2	2.27	0.49
1:B:111:VAL:HG11	1:B:187:LEU:HD22	1.94	0.49
1:A:219:LYS:HB3	1:A:219:LYS:NZ	2.27	0.49
1:A:369:THR:OG1	1:A:398:TRP:HZ3	1.96	0.49
1:A:509:GLN:N	1:A:510:PRO:HD3	2.26	0.49
1:B:183:TYR:CE2	1:B:184:MET:HG3	2.47	0.49
1:A:400:THR:O	1:A:404:GLU:HG2	2.12	0.49
1:A:542:ILE:O	1:A:543:GLY:C	2.56	0.49
1:B:205:LEU:HD22	1:B:209:LEU:CD2	2.42	0.49
1:A:289:LEU:HD23	1:A:289:LEU:N	2.28	0.49
1:B:274:ILE:HA	1:B:306:ASN:ND2	2.27	0.48
1:B:373:GLN:NE2	1:B:408:ALA:H	2.11	0.48
1:A:161:GLN:NE2	1:A:184:MET:O	2.43	0.48
1:B:373:GLN:HE22	1:B:407:GLN:N	1.99	0.48
1:A:552:VAL:HG23	1:A:552:VAL:O	2.13	0.48
1:B:194:GLU:H	1:B:197:GLN:HE21	1.62	0.48
1:A:194:GLU:CD	1:A:194:GLU:N	2.61	0.48
1:B:85:GLN:HA	1:B:88:TRP:NE1	2.28	0.48
1:A:298:GLU:CD	1:A:298:GLU:N	2.71	0.48
1:B:28:GLU:CB	1:B:135:ILE:HD11	2.44	0.48
1:B:387:PRO:HG2	1:B:389:PHE:CE1	2.49	0.47
1:A:450:THR:O	1:A:451:LYS:HB2	2.14	0.47
1:A:451:LYS:HB3	1:A:471:ASP:HA	1.96	0.47
1:B:260:LEU:HD22	1:B:264:LEU:CD1	2.45	0.47
1:A:21:VAL:O	1:A:57:ASN:HB3	2.15	0.47
1:A:289:LEU:HD23	1:A:290:THR:H	1.78	0.47
1:B:422:LEU:HD12	1:B:422:LEU:HA	1.73	0.47
1:B:429:LEU:HD12	1:B:429:LEU:H	1.79	0.47
1:A:183:TYR:CE2	1:A:184:MET:HG3	2.50	0.47
1:A:497:THR:O	1:A:535:TRP:HA	2.15	0.47
1:A:282:LEU:HD11	1:A:296:THR:HG23	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:HIS:HE1	1:B:381:VAL:O	1.98	0.47
1:A:255:ASN:HB2	1:A:289:LEU:HB2	1.97	0.46
1:A:345:PRO:HA	1:A:346:PHE:HA	1.58	0.46
1:B:340:GLN:NE2	4:B:580:HOH:O	2.42	0.46
1:A:228:LEU:HG	1:A:242:GLN:HE21	1.81	0.46
1:A:501:TYR:CE1	1:A:505:ILE:HD11	2.50	0.46
1:B:276:VAL:HG12	1:B:276:VAL:O	2.15	0.46
1:A:289:LEU:CD2	1:A:289:LEU:N	2.78	0.46
1:A:52:PRO:HA	1:A:54:ASN:N	2.30	0.46
1:A:435:VAL:HA	1:B:290:THR:HG21	1.97	0.46
1:A:542:ILE:HG23	1:B:283:LEU:HD13	1.97	0.46
1:A:66:LYS:CD	1:A:67:ASP:H	2.25	0.46
1:A:244:ILE:HD13	1:A:267:ALA:HB2	1.98	0.46
1:A:228:LEU:CB	1:A:242:GLN:NE2	2.77	0.46
1:A:518:VAL:O	1:A:522:ILE:HG13	2.15	0.46
1:A:396:GLU:HA	1:A:399:GLU:HG2	1.98	0.46
1:B:88:TRP:CZ2	1:B:154:LYS:HD3	2.51	0.45
