



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:57 PM UTC

PDB ID : 4WE1 / pdb_00004we1
Title : Crystal Structure of HIV-1 Reverse Transcriptase in Complex with 5-(2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy)phenoxy)-2-naphthonitrile (JLJ600)
Authors : Frey, K.M.; Anderson, K.S.
Deposited on : 2014-09-09
Resolution : 2.49 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

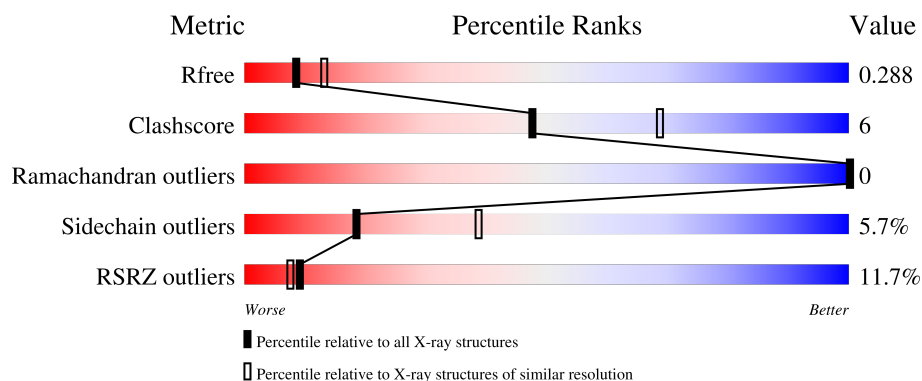
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	557	12% 78% 18% • •
2	B	428	11% 80% 12% • 6%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7807 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 Gag-Pol polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	542	Total	C	N	O	S	0	0	0
			4375	2832	728	808	7			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	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

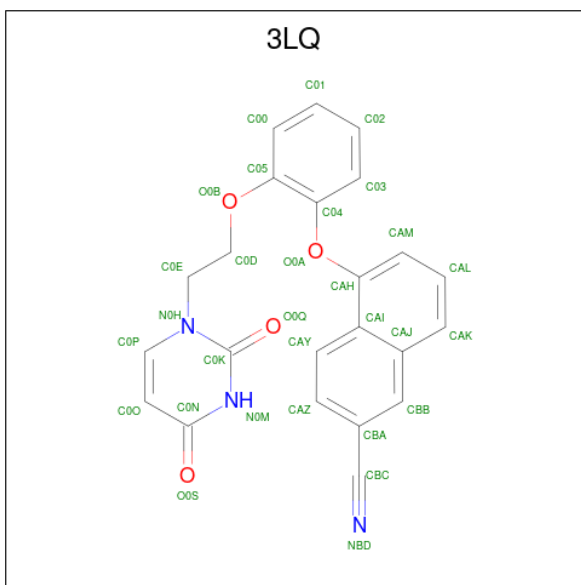
- Molecule 2 is a protein called Gag-Pol polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	402	Total	C	N	O	S	0	1	0
			3296	2145	545	600	6			

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 5-{2-[2-(2,4-dioxo-3,4-dihydropyrimidin-1(2H)-yl)ethoxy]phenoxy}naphthalen e-2-carbonitrile (CCD ID: 3LQ) (formula: C₂₃H₁₇N₃O₄).

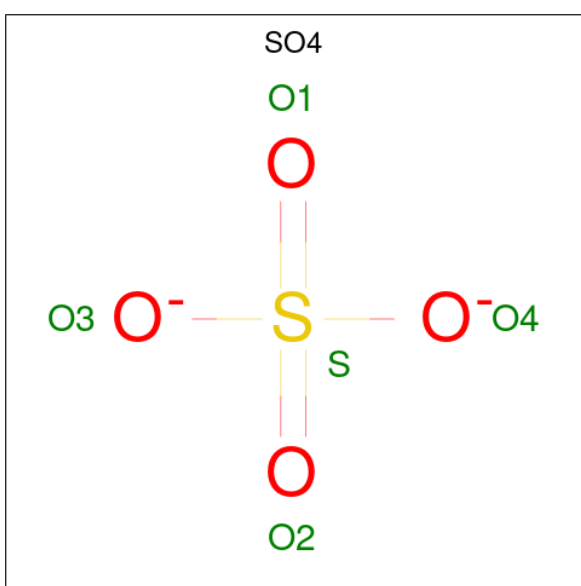


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			30	23	3	4		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

