



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 02:12 PM UTC

PDB ID : 3S3O / pdb_00003s3o
Title : Crystal structure of the Prototype Foamy Virus (PFV) N224H mutant in-tosome in complex with magnesium and Dolutegravir (S/GSK1349572)
Authors : Hare, S.; Cherepanov, P.
Deposited on : 2011-05-18
Resolution : 2.55 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

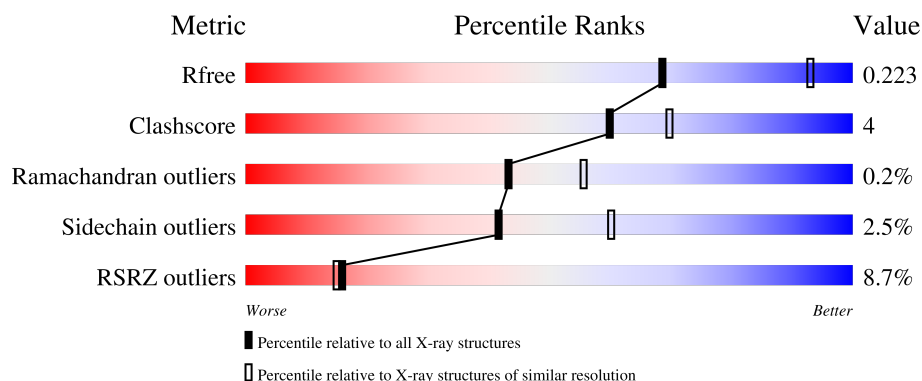
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	395	5% 81% 11% 7%
1	B	395	7% 41% 5% 53%
2	C	19	5% 68% 32%
3	D	17	71% 24% 6%

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 5396 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 PFV integrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	1	0
			2934	1881	517	532	4			
1	B	184	Total	C	N	O	S	0	0	0
			1450	941	236	272	1			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P14350
A	-1	PRO	-	expression tag	UNP P14350
A	0	GLY	-	expression tag	UNP P14350
A	217	SER	GLY	engineered mutation	UNP P14350
A	218	GLY	SER	variant	UNP P14350
A	224	HIS	ASN	engineered mutation	UNP P14350
B	-2	GLY	-	expression tag	UNP P14350
B	-1	PRO	-	expression tag	UNP P14350
B	0	GLY	-	expression tag	UNP P14350
B	217	SER	GLY	variant	UNP P14350
B	218	GLY	SER	variant	UNP P14350
B	224	HIS	ASN	variant	UNP P14350

- Molecule 2 is a DNA chain called 5'-D(*AP*TP*TP*GP*TP*CP*AP*TP*GP*GP*AP*A P*TP*TP*TP*CP*GP*CP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	19	Total	C	N	O	P	0	0	0
			387	187	68	114	18			

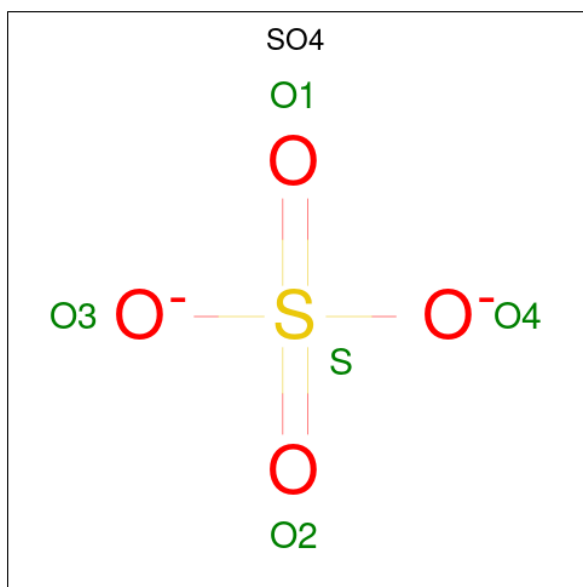
- Molecule 3 is a DNA chain called 5'-D(*TP*GP*CP*GP*AP*AP*AP*TP*TP*CP*CP*A P*TP*GP*AP*CP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	17	Total	C	N	O	P	0	1	0
			345	166	65	98	16			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