1:B:373:GLN:HE21	1:B:408:ALA:H	1.64	0.45
1:A:552:VAL:O	1:A:553:SER:HB3	2.16	0.45
1:A:511:ASP:HB3	4:A:616:HOH:O	2.16	0.45
1:B:120:LEU:HD23	1:B:125:ARG:HG2	1.98	0.45
1:B:261:VAL:HG13	1:B:276:VAL:HG11	1.99	0.45
1:A:131:THR:HG22	1:A:143:ARG:HH11	1.76	0.45
1:A:131:THR:CG2	1:A:143:ARG:HG2	2.18	0.45
1:B:125:ARG:HD3	1:B:147:ASN:HA	1.99	0.44
1:B:249:LYS:CG	1:B:250:ASP:H	2.29	0.44
1:A:21:VAL:HG12	1:A:22:LYS:N	2.31	0.44
1:A:135:ILE:C	1:A:137:ASN:H	2.25	0.44
1:B:353:LYS:HD3	1:B:429:LEU:HD22	1.99	0.44
1:A:54:ASN:O	1:A:143:ARG:NH2	2.51	0.44
1:A:105:SER:HB2	1:A:198:HIS:ND1	2.33	0.44
1:A:129:ALA:HA	1:A:144:TYR:O	2.18	0.44
1:B:175:ASN:ND2	1:B:201:LYS:NZ	2.63	0.44
1:B:65:LYS:CG	1:B:68:SER:HB2	2.42	0.44
1:A:48:SER:O	1:A:144:TYR:HA	2.18	0.44
1:A:139:THR:HB	1:A:140:PRO:HD2	2.00	0.44
1:A:235:HIS:HB3	1:A:236:PRO:HD2	2.00	0.44
1:B:330:GLN:HE21	1:B:338:THR:HG1	1.62	0.43
1:B:407:GLN:HE21	1:B:407:GLN:HB2	1.50	0.43
1:A:131:THR:HG23	1:A:143:ARG:NH1	2.32	0.43
1:B:328:GLU:O	1:B:339:TYR:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286:THR:HG23	1:A:286:THR:O	2.17	0.43
1:B:257:ILE:O	1:B:261:VAL:HG23	2.18	0.43
1:A:199:ARG:HD3	1:A:219:LYS:HE2	1.99	0.43
1:A:457:TYR:CD1	1:A:457:TYR:C	2.97	0.43
1:B:86:ASP:OD1	1:B:90:VAL:HG21	2.18	0.43
1:A:179:VAL:O	1:A:189:VAL:HA	2.19	0.43
1:B:21:VAL:HB	1:B:59:PRO:HD3	2.01	0.43
1:B:47:ILE:HD12	1:B:144:TYR:CG	2.54	0.43
1:B:354:TYR:HD2	1:B:374:LYS:HD2	1.83	0.43
1:A:246:LEU:HD22	1:A:260:LEU:CD1	2.49	0.43
1:A:406:TRP:CE3	1:A:407:GLN:NE2	2.86	0.43
1:B:24:TRP:CZ3	1:B:59:PRO:CG	2.97	0.43
1:A:486:LEU:CD1	1:A:521:ILE:HG23	2.49	0.43
1:B:269:GLN:H	1:B:269:GLN:HG2	1.65	0.43
1:A:52:PRO:HA	1:A:53:GLU:C	2.44	0.43
1:A:195:ILE:HD12	1:A:199:ARG:NH2	2.34	0.43
1:B:82:LYS:HE3	1:B:82:LYS:HB2	1.72	0.43
1:B:388:LYS:HA	1:B:413:GLU:O	2.19	0.43
1:B:275:LYS:H	1:B:306:ASN:HD21	1.67	0.43
1:A:308:GLU:O	1:A:311:LYS:HB2	2.19	0.42
1:B:296:THR:HG22	1:B:297:GLU:N	2.33	0.42
1:A:228:LEU:CB	1:A:242:GLN:HE21	2.32	0.42
1:A:534:ALA:HB1	4:A:594:HOH:O	2.18	0.42
1:B:175:ASN:ND2	1:B:201:LYS:HD2	2.34	0.42