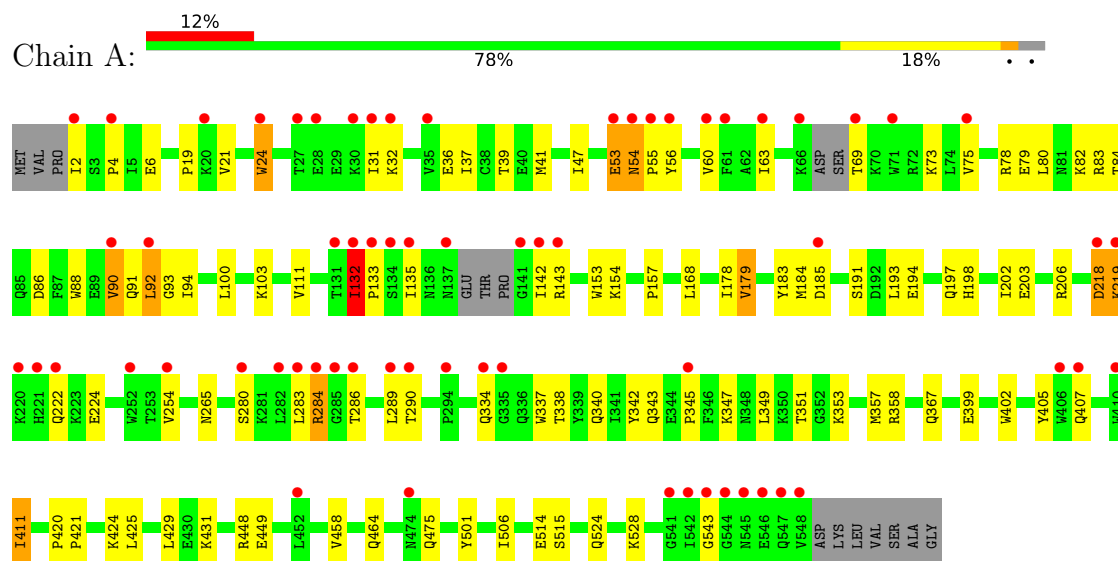
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	44	Total O 44 44	0	0
6	B	40	Total O 40 40	0	0

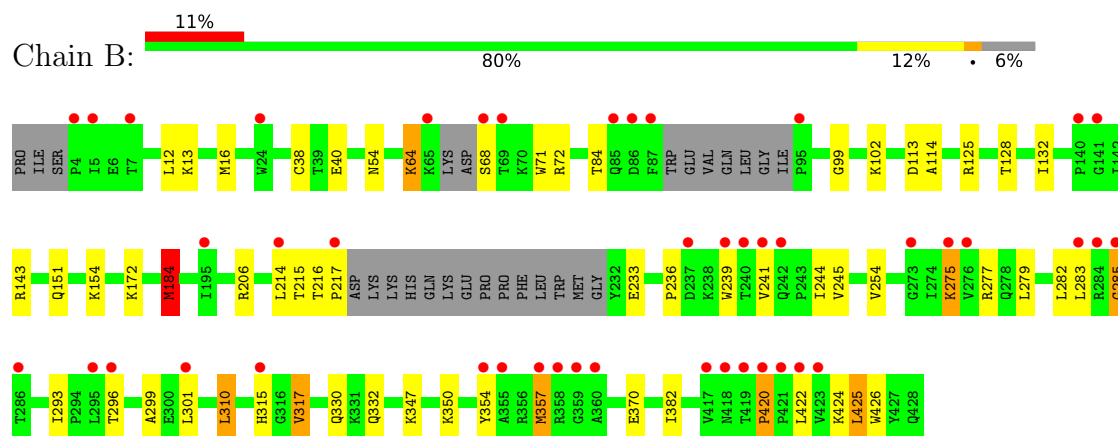
3 Residue-property plots [i](#)