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



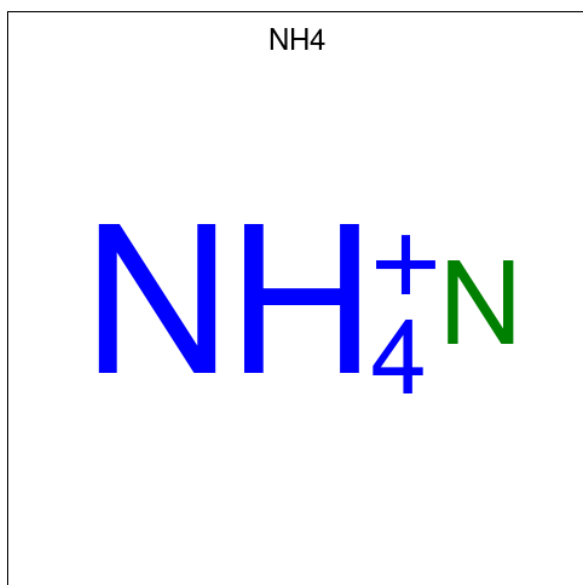
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is AMMONIUM ION (CCD ID: NH4) (formula: H_4N).

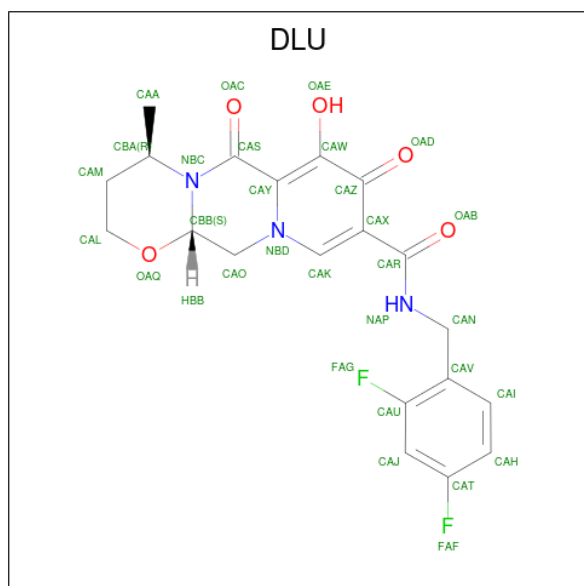


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total N 1 1	0	0

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

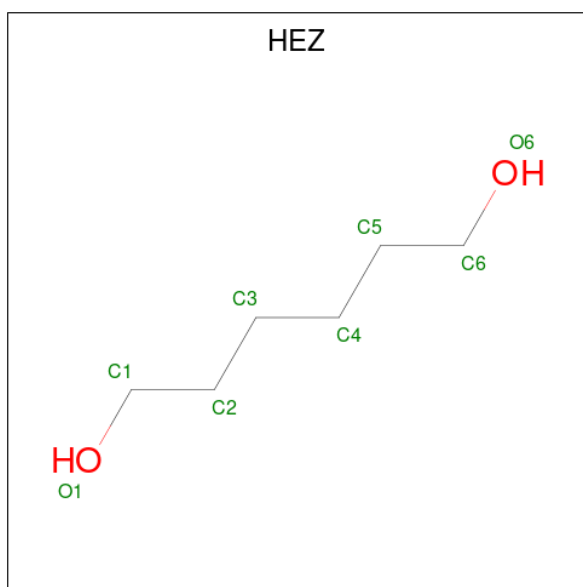
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	2	Total Mg 2 2	0	0

- Molecule 9 is (4R,12aS)-N-(2,4-difluorobenzyl)-7-hydroxy-4-methyl-6,8-dioxo-3,4,6,8,12,12a-hexahydro-2H-pyrido[1',2':4,5]pyrazino[2,1-b][1,3]oxazine-9-carboxamide (CCD ID: DLU) (formula: C₂₀H₁₉F₂N₃O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	F	N	O	
			30	20	2	3	5	

