



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:05 AM UTC

PDB ID : 8CYU / pdb_00008cyu
Title : Crystal structure of SARS-CoV-2 Mpro with compound C5
Authors : Worrall, L.J.; Lee, J.; Strynadka, N.C.J.
Deposited on : 2022-05-24
Resolution : 1.80 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

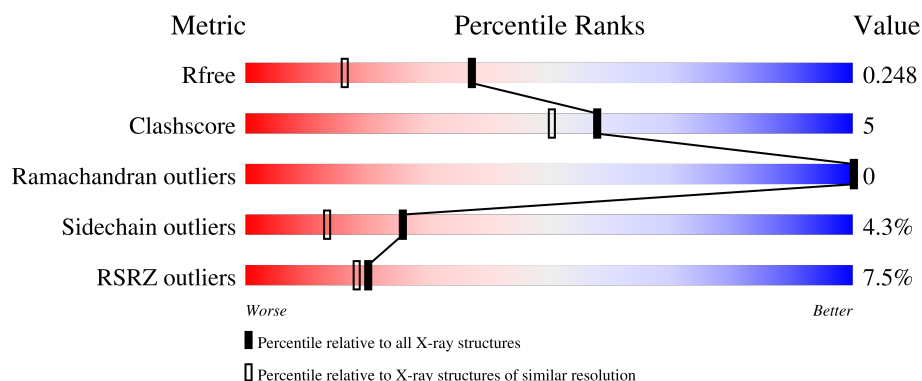
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	306	5% 86% 12% .
1	B	306	7% 88% 11% .
1	C	306	12% 86% 12% .
1	D	306	6% 86% 12% .

2 Entry composition [i](#)

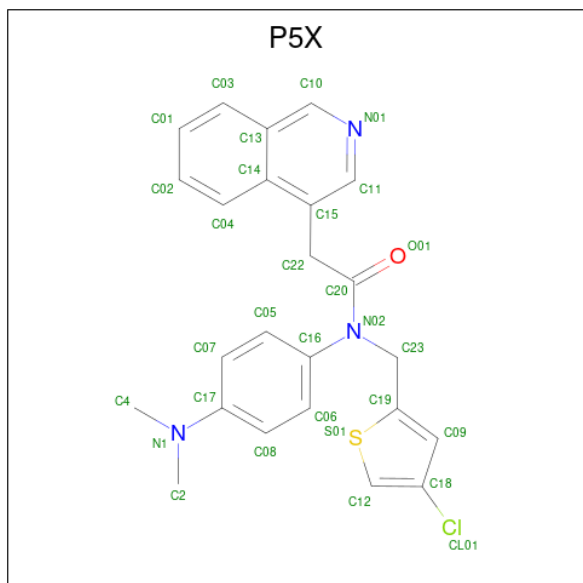
There are 3 unique types of molecules in this entry. The entry contains 10298 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 3C-like proteinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	3	0
			2378	1506	403	447	22			
1	B	306	Total	C	N	O	S	0	3	0
			2378	1507	401	447	23			
1	C	305	Total	C	N	O	S	0	4	0
			2386	1511	404	448	23			
1	D	306	Total	C	N	O	S	0	2	0
			2370	1502	400	446	22			

- Molecule 2 is N-[(4-chlorothiophen-2-yl)methyl]-N-[4-(dimethylamino)phenyl]-2-(isoquinolin-4-yl)acetamide (CCD ID: P5X) (formula: C₂₄H₂₂ClN₃OS) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	S	0	0
			30	24	1	3	1	1		

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	B	1	Total 30	C 24	Cl 1	N 3	O 1	S 1	0	0
2	C	1	Total 30	C 24	Cl 1	N 3	O 1	S 1	0	0
2	D	1	Total 30	C 24	Cl 1	N 3	O 1	S 1	0	0