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: Gag-Pol polyprotein



• Molecule 2: Gag-Pol polyprotein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	220.27Å 67.22Å 102.48Å 90.00° 108.08° 90.00°	Depositor
Resolution (Å)	38.77 – 2.49 38.77 – 2.49	Depositor EDS
% Data completeness (in resolution range)	98.5 (38.77-2.49) 98.6 (38.77-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.62 (at 2.48Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.238 , 0.287 0.245 , 0.288	Depositor DCC
R_{free} test set	2000 reflections (4.04%)	wwPDB-VP
Wilson B-factor (Å ²)	50.9	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 41.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7807	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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, MG, 3LQ

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	0.32	0/4489	0.77	2/6106 (0.0%)
2	B	0.30	0/3391	0.81	8/4609 (0.2%)
All	All	0.31	0/7880	0.78	10/10715 (0.1%)

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
2	B	0	1
All	All	0	2

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	285	GLY	N-CA-C	8.14	123.41	110.87
2	B	420	PRO	N-CA-C	6.61	118.76	110.70
2	B	420	PRO	CA-C-N	-6.50	112.21	118.97
2	B	420	PRO	C-N-CA	-6.50	112.21	118.97
1	A	132	ILE	CA-C-N	6.01	126.32	119.83
1	A	132	ILE	C-N-CA	6.01	126.32	119.83
2	B	54	ASN	CA-C-N	5.62	125.24	119.56
2	B	54	ASN	C-N-CA	5.62	125.24	119.56
2	B	382	ILE	N-CA-C	5.56	115.64	110.42
2	B	184	MET	CB-CA-C	-5.35	109.95	117.23

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	54	ASN	Peptide
2	B	420	PRO	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	4375	0	4369	59	0
2	B	3296	0	3309	30	0
3	A	30	0	17	3	0
4	A	2	0	0	0	0
5	A	20	0	0	1	0
6	A	44	0	0	0	0
6	B	40	0	0	2	0
All	All	7807	0	7695	87	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 6.