- Molecule 10 is HEXANE-1,6-DIOL (CCD ID: HEZ) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	C	O	0	0
			8	6	2		

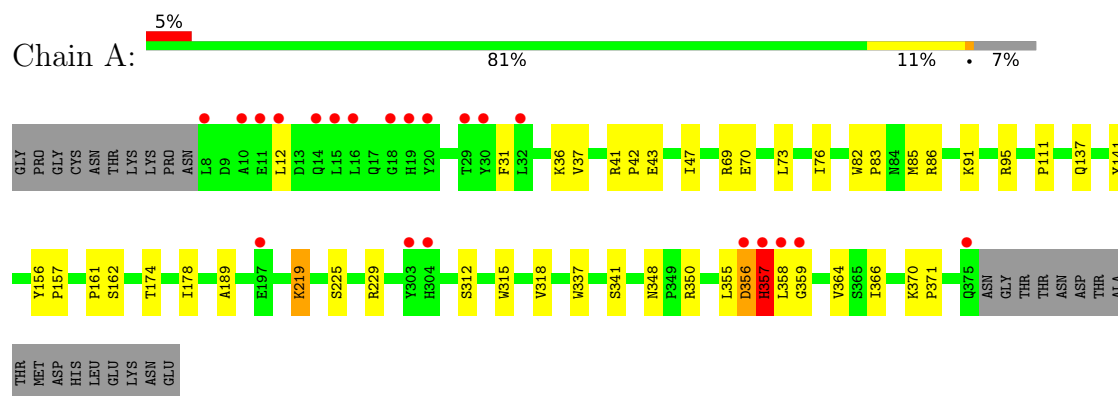
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	117	Total	O	0	0
			117	117		
11	B	36	Total	O	0	0
			36	36		
11	C	28	Total	O	0	0
			28	28		
11	D	18	Total	O	0	0
			18	18		

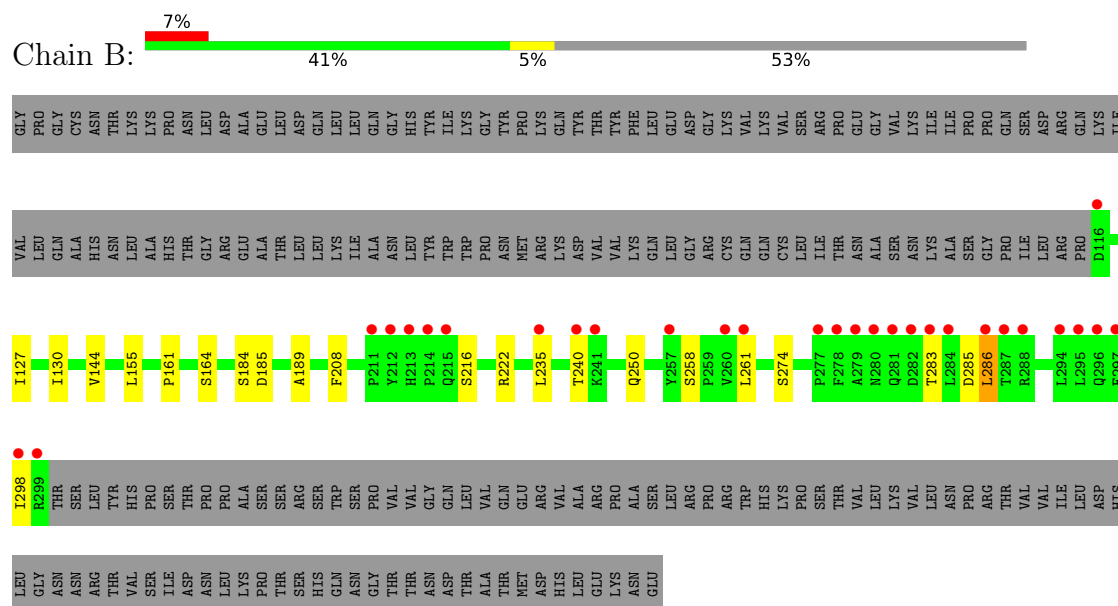
3 Residue-property plots [i](#)