1:A:180:ILE:CG2	1:A:187:LEU:HD22	2.49	0.42
1:A:494:ASN:HB3	1:B:289:LEU:HD12	2.01	0.42
1:B:325:LEU:HD12	1:B:385:LYS:HB3	2.02	0.42
1:A:63:ILE:HG22	1:A:64:LYS:N	2.34	0.42
1:A:271:TYR:CE1	1:A:314:VAL:HG23	2.55	0.42
1:A:429:LEU:HD11	1:A:506:ILE:HG22	2.02	0.42
1:B:12:LEU:HD23	1:B:84:THR:HG22	2.01	0.42
1:B:344:GLU:HA	1:B:345:PRO:HD3	1.95	0.42
1:B:236:PRO:HA	1:B:239:TRP:CD2	2.54	0.42
1:A:312:GLU:OE2	1:A:312:GLU:N	2.52	0.41
1:A:209:LEU:HD23	1:A:216:THR:HG21	2.01	0.41
1:B:249:LYS:HG2	1:B:250:ASP:N	2.32	0.41
1:A:458:VAL:HG11	1:B:286:THR:HG21	2.01	0.41
1:A:460:ASN:ND2	1:B:288:ALA:HB2	2.34	0.41
1:B:80:LEU:HD13	1:B:153:TRP:CD1	2.55	0.41
1:A:431:LYS:HB2	1:A:431:LYS:NZ	2.35	0.41
1:B:296:THR:O	1:B:300:GLU:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:LYS:HZ1	1:A:69:THR:HA	1.84	0.41
1:A:108:VAL:HG13	1:A:108:VAL:O	2.20	0.41
1:A:253:THR:HA	1:A:291:GLU:O	2.21	0.41
1:B:388:LYS:HE2	1:B:415:GLU:OE1	2.21	0.41
1:A:424:LYS:HE3	1:A:424:LYS:HB3	1.77	0.41
1:A:103:LYS:HD3	1:A:103:LYS:HA	1.85	0.41
1:A:218:ASP:OD2	1:A:221:HIS:ND1	2.53	0.41
1:A:484:LEU:HD12	1:A:484:LEU:HA	1.83	0.41
1:B:260:LEU:HD22	1:B:264:LEU:HD11	2.02	0.41
1:A:339:TYR:CZ	1:A:352:GLY:HA3	2.56	0.41
1:A:473:THR:C	1:A:475:GLN:H	2.29	0.41
1:B:172:ARG:HD2	4:B:564:HOH:O	2.20	0.41
1:A:8:VAL:O	1:A:10:VAL:HG23	2.21	0.40
1:A:126:LYS:HA	1:A:145:GLN:NE2	2.37	0.40
1:A:320:ASP:HA	1:A:321:PRO:HD3	1.84	0.40
1:B:245:VAL:O	1:B:245:VAL:HG23	2.21	0.40
1:B:270:ILE:HG12	1:B:346:PHE:O	2.21	0.40
1:A:3:SER:OG	1:A:5:ILE:HG22	2.21	0.40
1:B:338:THR:HG21	1:B:427:TYR:O	2.21	0.40
1:A:195:ILE:HG23	1:A:199:ARG:NE	2.37	0.40
1:A:209:LEU:HD12	1:A:209:LEU:HA	1.92	0.40
1:A:210:LEU:C	1:A:212:TRP:H	2.29	0.40
1:A:225:PRO:HA	1:A:226:PRO:C	2.47	0.40
1:B:368:LEU:HD13	1:B:398:TRP:CZ3	2.56	0.40
1:A:24:TRP:CD2	1:A:25:PRO:HD2	2.57	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%)	522 (95%)	22 (4%)	7 (1%)	9 18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	395/560 (70%)	384 (97%)	11 (3%)	0	100	100
All	All	946/1120 (84%)	906 (96%)	33 (4%)	7 (1%)	18	34