All (87) 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:543:GLY:HA2	2:B:285:GLY:HA2	1.71	0.73
2:B:12:LEU:HG	2:B:84:THR:HG22	1.71	0.71
1:A:265:ASN:OD1	1:A:353:LYS:NZ	2.24	0.69
1:A:191:SER:OG	1:A:198:HIS:ND1	2.26	0.67
2:B:114:ALA:HB2	2:B:214:LEU:HG	1.80	0.64
1:A:90:VAL:HG21	1:A:157:PRO:HB2	1.80	0.63
1:A:194:GLU:H	1:A:197:GLN:HE21	1.47	0.62
1:A:54:ASN:HD22	1:A:56:TYR:HB2	1.65	0.62
2:B:357:MET:N	2:B:357:MET:SD	2.73	0.61
2:B:279:LEU:HD23	2:B:299:ALA:HB1	1.86	0.58
2:B:254:VAL:HG13	2:B:283:LEU:HD22	1.86	0.57
2:B:241:VAL:HG22	2:B:350:LYS:HA	1.86	0.57
1:A:219:LYS:HD2	1:A:222:GLN:HG2	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:244:ILE:HB	2:B:310:LEU:HG	1.88	0.56
2:B:72:ARG:NH2	2:B:151:GLN:OE1	2.38	0.56
2:B:296:THR:HG23	2:B:299:ALA:H	1.70	0.56
1:A:193:LEU:HB3	1:A:197:GLN:HG3	1.88	0.55
1:A:448:ARG:NH1	5:A:604:SO4:O2	2.39	0.55
1:A:31:ILE:HG23	1:A:132:ILE:HD11	1.89	0.55
2:B:99:GLY:HA2	2:B:102:LYS:HD3	1.89	0.54
1:A:19:PRO:HG3	1:A:80:LEU:HB2	1.90	0.54
1:A:84:THR:HG21	1:A:153:TRP:HE1	1.73	0.53
1:A:86:ASP:OD1	1:A:154:LYS:NZ	2.33	0.52
2:B:172:LYS:NZ	6:B:535:HOH:O	2.41	0.52
1:A:24:TRP:H	1:A:24:TRP:CD1	2.27	0.52
1:A:343:GLN:HG3	1:A:349:LEU:HD21	1.91	0.52
1:A:132:ILE:HD12	1:A:133:PRO:HD2	1.92	0.52
2:B:13:LYS:HB2	2:B:16:MET:HE2	1.92	0.51
1:A:337:TRP:HE1	1:A:367:GLN:NE2	2.08	0.51
2:B:206:ARG:NH1	2:B:216:THR:O	2.43	0.51
1:A:37:ILE:HG22	1:A:41:MET:HE2	1.93	0.51
2:B:275:LYS:HD3	2:B:277:ARG:H	1.74	0.51
1:A:218:ASP:OD1	1:A:218:ASP:N	2.33	0.50
2:B:236:PRO:HA	2:B:239:TRP:CD2	2.47	0.50
1:A:203:GLU:OE2	1:A:206:ARG:NH1	2.45	0.49
1:A:60:VAL:HG23	1:A:75:VAL:HG22	1.95	0.49
1:A:405:TYR:CE2	1:A:407:GLN:HB3	2.47	0.49
2:B:354:TYR:OH	2:B:370:GLU:OE1	2.30	0.48
1:A:340:GLN:HG3	1:A:351:THR:HG22	1.94	0.48
1:A:78:ARG:O	1:A:82:LYS:HG3	2.13	0.48
1:A:2:ILE:C	1:A:4:PRO:HD2	2.39	0.48
1:A:36:GLU:O	1:A:39:THR:OG1	2.22	0.48
1:A:342:TYR:HA	1:A:349:LEU:HD23	1.95	0.48
1:A:132:ILE:HG23	1:A:142:ILE:HB	1.97	0.47
1:A:179:VAL:HG13	3:A:601:3LQ:H2	1.97	0.47
2:B:424:LYS:NZ	6:B:518:HOH:O	2.46	0.47
1:A:357:MET:O	1:A:358:ARG:HG2	2.16	0.46
1:A:284:ARG:NH2	1:A:514:GLU:OE1	2.48	0.46
1:A:280:SER:HA	1:A:283:LEU:HB2	1.97	0.46
2:B:425:LEU:HD12	2:B:426:TRP:CD1	2.51	0.46
1:A:54:ASN:ND2	1:A:56:TYR:HB2	2.29	0.46
1:A:342:TYR:CE1	1:A:345:PRO:HA	2.51	0.46
1:A:458:VAL:HG13	1:A:464:GLN:HG2	1.98	0.45
1:A:55:PRO:O	1:A:56:TYR:CG	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:GLU:OE2	1:A:83:ARG:NH1	2.44	0.45
1:A:32:LYS:HA	1:A:32:LYS:HD3	1.75	0.45
2:B:215:THR:HG22	2:B:217:PRO:HD3	1.98	0.44
1:A:100:LEU:HB3	3:A:601:3LQ:H6	1.98	0.44
1:A:103:LYS:HE3	1:A:179:VAL:HG11	2.00	0.44
2:B:282:LEU:HB3	2:B:293:ILE:HG21	2.00	0.44
1:A:41:MET:HE3	1:A:47:ILE:HD13	2.00	0.43
1:A:429:LEU:HD11	1:A:506:ILE:HG22	1.99	0.43
1:A:88:TRP:CE2	2:B:143:ARG:HD2	2.53	0.43
2:B:317:VAL:HG12	2:B:347:LYS:HB3	2.00	0.43
2:B:154:LYS:HE3	2:B:184:MET:HE1	2.01	0.43
1:A:424:LYS:NZ	1:A:425:LEU:H	2.16	0.43
2:B:125:ARG:O	2:B:128:THR:OG1	2.36	0.43
2:B:332:GLN:HA	2:B:424:LYS:HE3	2.01	0.43
1:A:79:GLU:O	1:A:83:ARG:HG3	2.19	0.42
1:A:399:GLU:HA	1:A:402:TRP:CE2	2.54	0.42
2:B:38:CYS:SG	2:B:132:ILE:HD11	2.59	0.42
1:A:402:TRP:CE3	1:A:411:ILE:HG21	2.55	0.41
1:A:56:TYR:O	1:A:143:ARG:NH2	2.52	0.41
2:B:422:LEU:HD23	2:B:422:LEU:HA	1.84	0.41
1:A:41:MET:HE1	1:A:73:LYS:NZ	2.35	0.41
1:A:92:LEU:HD12	1:A:93:GLY:H	1.86	0.41
1:A:94:ILE:HG22	1:A:183:TYR:CE1	2.56	0.41
1:A:420:PRO:HA	1:A:421:PRO:C	2.45	0.41
1:A:53:GLU:H	1:A:53:GLU:CD	2.29	0.41
2:B:64:LYS:HE2	2:B:71:TRP:CZ2	2.56	0.41
1:A:475:GLN:HB3	1:A:501:TYR:CE2	2.56	0.40
2:B:310:LEU:HD12	2:B:310:LEU:HA	1.84	0.40
1:A:202:ILE:O	1:A:206:ARG:HG3	2.21	0.40
1:A:524:GLN:O	1:A:528:LYS:HG2	2.21	0.40
1:A:132:ILE:HA	1:A:133:PRO:HD3	1.77	0.40
1:A:179:VAL:CG1	3:A:601:3LQ:H2	2.51	0.40
1:A:80:LEU:O	1:A:84:THR:HG23	2.22	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	536/557 (96%)	525 (98%)	11 (2%)	0	100	100
2	B	395/428 (92%)	389 (98%)	6 (2%)	0	100	100
All	All	931/985 (94%)	914 (98%)	17 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	471/495 (95%)	439 (93%)	32 (7%)	14	31
2	B	360/390 (92%)	344 (96%)	16 (4%)	25	50
All	All	831/885 (94%)	783 (94%)	48 (6%)	18	38