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: PFV integrase



• Molecule 1: PFV integrase



• Molecule 2: 5'-D(*AP*TP*TP*GP*TP*CP*AP*TP*GP*GP*AP*AP*TP*TP*TP*CP*GP*C P*A)-3'





● Molecule 3: 5'-D(*TP*GP*CP*GP*AP*AP*AP*TP*TP*CP*CP*AP*TP*GP*AP*CP*A)-3',

Chain D:
71% 24% 6%

Chain D quality bar chart showing the percentage of residues in different quality categories: 71% (green), 24% (yellow), and 6% (orange).



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	160.44Å 160.44Å 123.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.14 – 2.55 39.14 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.1 (39.14-2.55) 99.0 (39.14-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.92 (at 2.54Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.208 , 0.230 0.201 , 0.223	Depositor DCC
R_{free} test set	2643 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	62.2	Xtriage
Anisotropy	0.271	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.6	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	0.95	EDS
Total number of atoms	5396	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, DLU, GOL, NH4, ZN, HEZ, SO4

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	0.69	0/3017	0.89	4/4117 (0.1%)
1	B	0.69	0/1491	0.88	1/2039 (0.0%)
2	C	0.49	0/433	1.34	3/667 (0.4%)
3	D	0.49	0/387	1.33	5/595 (0.8%)
All	All	0.66	0/5328	0.98	13/7418 (0.2%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	12	DA	C4'-C3'-O3'	-6.61	100.09	110.00
1	A	348	ASN	CA-C-N	-6.30	113.32	119.87
1	A	348	ASN	C-N-CA	-6.30	113.32	119.87
1	A	356	ASP	N-CA-C	6.16	119.99	112.23
3	D	17[A]	DA	C5'-C4'-C3'	-6.01	105.89	114.90
3	D	1	DT	P-O3'-C3'	5.74	128.82	120.20
2	C	14	DT	C4'-C3'-O3'	-5.68	101.48	110.00
2	C	1	DA	P-O3'-C3'	5.66	128.69	120.20
3	D	17[A]	DA	P-O5'-C5'	-5.51	111.73	120.00
2	C	5	DT	O4'-C1'-N1	-5.46	100.22	108.40
3	D	8	DT	N1-C1'-C2'	5.41	121.61	113.50
1	A	31	PHE	N-CA-C	5.13	116.60	108.96
1	B	216	SER	CB-CA-C	-5.11	109.08	116.54

There are no chirality outliers.

There are no planarity outliers.

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	2934	0	2965	28	0
1	B	1450	0	1432	12	0
2	C	387	0	218	2	0
3	D	345	0	193	2	0
4	A	1	0	0	0	0
5	A	10	0	0	0	0
5	B	5	0	0	1	0
6	A	24	0	32	2	0
7	A	1	0	0	0	0
8	A	2	0	0	0	0
9	A	30	0	18	0	0
10	B	8	0	14	3	0
11	A	117	0	0	0	0
11	B	36	0	0	0	0
11	C	28	0	0	0	0
11	D	18	0	0	0	0
All	All	5396	0	4872	43	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 4.