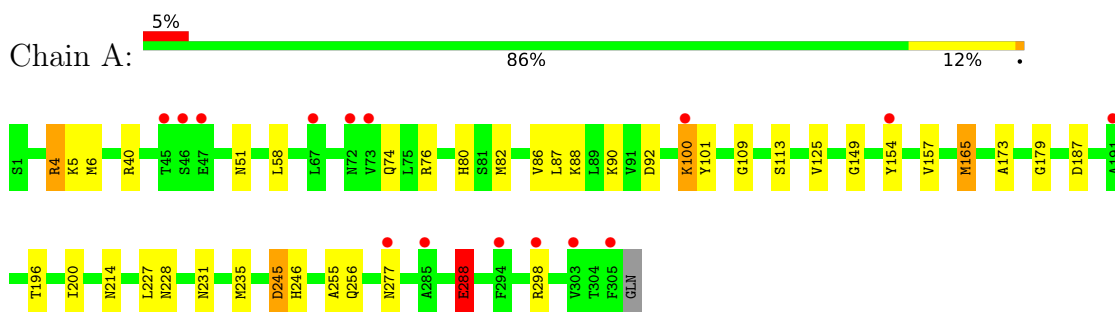
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	179	Total 179	O 179	0	0
3	B	152	Total 152	O 152	0	0
3	C	140	Total 140	O 140	0	0
3	D	195	Total 195	O 195	0	0

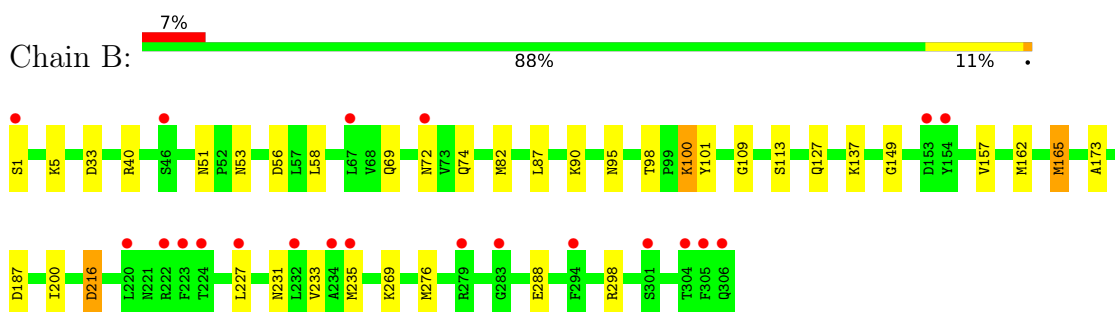
3 Residue-property plots [i](#)

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.