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLU
1	A	21	VAL
1	A	24	TRP
1	A	53	GLU
1	A	63	ILE
1	A	69	THR
1	A	90	VAL
1	A	91	GLN

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Mol	Chain	Res	Type
1	A	92	LEU
1	A	111	VAL
1	A	132	ILE
1	A	135	ILE
1	A	168	LEU
1	A	178	ILE
1	A	179	VAL
1	A	184	MET
1	A	185	ASP
1	A	218	ASP
1	A	219	LYS
1	A	224	GLU
1	A	254	VAL
1	A	284	ARG
1	A	286	THR
1	A	289	LEU
1	A	290	THR
1	A	334	GLN
1	A	338	THR
1	A	347	LYS
1	A	411	ILE
1	A	431	LYS
1	A	449	GLU
1	A	515	SER
2	B	40	GLU
2	B	64	LYS
2	B	68	SER
2	B	113	ASP
2	B	184	MET
2	B	233[A]	GLU
2	B	233[B]	GLU
2	B	245	VAL
2	B	275	LYS
2	B	301	LEU
2	B	310	LEU
2	B	315	HIS
2	B	317	VAL
2	B	330	GLN
2	B	357	MET
2	B	425	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	161	GLN
1	A	182	GLN
1	A	197	GLN
1	A	208	HIS
1	A	242	GLN
1	A	340	GLN
1	A	361	HIS
1	A	367	GLN
1	A	407	GLN
1	A	428	GLN
1	A	500	GLN
2	B	147	ASN
2	B	161	GLN
2	B	174	GLN
2	B	235	HIS
2	B	255	ASN
2	B	278	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](#)