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	195	ILE
1	A	543	GLY
1	A	334	GLN
1	A	410	TRP
1	A	136	ASN
1	A	356	ARG
1	A	412	PRO

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%)	453 (92%)	42 (8%)	10	22
1	B	367/500 (73%)	338 (92%)	29 (8%)	11	24
All	All	862/1000 (86%)	791 (92%)	71 (8%)	10	23

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	66	LYS
1	A	73	LYS
1	A	74	LEU
1	A	78	ARG
1	A	89	GLU
1	A	120	LEU
1	A	137	ASN
1	A	194	GLU

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Mol	Chain	Res	Type
1	A	205	LEU
1	A	209	LEU
1	A	219	LYS
1	A	242	GLN
1	A	264	LEU
1	A	282	LEU
1	A	283	LEU
1	A	289	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	347	LYS
1	A	349	LEU
1	A	361	HIS
1	A	379	SER
1	A	402	TRP
1	A	406	TRP
1	A	423	VAL
1	A	428	GLN
1	A	431	LYS
1	A	448	ARG
1	A	479	LEU
1	A	500	GLN
1	A	507	GLN
1	A	517	LEU
1	A	529	GLU
1	A	533	LEU
1	A	551	LEU
1	B	11	LYS
1	B	22	LYS
1	B	48	SER
1	B	80	LEU
1	B	83	ARG
1	B	91	GLN
1	B	184	MET
1	B	203	GLU
1	B	205	LEU

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Mol	Chain	Res	Type
1	B	209	LEU
1	B	234	LEU
1	B	238	LYS
1	B	248	GLU
1	B	249	LYS
1	B	250	ASP
1	B	260	LEU
1	B	269	GLN
1	B	287	LYS
1	B	293	ILE
1	B	308	GLU
1	B	310	LEU
1	B	323	LYS
1	B	336	GLN
1	B	368	LEU
1	B	385	LYS
1	B	403	THR
1	B	407	GLN
1	B	422	LEU
1	B	429	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	GLN
1	A	242	GLN
1	A	258	GLN
1	A	332	GLN
1	A	336	GLN
1	A	348	ASN
1	A	373	GLN
1	A	407	GLN
1	A	428	GLN
1	A	487	GLN
1	A	500	GLN
1	A	507	GLN
1	A	509	GLN
1	A	512	GLN
1	A	519	ASN
1	A	520	GLN
1	A	524	GLN
1	B	85	GLN

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Mol	Chain	Res	Type
1	B	147	ASN
1	B	161	GLN
1	B	175	ASN
1	B	182	GLN
1	B	197	GLN
1	B	235	HIS
1	B	255	ASN
1	B	306	ASN
1	B	330	GLN
1	B	336	GLN
1	B	340	GLN
1	B	348	ASN
1	B	373	GLN
1	B	407	GLN
1	B	418	ASN
1	B	428	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	SO4	B	561	-	4,4,4	0.29	0	6,6,6	0.07	0
3	SO4	A	563	-	4,4,4	0.37	0	6,6,6	0.17	0
3	SO4	A	562	-	4,4,4	0.35	0	6,6,6	0.15	0
2	KRV	A	561	-	25,25,25	2.22	8 (32%)	32,36,36	1.78	8 (25%)
3	SO4	A	564	-	4,4,4	0.41	0	6,6,6	0.07	0
3	SO4	A	565	-	4,4,4	0.37	0	6,6,6	0.13	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
2	KRV	A	561	-	-	2/16/16/16	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	561	KRV	C11-C12	4.39	1.53	1.43
2	A	561	KRV	C21-C22	4.32	1.45	1.39
2	A	561	KRV	C13-C12	4.04	1.42	1.35
2	A	561	KRV	C21-C18	3.74	1.46	1.39
2	A	561	KRV	C23-C22	3.31	1.44	1.39
2	A	561	KRV	C23-C16	3.12	1.44	1.39
2	A	561	KRV	C17-C16	2.67	1.43	1.39
2	A	561	KRV	C9-N8	-2.39	1.35	1.39

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	561	KRV	C16-C14-C13	4.43	127.39	119.34
2	A	561	KRV	O15-C14-C13	-3.61	113.93	120.35
2	A	561	KRV	C25-C12-C13	3.60	125.25	119.06
2	A	561	KRV	C1-N8-C13	3.03	122.85	116.64
2	A	561	KRV	N10-C9-N8	2.56	120.04	115.45
2	A	561	KRV	C21-C18-C19	2.38	122.76	119.55
2	A	561	KRV	C11-N10-C9	-2.31	124.31	127.34
2	A	561	KRV	O29-C9-N10	-2.03	117.74	121.49

There are no chirality outliers.