All (43) 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:137:GLN:H	6:A:802:GOL:H2	1.40	0.84
1:A:162:SER:HB2	6:A:801:GOL:H2	1.60	0.82
1:B:258:SER:O	1:B:261:LEU:O	2.08	0.71
1:A:69:ARG:HD2	1:A:70:GLU:OE2	1.93	0.67
1:A:358:LEU:N	1:A:359:GLY:HA2	2.10	0.66
1:A:356:ASP:O	1:A:357:HIS:HB3	1.96	0.66
1:A:73:LEU:HD13	1:A:86:ARG:HG3	1.83	0.60
1:A:91:LYS:HE2	1:A:95:ARG:HH21	1.66	0.60
1:A:219:LYS:HE3	2:C:6:DC:OP1	2.02	0.60
3:D:16:DC:H3'	3:D:17[A]:DA:H5'	1.83	0.60
1:B:250:GLN:HE22	10:B:394:HEZ:H12	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:SER:HB2	1:A:355:LEU:O	2.07	0.55
1:B:250:GLN:HE22	10:B:394:HEZ:C1	2.21	0.54
1:B:283:THR:C	1:B:285:ASP:H	2.18	0.51
1:B:286:LEU:H	1:B:286:LEU:HD12	1.78	0.48
1:A:341:SER:CB	1:A:356:ASP:HA	2.44	0.48
1:A:315:TRP:CE2	1:A:371:PRO:HD3	2.48	0.48
1:A:76:ILE:HG22	1:A:85:MET:CE	2.44	0.47
1:A:174:THR:HB	1:A:178:ILE:HD13	1.95	0.47
1:B:130:ILE:HD12	1:B:144:VAL:HG21	1.96	0.47
1:A:43:GLU:H	1:A:43:GLU:CD	2.23	0.47
1:A:337:TRP:CG	1:A:370:LYS:HE2	2.50	0.47
3:D:16:DC:H3'	3:D:17[A]:DA:C5'	2.47	0.45
1:A:36:LYS:HE2	1:A:47:ILE:HD13	1.99	0.44
1:A:161:PRO:O	1:A:189:ALA:HB2	2.18	0.44
1:A:225:SER:O	1:A:229:ARG:HB2	2.17	0.44
1:A:156:TYR:HA	1:A:157:PRO:HD2	1.82	0.43
2:C:10:DG:H2''	2:C:11:DA:C8	2.54	0.43
1:B:184:SER:O	1:B:208:PHE:HA	2.19	0.43
1:A:91:LYS:HE2	1:A:95:ARG:NH2	2.34	0.42
1:B:222:ARG:NE	5:B:393:SO4:O3	2.42	0.42
1:B:185:ASP:C	1:B:185:ASP:OD2	2.63	0.42
1:A:82:TRP:HB2	1:A:83:PRO:HD2	2.01	0.42
1:B:127:ILE:HA	1:B:144:VAL:O	2.20	0.41
1:A:12:LEU:HD23	1:A:37:VAL:HG21	2.01	0.41
1:A:141:TYR:CE2	1:A:161:PRO:HD3	2.56	0.41
1:A:356:ASP:O	1:A:357:HIS:CB	2.67	0.41
1:A:41:ARG:O	1:A:42:PRO:C	2.63	0.41
1:A:358:LEU:H	1:A:359:GLY:HA2	1.83	0.41
1:A:111:PRO:O	1:A:350:ARG:HD3	2.20	0.41
1:B:155:LEU:O	10:B:394:HEZ:H12	2.20	0.40
1:B:161:PRO:O	1:B:189:ALA:HB2	2.21	0.40
1:A:341:SER:HB3	1:A:356:ASP:HA	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	367/395 (93%)	355 (97%)	11 (3%)	1 (0%)	36	45
1	B	182/395 (46%)	176 (97%)	6 (3%)	0	100	100
All	All	549/790 (70%)	531 (97%)	17 (3%)	1 (0%)	43	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	357	HIS

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	329/354 (93%)	323 (98%)	6 (2%)	51	70
1	B	161/354 (46%)	155 (96%)	6 (4%)	30	45
All	All	490/708 (69%)	478 (98%)	12 (2%)	42	61

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	219	LYS
1	A	312	SER
1	A	318	VAL
1	A	357	HIS
1	A	364	VAL
1	A	366	ILE
1	B	164	SER
1	B	235	LEU
1	B	240	THR
1	B	274	SER
1	B	286	LEU