Of 7 ligands modelled in this entry, 2 are 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
5	SO4	A	606	-	4,4,4	0.23	0	6,6,6	0.08	0
3	3LQ	A	601	-	33,33,33	1.57	5 (15%)	45,45,45	1.51	4 (8%)
5	SO4	A	607	-	4,4,4	0.25	0	6,6,6	0.08	0
5	SO4	A	605	-	4,4,4	0.24	0	6,6,6	0.08	0
5	SO4	A	604	-	4,4,4	0.25	0	6,6,6	0.07	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	3LQ	A	601	-	-	0/12/12/12	0/4/4/4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	3LQ	C0P-N0H	4.06	1.42	1.37
3	A	601	3LQ	CBA-CBC	3.59	1.52	1.44
3	A	601	3LQ	CAZ-CBA	2.71	1.44	1.39
3	A	601	3LQ	O0A-CAH	2.16	1.45	1.39
3	A	601	3LQ	CAI-CAJ	-2.09	1.39	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	3LQ	C0E-N0H-C0K	6.82	125.26	118.48
3	A	601	3LQ	C0E-N0H-C0P	-3.77	114.41	119.75
3	A	601	3LQ	C0D-O0B-C05	2.93	124.77	117.69
3	A	601	3LQ	O0Q-C0K-N0H	2.05	124.61	122.87

There are no chirality outliers.

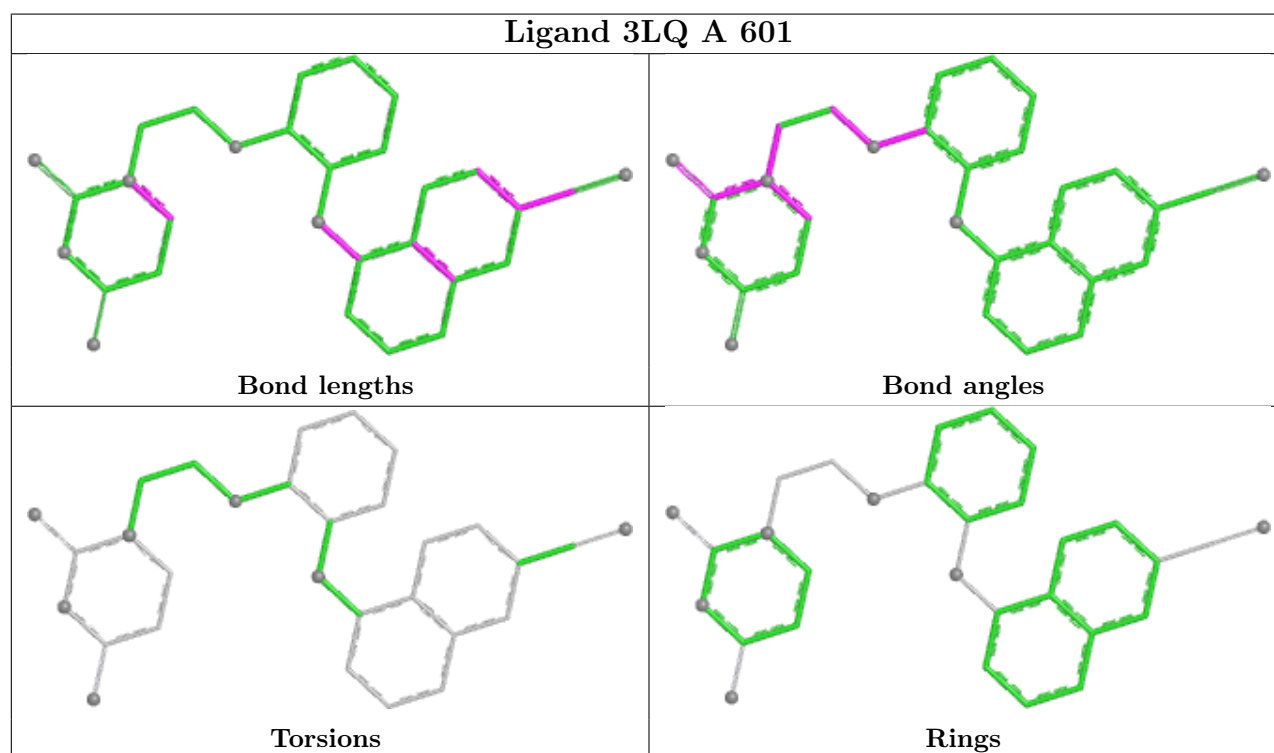
There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	3LQ	3	0
5	A	604	SO4	1	0