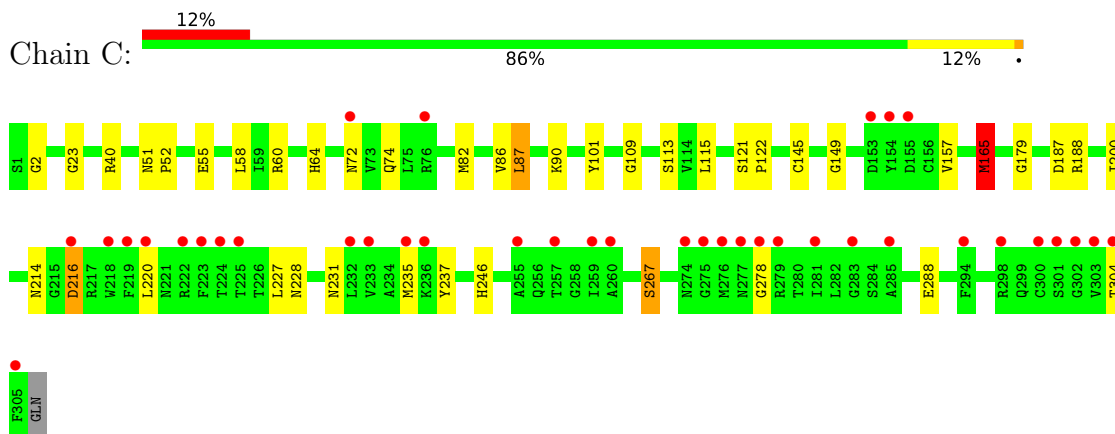
- Molecule 1: 3C-like proteinase



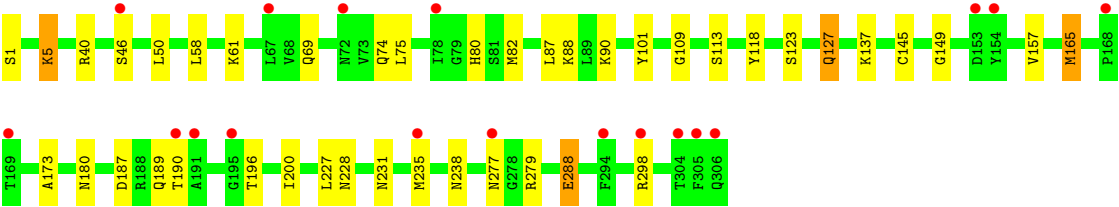
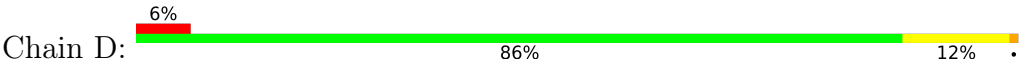
- Molecule 1: 3C-like proteinase



- Molecule 1: 3C-like proteinase



- Molecule 1: 3C-like proteinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.29Å 102.64Å 102.20Å 90.00° 100.25° 90.00°	Depositor
Resolution (Å)	45.75 – 1.80 45.75 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.75-1.80) 99.9 (45.75-1.80)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.54 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.212 , 0.245 0.223 , 0.248	Depositor DCC
R_{free} test set	6360 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.444	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10298	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.09 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.7943e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: P5X

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	1.14	4/2432 (0.2%)	1.32	5/3306 (0.2%)
1	B	1.10	1/2432 (0.0%)	1.29	2/3306 (0.1%)
1	C	1.08	4/2440 (0.2%)	1.30	3/3316 (0.1%)
1	D	1.13	1/2424 (0.0%)	1.30	6/3296 (0.2%)
All	All	1.11	10/9728 (0.1%)	1.30	16/13224 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	246	HIS	CE1-NE2	6.69	1.39	1.32
1	A	125	VAL	C-O	6.44	1.30	1.24
1	A	4	ARG	NE-CZ	5.56	1.39	1.33
1	C	165	MET	CG-SD	5.45	1.94	1.80
1	D	75	LEU	C-O	5.42	1.30	1.24
1	A	6	MET	C-O	5.17	1.30	1.24
1	C	87	LEU	C-O	5.17	1.30	1.24
1	C	246	HIS	CE1-NE2	5.13	1.37	1.32
1	C	64	HIS	CE1-NE2	5.10	1.37	1.32
1	B	162	MET	C-O	5.01	1.29	1.24

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	214	ASN	CA-CB-CG	-6.28	106.32	112.60
1	C	216	ASP	CA-CB-CG	6.22	118.82	112.60
1	B	33	ASP	CA-CB-CG	5.77	118.37	112.60
1	D	127	GLN	CB-CA-C	-5.66	99.93	109.50
1	C	2	GLY	CA-C-O	-5.57	118.12	122.52
1	D	196	THR	CA-CB-OG1	-5.56	101.27	109.60
1	A	245	ASP	CA-CB-CG	-5.54	107.06	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	216	ASP	CA-CB-CG	5.44	118.04	112.60
1	A	288	GLU	CB-CG-CD	5.43	121.84	112.60
1	A	228	ASN	CA-CB-CG	-5.32	107.28	112.60
1	D	238	ASN	CB-CA-C	5.31	119.28	111.73
1	D	190	THR	CB-CA-C	5.30	119.25	109.71
1	D	228	ASN	CA-CB-CG	-5.28	107.32	112.60
1	C	214	ASN	CA-CB-CG	-5.12	107.48	112.60
1	D	123	SER	N-CA-C	-5.11	106.73	113.12
1	A	196	THR	CA-CB-OG1	-5.06	102.01	109.60

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	2378	0	2322	25	0
1	B	2378	0	2315	18	0
1	C	2386	0	2330	28	0
1	D	2370	0	2307	24	0
2	A	30	0	0	3	0
2	B	30	0	0	1	0
2	C	30	0	0	1	0
2	D	30	0	0	1	0
3	A	179	0	0	4	0
3	B	152	0	0	4	0
3	C	140	0	0	5	0
3	D	195	0	0	11	1
All	All	10298	0	9274	86	1

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 5.