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Mol	Chain	Res	Type
1	B	298	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	304	HIS
1	A	338	HIS
1	B	205	HIS
1	B	250	GLN
1	B	266	HIS
1	B	296	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 13 ligands modelled in this entry, 3 are monoatomic and 1 is modelled with single atom - leaving 9 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	398	-	4,4,4	0.32	0	6,6,6	0.16	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	DLU	A	399	8	30,33,33	1.78	5 (16%)	34,49,49	1.52	6 (17%)
6	GOL	A	801	-	5,5,5	0.50	0	5,5,5	0.66	0
5	SO4	B	393	-	4,4,4	0.26	0	6,6,6	0.19	0
6	GOL	A	803	-	5,5,5	0.36	0	5,5,5	0.36	0
6	GOL	A	802	-	5,5,5	0.52	0	5,5,5	0.58	0
6	GOL	A	805	-	5,5,5	0.42	0	5,5,5	0.48	0
5	SO4	A	394	-	4,4,4	0.23	0	6,6,6	0.30	0
10	HEZ	B	394	-	7,7,7	0.50	0	6,6,6	0.88	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
9	DLU	A	399	8	-	0/9/35/35	1/4/4/4
6	GOL	A	801	-	-	3/4/4/4	-
6	GOL	A	803	-	-	0/4/4/4	-
6	GOL	A	802	-	-	0/4/4/4	-
6	GOL	A	805	-	-	2/4/4/4	-
10	HEZ	B	394	-	-	2/5/5/5	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	399	DLU	CAS-NBC	6.22	1.42	1.36
9	A	399	DLU	CAY-CAW	3.96	1.41	1.36
9	A	399	DLU	CAW-CAZ	-3.63	1.40	1.50
9	A	399	DLU	OAQ-CAL	2.08	1.48	1.44
9	A	399	DLU	CAK-NBD	2.04	1.41	1.37

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	399	DLU	CAN-NAP-CAR	-3.94	116.43	122.07
9	A	399	DLU	CAA-CBA-CAM	-3.57	106.10	112.74
9	A	399	DLU	OAQ-CAL-CAM	2.73	115.82	110.75
9	A	399	DLU	OAB-CAR-CAX	-2.72	116.31	121.04
9	A	399	DLU	CAN-CAV-CAI	-2.05	116.45	120.98
9	A	399	DLU	OAC-CAS-NBC	-2.04	119.05	121.94

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	801	GOL	C1-C2-C3-O3
6	A	805	GOL	C1-C2-C3-O3
6	A	801	GOL	O2-C2-C3-O3
6	A	805	GOL	O2-C2-C3-O3
6	A	801	GOL	O1-C1-C2-O2
10	B	394	HEZ	C1-C2-C3-C4
10	B	394	HEZ	C4-C5-C6-O6

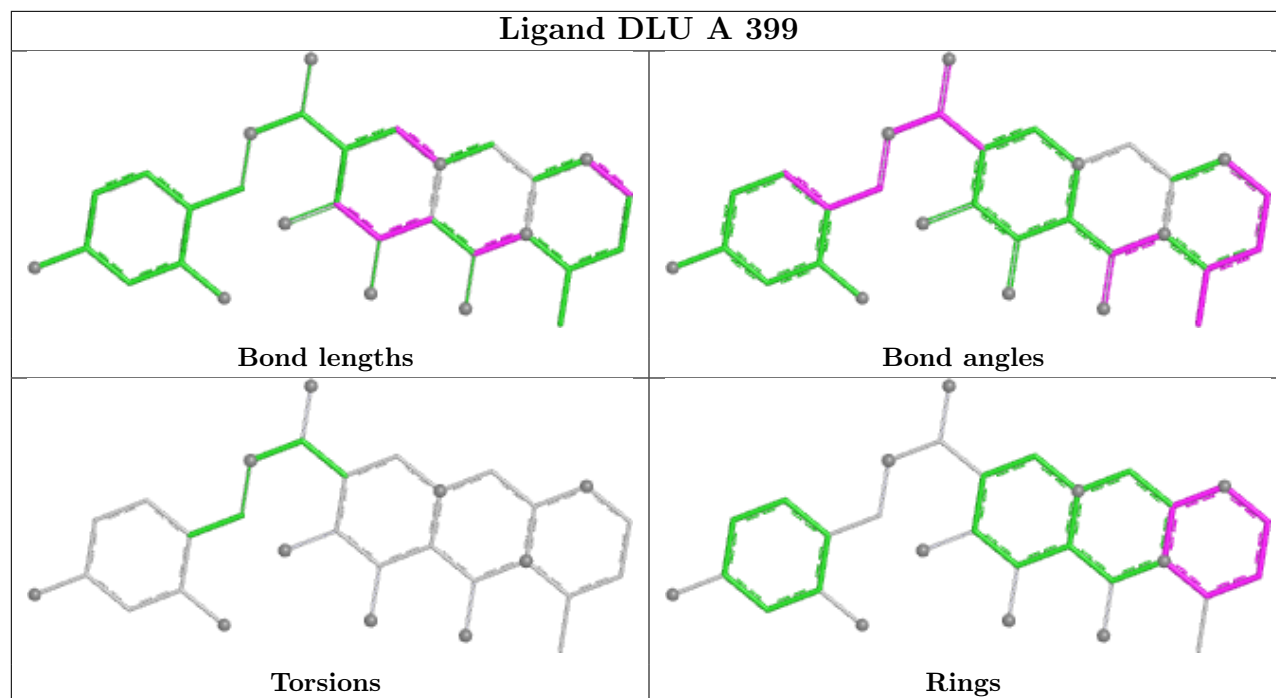
All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	A	399	DLU	CAL-CAM-CBA-CBB-NBC-OAQ