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

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	542/557 (97%)	0.84	65 (11%) 9 7	34, 59, 96, 115	0
2	B	402/428 (93%)	0.68	45 (11%) 10 8	36, 56, 85, 130	1 (0%)
All	All	944/985 (95%)	0.77	110 (11%) 9 7	34, 58, 92, 130	1 (0%)

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	542	ILE	8.5
1	A	137	ASN	6.7
1	A	54	ASN	6.6
1	A	545	ASN	5.6
2	B	285	GLY	5.4
2	B	423	VAL	5.4
2	B	5	ILE	5.4
1	A	544	GLY	5.3
1	A	335	GLY	5.2
1	A	32	LYS	5.2
2	B	360	ALA	5.2
1	A	2	ILE	5.0
2	B	241	VAL	4.9
1	A	132	ILE	4.8
1	A	27	THR	4.7
1	A	543	GLY	4.7
2	B	419	THR	4.5
2	B	240	THR	4.5
1	A	294	PRO	4.3
1	A	55	PRO	4.2
2	B	87	PHE	4.2
1	A	69	THR	4.1
1	A	541	GLY	3.9
2	B	286	THR	3.9

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Mol	Chain	Res	Type	RSRZ
2	B	86	ASP	3.9
1	A	547	GLN	3.8
1	A	219	LYS	3.7
2	B	296	THR	3.7
1	A	133	PRO	3.7
2	B	237	ASP	3.7
1	A	24	TRP	3.6
2	B	65	LYS	3.6
1	A	4	PRO	3.6
1	A	406	TRP	3.4
2	B	421	PRO	3.4
1	A	134	SER	3.4
1	A	254	VAL	3.3
1	A	135	ILE	3.3
1	A	221	HIS	3.2
2	B	359	GLY	3.2
2	B	95	PRO	3.2
2	B	85	GLN	3.1
2	B	284	ARG	3.1
1	A	548	VAL	3.1
2	B	420	PRO	3.0
1	A	474	ASN	3.0
1	A	31	ILE	3.0
1	A	28	GLU	3.0
1	A	142	ILE	3.0
2	B	242	GLN	3.0
1	A	546	GLU	2.9
1	A	222	GLN	2.9
2	B	24	TRP	2.9
1	A	220	LYS	2.9
2	B	358	ARG	2.8
2	B	422	LEU	2.8
2	B	417	VAL	2.8
1	A	66	LYS	2.8
1	A	92	LEU	2.7
1	A	280	SER	2.6
1	A	131	THR	2.6
1	A	71	TRP	2.6
2	B	355	ALA	2.6
2	B	217	PRO	2.5
1	A	61	PHE	2.5
2	B	140	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	283	LEU	2.5
1	A	452	LEU	2.5
1	A	286	THR	2.5
2	B	4	PRO	2.4
2	B	283	LEU	2.4
1	A	252	TRP	2.4
2	B	195	ILE	2.4
1	A	407	GLN	2.4
1	A	218	ASP	2.4
1	A	282	LEU	2.4
2	B	295	LEU	2.4
2	B	239	TRP	2.4
2	B	273	GLY	2.4
2	B	418	ASN	2.4
1	A	345	PRO	2.3
2	B	69	THR	2.3
1	A	60	VAL	2.3
1	A	20	LYS	2.3
1	A	75	VAL	2.3
2	B	276	VAL	2.3
1	A	141	GLY	2.3
1	A	290	THR	2.2
2	B	315	HIS	2.2
1	A	289	LEU	2.2
2	B	214	LEU	2.2
1	A	410	TRP	2.2
1	A	53	GLU	2.2
1	A	143	ARG	2.2
1	A	90	VAL	2.2
1	A	63	ILE	2.2
1	A	284	ARG	2.2
2	B	141	GLY	2.2
1	A	285	GLY	2.1
2	B	301	LEU	2.1
2	B	7	THR	2.1
2	B	354	TYR	2.1
1	A	30	LYS	2.1
2	B	357	MET	2.1
1	A	334	GLN	2.1
1	A	35	VAL	2.1
1	A	185	ASP	2.1
2	B	275	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	56	TYR	2.0
2	B	68	SER	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