All (86) 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:256:GLN:N	1:C:72:ASN:OD1	1.94	0.99
1:D:288:GLU:HG2	3:D:517:HOH:O	1.73	0.88
1:A:288:GLU:HG2	3:A:525:HOH:O	1.78	0.81
1:A:255:ALA:C	1:C:72:ASN:OD1	2.24	0.80
1:B:56:ASP:HB2	3:B:504:HOH:O	1.85	0.77
1:C:145:CYS:SG	3:C:616:HOH:O	2.43	0.75
1:D:231:ASN:HB3	1:D:235:MET:HE3	1.68	0.75
1:B:72:ASN:OD1	1:C:60:ARG:HA	1.89	0.72
1:D:58:LEU:HD22	1:D:82:MET:HE3	1.71	0.71
1:A:58:LEU:HD22	1:A:82:MET:HE3	1.73	0.71
1:B:231:ASN:HB3	1:B:235:MET:HE3	1.76	0.68
1:C:237:TYR:OH	3:C:501:HOH:O	2.11	0.68
1:B:165:MET:HG2	1:B:173:ALA:HB3	1.76	0.67
1:D:137:LYS:HE2	3:D:637:HOH:O	1.95	0.66
1:C:231:ASN:HB3	1:C:235:MET:HE3	1.78	0.66
1:A:298:ARG:HD3	3:A:508:HOH:O	1.95	0.66
1:C:58:LEU:HD22	1:C:82:MET:HE3	1.78	0.64
1:D:165:MET:HG2	1:D:173:ALA:HB3	1.79	0.63
1:A:165:MET:HE2	2:A:401:P5X:CL01	2.36	0.62
1:B:58:LEU:HD22	1:B:82:MET:HE3	1.81	0.62
1:D:298:ARG:HD3	3:D:507:HOH:O	1.99	0.61
1:A:231:ASN:HB3	1:A:235:MET:HE3	1.80	0.61
1:C:72:ASN:HB2	3:C:617:HOH:O	2.01	0.59
1:D:61:LYS:HE2	3:D:558:HOH:O	2.02	0.59
1:A:277:ASN:HB3	3:A:647:HOH:O	2.04	0.58
1:B:100:LYS:NZ	1:B:100:LYS:HB3	2.18	0.58
1:B:40:ARG:HB2	1:B:82:MET:HE2	1.86	0.56
1:A:165:MET:HG2	1:A:173:ALA:HB3	1.87	0.55
1:C:52:PRO:HD2	1:C:188:ARG:HG2	1.89	0.55
1:B:233:VAL:HG21	1:B:269:LYS:HD2	1.89	0.54
1:A:255:ALA:HB3	1:C:72:ASN:OD1	2.08	0.52
1:A:255:ALA:CB	1:C:72:ASN:OD1	2.58	0.52
1:D:5:LYS:HE3	3:D:666:HOH:O	2.10	0.51
1:D:40:ARG:HB2	1:D:82:MET:HE2	1.92	0.51
1:A:100:LYS:HB3	1:A:100:LYS:NZ	2.26	0.50
1:B:109:GLY:HA2	1:B:200:ILE:HD13	1.93	0.50
1:B:53:ASN:ND2	3:B:504:HOH:O	2.45	0.50
1:C:40:ARG:HB2	1:C:82:MET:HE2	1.94	0.50
1:D:145:CYS:SG	3:D:639:HOH:O	2.43	0.50
1:A:154:TYR:HB2	1:C:23:GLY:HA2	1.94	0.49
1:A:231:ASN:O	1:A:235:MET:HG3	2.14	0.48
1:A:113:SER:O	1:A:149:GLY:HA2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:MET:HE2	2:A:401:P5X:C18	2.43	0.48