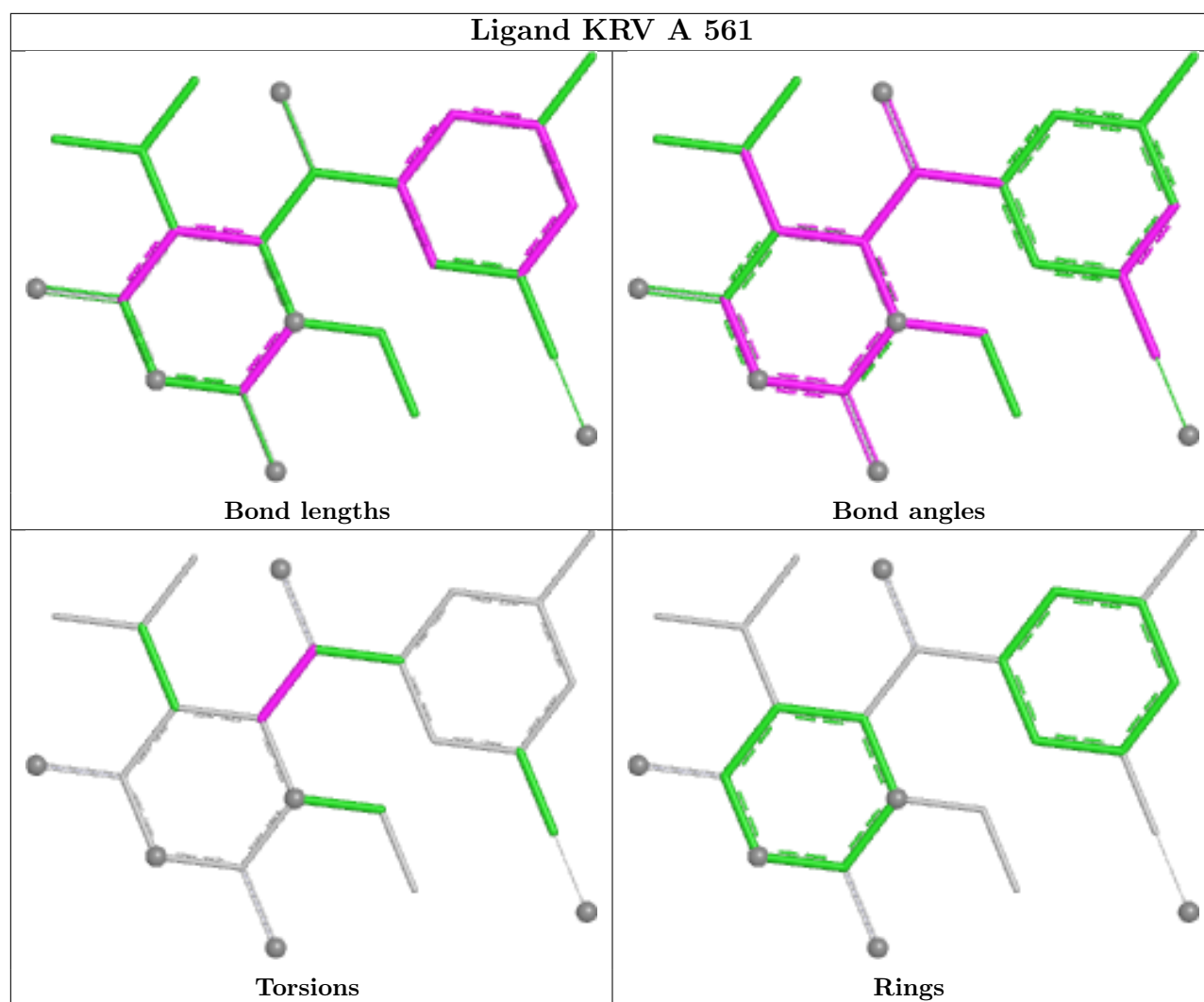
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	561	KRV	C12-C13-C14-C16
2	A	561	KRV	N8-C13-C14-C16

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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%)	0.03	10 (1%) 67 64	40, 62, 93, 117	0
1	B	403/560 (71%)	-0.08	6 (1%) 72 68	36, 57, 93, 108	0
All	All	956/1120 (85%)	-0.02	16 (1%) 69 65	36, 60, 93, 117	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	553	SER	3.9
1	A	551	LEU	3.5
1	B	232	TYR	3.0
1	B	361	HIS	2.9
1	B	5	ILE	2.9
1	B	346	PHE	2.9
1	A	52	PRO	2.7
1	A	24	TRP	2.7
1	A	181	TYR	2.7
1	B	24	TRP	2.7
1	A	291	GLU	2.5
1	B	429	LEU	2.5
1	A	446	ALA	2.4
1	A	402	TRP	2.2
1	A	445	ALA	2.2
1	A	452	LEU	2.1