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	801	GOL	1	0
5	B	393	SO4	1	0
6	A	802	GOL	1	0
10	B	394	HEZ	3	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	368/395 (93%)	0.02	21 (5%) 29 29	23, 50, 95, 122	1 (0%)
1	B	184/395 (46%)	0.68	29 (15%) 5 4	44, 61, 137, 146	0
2	C	19/19 (100%)	-0.29	1 (5%) 32 31	38, 55, 81, 104	0
3	D	17/17 (100%)	-0.41	0 100 100	43, 48, 71, 103	0
All	All	588/826 (71%)	0.20	51 (8%) 16 15	23, 54, 115, 146	1 (0%)

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	LEU	7.9
1	B	299	ARG	7.2
1	B	295	LEU	6.2
1	B	298	ILE	6.0
1	B	288	ARG	5.3
1	B	279	ALA	5.2
1	B	283	THR	5.0
1	B	286	LEU	4.8
1	A	358	LEU	4.4
1	B	235	LEU	3.8
1	B	261	LEU	3.8
1	B	213	HIS	3.7
1	A	20	TYR	3.7
1	B	284	LEU	3.5
2	C	1	DA	3.3
1	A	19	HIS	3.3
1	B	212	TYR	3.3
1	B	215	GLN	3.2
1	A	14	GLN	3.2
1	A	359	GLY	3.2
1	B	278	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	281	GLN	3.0
1	B	296	GLN	2.9
1	A	10	ALA	2.8
1	A	32	LEU	2.7
1	A	15	LEU	2.6
1	B	214	PRO	2.5
1	B	282	ASP	2.5
1	A	29	THR	2.5
1	A	16	LEU	2.4
1	B	240	THR	2.4
1	B	287	THR	2.4
1	B	280	ASN	2.4
1	A	357	HIS	2.4
1	A	375	GLN	2.3
1	B	294	LEU	2.3
1	A	12	LEU	2.3
1	A	30	TYR	2.3
1	A	356	ASP	2.2
1	B	257	TYR	2.2
1	A	11	GLU	2.2
1	B	116	ASP	2.2
1	B	297	GLU	2.2
1	B	277	PRO	2.2
1	A	18	GLY	2.1
1	B	260	VAL	2.1
1	A	303	TYR	2.1
1	B	241	LYS	2.0
1	B	211	PRO	2.0
1	A	197	GLU	2.0
1	A	304	HIS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