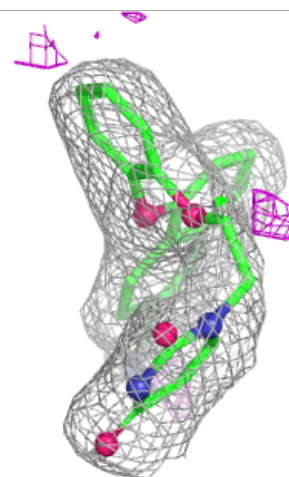
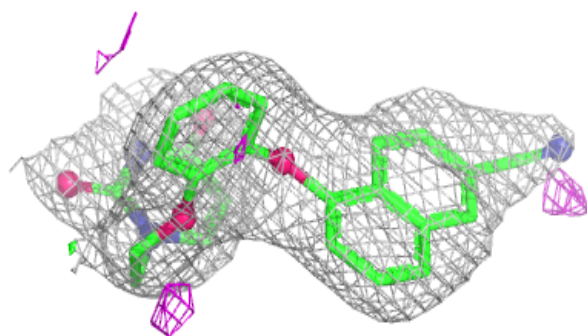
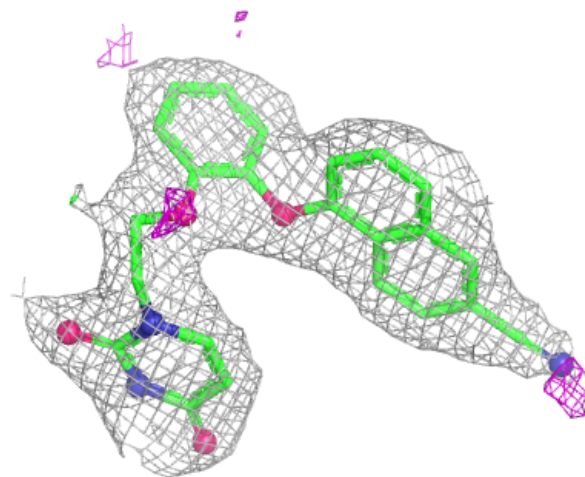
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	A	602	1/1	0.67	0.17	75,75,75,75	0
5	SO4	A	604	5/5	0.77	0.17	75,79,99,99	0
4	MG	A	603	1/1	0.82	0.11	58,58,58,58	0
5	SO4	A	605	5/5	0.89	0.09	69,74,85,88	0
5	SO4	A	606	5/5	0.90	0.09	58,64,72,75	0
3	3LQ	A	601	30/30	0.92	0.10	38,47,52,56	0
5	SO4	A	607	5/5	0.93	0.18	61,64,69,70	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.

Electron density around 3LQ A 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.