1:A:76:ARG:NH1	1:A:92:ASP:OD2	2.47	0.48
1:B:137:LYS:HE3	3:B:634:HOH:O	2.14	0.48
1:A:109:GLY:HA2	1:A:200:ILE:HD13	1.96	0.47
1:D:187:ASP:HA	2:D:401:P5X:CL01	2.52	0.46
1:C:231:ASN:O	1:C:235:MET:HG3	2.15	0.46
1:D:109:GLY:HA2	1:D:200:ILE:HD13	1.97	0.46
1:B:113:SER:OG	1:B:127:GLN:NE2	2.49	0.46
1:A:40:ARG:HB2	1:A:82:MET:HE2	1.98	0.46
1:C:187:ASP:HA	2:C:401:P5X:CL01	2.53	0.46
1:D:113:SER:O	1:D:149:GLY:HA2	2.15	0.46
1:C:72:ASN:CB	3:C:617:HOH:O	2.61	0.46
1:C:109:GLY:HA2	1:C:200:ILE:HD13	1.97	0.46
1:C:86:VAL:HG13	1:C:179:GLY:HA2	1.99	0.45
1:D:5:LYS:HD2	3:D:504:HOH:O	2.16	0.45
1:C:304:THR:CG2	1:D:118:TYR:HB2	2.47	0.44
1:C:278:GLY:HA3	3:C:586:HOH:O	2.18	0.44
1:D:101:TYR:HA	1:D:157:VAL:O	2.18	0.44
1:A:101:TYR:HA	1:A:157:VAL:O	2.18	0.44
1:B:113:SER:O	1:B:149:GLY:HA2	2.18	0.44
1:B:101:TYR:HA	1:B:157:VAL:O	2.18	0.43
1:C:113:SER:O	1:C:149:GLY:HA2	2.17	0.43
1:A:86:VAL:HG13	1:A:179:GLY:HA2	2.00	0.42
1:B:5:LYS:HE2	3:B:579:HOH:O	2.19	0.42
1:D:46:SER:HB2	3:D:653:HOH:O	2.19	0.42
1:D:113:SER:OG	1:D:127:GLN:NE2	2.52	0.42
1:C:101:TYR:HA	1:C:157:VAL:O	2.20	0.42
1:D:180:ASN:ND2	3:D:506:HOH:O	2.48	0.42
1:A:256:GLN:CA	1:C:72:ASN:OD1	2.67	0.42
1:D:277:ASN:HB3	3:D:524:HOH:O	2.20	0.42
1:D:279:ARG:HD3	3:D:655:HOH:O	2.20	0.42
1:B:231:ASN:O	1:B:235:MET:HG3	2.19	0.41
1:A:187:ASP:HA	2:A:401:P5X:CL01	2.57	0.41
1:A:80:HIS:HA	1:A:88:LYS:O	2.20	0.41
1:C:165:MET:HE2	1:C:165:MET:HB2	1.99	0.41
1:C:304:THR:HG21	1:D:118:TYR:HB2	2.03	0.41
1:C:220:LEU:HA	1:C:267:SER:OG	2.21	0.41
1:C:121:SER:HA	1:C:122:PRO:HD3	1.96	0.41
1:D:80:HIS:HA	1:D:88:LYS:O	2.21	0.41
1:A:51:ASN:ND2	3:A:506:HOH:O	2.49	0.41
1:B:187:ASP:HA	2:B:401:P5X:CL01	2.58	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:ASN:HB3	1:B:98:THR:OG1	2.21	0.40
1:C:115:LEU:HD11	1:C:122:PRO:HB3	2.02	0.40
1:D:231:ASN:O	1:D:235:MET:HG3	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:602:HOH:O	3:D:647:HOH:O[1_655]	2.11	0.09