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

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

6.3 Carbohydrates [i](#)

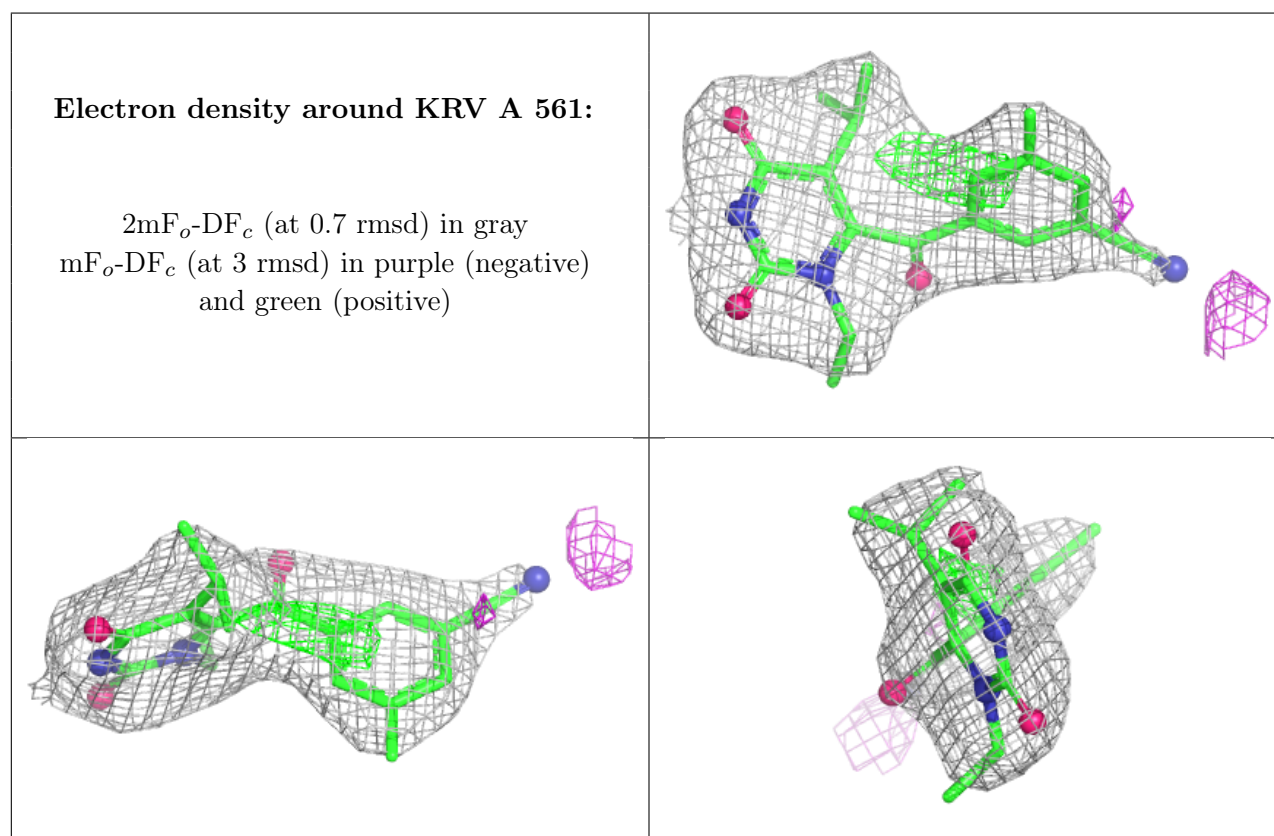
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	A	565	5/5	0.65	0.11	120,120,121,121	0
3	SO4	A	564	5/5	0.72	0.15	121,121,121,121	0
3	SO4	A	562	5/5	0.90	0.09	77,78,80,82	0
3	SO4	B	561	5/5	0.90	0.14	82,83,84,85	0
3	SO4	A	563	5/5	0.92	0.07	76,77,78,79	0
2	KRV	A	561	24/24	0.93	0.14	55,59,66,67	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.