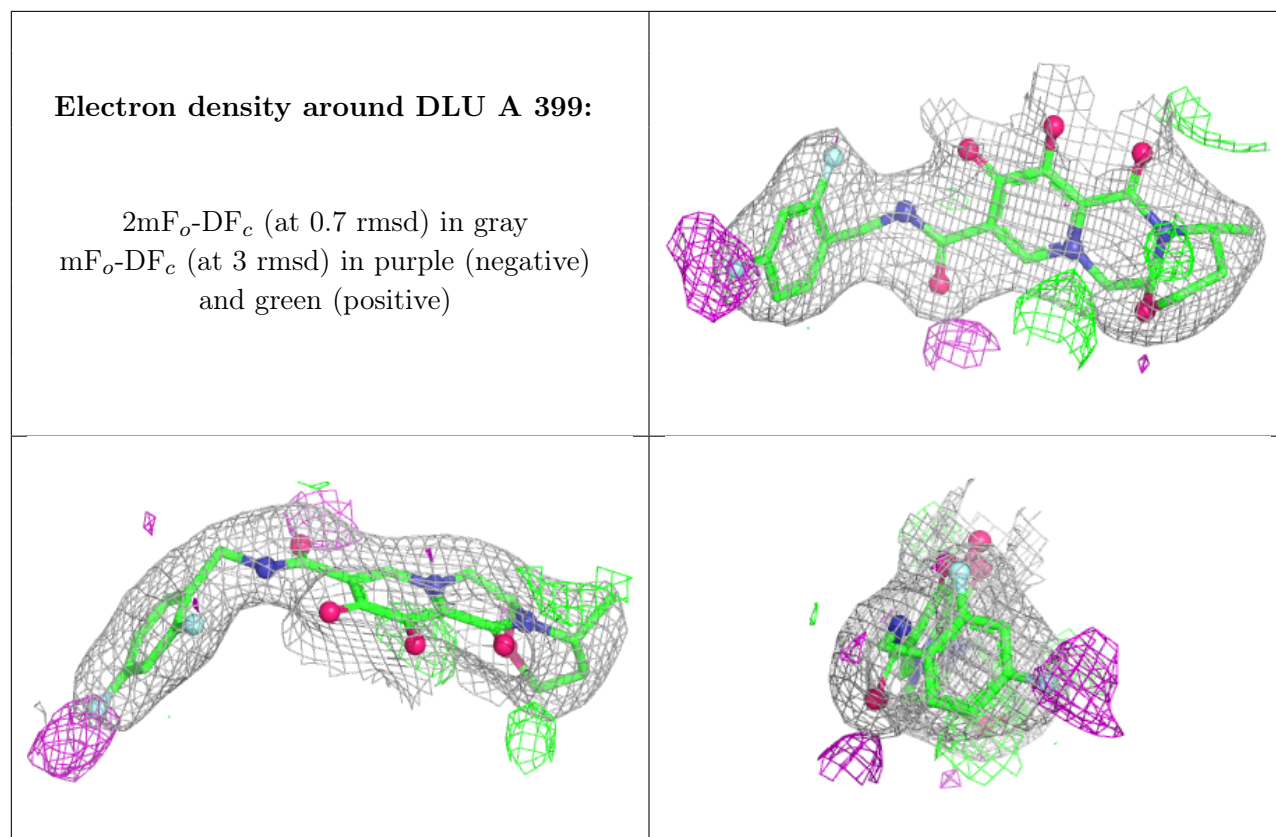
There are no oligosaccharides in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
6	GOL	A	802	6/6	0.58	0.31	66,72,76,76	0
6	GOL	A	801	6/6	0.75	0.30	78,81,81,81	0
6	GOL	A	805	6/6	0.87	0.20	86,88,88,90	0
10	HEZ	B	394	8/8	0.87	0.16	60,62,64,65	0
7	NH4	A	395	1/1	0.91	0.32	32,32,32,32	0
6	GOL	A	803	6/6	0.93	0.17	74,75,75,78	0
5	SO4	B	393	5/5	0.93	0.15	86,86,87,88	0
5	SO4	A	398	5/5	0.94	0.12	84,86,86,86	0
5	SO4	A	394	5/5	0.95	0.10	74,76,78,78	0
9	DLU	A	399	30/30	0.97	0.06	29,36,40,40	0
8	MG	A	397	1/1	1.00	0.04	35,35,35,35	0
4	ZN	A	393	1/1	1.00	0.01	46,46,46,46	0
8	MG	A	396	1/1	1.00	0.03	35,35,35,35	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.