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	306/306 (100%)	299 (98%)	7 (2%)	0	100	100
1	B	307/306 (100%)	302 (98%)	5 (2%)	0	100	100
1	C	307/306 (100%)	301 (98%)	6 (2%)	0	100	100
1	D	306/306 (100%)	299 (98%)	7 (2%)	0	100	100
All	All	1226/1224 (100%)	1201 (98%)	25 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	265/263 (101%)	255 (96%)	10 (4%)	29	17
1	B	264/263 (100%)	250 (95%)	14 (5%)	20	9
1	C	266/263 (101%)	255 (96%)	11 (4%)	27	15
1	D	263/263 (100%)	252 (96%)	11 (4%)	26	14
All	All	1058/1052 (101%)	1012 (96%)	46 (4%)	26	13

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	5	LYS
1	A	74	GLN
1	A	87	LEU
1	A	90	LYS
1	A	100	LYS
1	A	165	MET
1	A	227	LEU
1	A	245	ASP
1	A	288	GLU
1	B	1	SER
1	B	51	ASN
1	B	69	GLN
1	B	74	GLN
1	B	87	LEU
1	B	90	LYS
1	B	100	LYS
1	B	165	MET
1	B	216	ASP
1	B	227	LEU
1	B	276[A]	MET
1	B	276[B]	MET
1	B	288	GLU
1	B	298	ARG
1	C	51	ASN
1	C	55	GLU
1	C	74	GLN
1	C	87	LEU
1	C	90	LYS
1	C	165	MET
1	C	216	ASP
1	C	227	LEU
1	C	228	ASN

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Mol	Chain	Res	Type
1	C	267	SER
1	C	288	GLU
1	D	1	SER
1	D	5	LYS
1	D	50	LEU
1	D	69	GLN
1	D	74	GLN
1	D	87	LEU
1	D	90	LYS
1	D	165	MET
1	D	189	GLN
1	D	227	LEU
1	D	288	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	HIS
1	B	41	HIS
1	B	127	GLN
1	B	189	GLN
1	C	41	HIS
1	C	51	ASN
1	C	84	ASN
1	C	228	ASN
1	D	41	HIS
1	D	51	ASN
1	D	65	ASN
1	D	72	ASN
1	D	119	ASN
1	D	127	GLN
1	D	180	ASN
1	D	189	GLN

5.3.3 RNA ⓘ

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](#)

4 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$
2	P5X	C	401	-	33,33,33	1.32	2 (6%)	43,46,46	3.02	15 (34%)
2	P5X	D	401	-	33,33,33	1.28	3 (9%)	43,46,46	3.05	14 (32%)
2	P5X	B	401	-	33,33,33	1.37	2 (6%)	43,46,46	2.68	10 (23%)
2	P5X	A	401	-	33,33,33	1.45	3 (9%)	43,46,46	2.75	14 (32%)

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	P5X	C	401	-	-	2/20/20/20	0/4/4/4
2	P5X	D	401	-	-	2/20/20/20	0/4/4/4
2	P5X	B	401	-	-	2/20/20/20	0/4/4/4
2	P5X	A	401	-	-	2/20/20/20	0/4/4/4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	P5X	C18-CL01	5.29	1.83	1.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	P5X	C18-CL01	4.73	1.82	1.73
2	B	401	P5X	C18-CL01	4.55	1.82	1.73
2	C	401	P5X	C18-CL01	4.12	1.81	1.73
2	A	401	P5X	C10-C13	-3.04	1.36	1.42
2	D	401	P5X	C10-C13	-2.89	1.36	1.42
2	B	401	P5X	C02-C04	2.41	1.41	1.36
2	A	401	P5X	C02-C04	2.13	1.41	1.36
2	C	401	P5X	C22-C20	2.11	1.54	1.52
2	D	401	P5X	C17-N1	2.07	1.42	1.37

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	401	P5X	C18-C12-S01	-11.97	102.99	110.87
2	C	401	P5X	C18-C12-S01	-11.78	103.12	110.87
2	B	401	P5X	C18-C12-S01	-9.89	104.36	110.87
2	A	401	P5X	C18-C12-S01	-9.78	104.44	110.87
2	B	401	P5X	C09-C18-C12	8.53	120.89	107.54
2	C	401	P5X	C09-C18-C12	8.52	120.88	107.54
2	D	401	P5X	C09-C18-C12	8.28	120.51	107.54
2	A	401	P5X	C09-C18-C12	7.60	119.44	107.54
2	D	401	P5X	C12-S01-C19	6.60	98.13	92.56
2	D	401	P5X	C12-C18-CL01	-6.08	119.39	126.58
2	A	401	P5X	C19-C23-N02	-5.93	106.35	112.92
2	D	401	P5X	C19-C23-N02	-5.45	106.88	112.92
2	B	401	P5X	C19-C23-N02	-5.39	106.94	112.92
2	C	401	P5X	C12-C18-CL01	-5.10	120.55	126.58
2	A	401	P5X	C12-C18-CL01	-4.99	120.67	126.58
2	A	401	P5X	C12-S01-C19	4.83	96.64	92.56
2	C	401	P5X	C10-N01-C11	4.67	121.47	117.03
2	C	401	P5X	C12-S01-C19	4.49	96.35	92.56
2	C	401	P5X	C23-N02-C20	3.86	122.67	117.14
2	B	401	P5X	C23-N02-C20	3.69	122.44	117.14
2	B	401	P5X	C12-C18-CL01	-3.67	122.24	126.58
2	C	401	P5X	C19-C23-N02	-3.67	108.86	112.92
2	B	401	P5X	C09-C18-CL01	-3.59	116.81	125.45
2	A	401	P5X	C23-N02-C20	3.57	122.27	117.14
2	C	401	P5X	C15-C11-N01	-3.28	120.27	124.57
2	B	401	P5X	C10-N01-C11	3.20	120.07	117.03
2	B	401	P5X	C12-S01-C19	3.10	95.17	92.56
2	D	401	P5X	C23-N02-C20	3.07	121.54	117.14
2	C	401	P5X	C09-C18-CL01	-2.86	118.57	125.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	P5X	C07-C17-N1	-2.85	117.84	121.65
2	A	401	P5X	C10-N01-C11	2.84	119.73	117.03
2	B	401	P5X	C22-C15-C11	-2.76	117.79	122.20
2	C	401	P5X	C23-N02-C16	-2.68	113.41	118.38
2	D	401	P5X	C15-C11-N01	-2.60	121.16	124.57
2	D	401	P5X	O01-C20-N02	-2.55	118.15	121.70
2	D	401	P5X	C09-C19-S01	-2.53	107.14	111.38
2	A	401	P5X	O01-C20-N02	-2.40	118.35	121.70
2	C	401	P5X	C10-C13-C14	2.39	120.16	117.84
2	D	401	P5X	C07-C17-N1	-2.39	118.46	121.65
2	A	401	P5X	C09-C18-CL01	-2.32	119.86	125.45
2	D	401	P5X	C10-N01-C11	2.28	119.19	117.03
2	A	401	P5X	C22-C15-C11	-2.26	118.58	122.20
2	C	401	P5X	C23-C19-C09	2.26	134.68	127.51
2	B	401	P5X	C15-C11-N01	-2.24	121.63	124.57
2	D	401	P5X	C09-C18-CL01	-2.23	120.08	125.45
2	C	401	P5X	C04-C14-C13	2.22	120.79	117.92
2	C	401	P5X	O01-C20-N02	-2.18	118.66	121.70
2	A	401	P5X	C06-C16-N02	-2.12	117.12	120.18
2	A	401	P5X	C15-C11-N01	-2.09	121.82	124.57
2	A	401	P5X	C09-C19-S01	-2.08	107.88	111.38
2	D	401	P5X	C10-C13-C14	2.07	119.85	117.84
2	A	401	P5X	C07-C17-N1	-2.01	118.97	121.65
2	D	401	P5X	C01-C02-C04	-2.00	117.72	120.40

There are no chirality outliers.

All (8) torsion outliers are listed below:

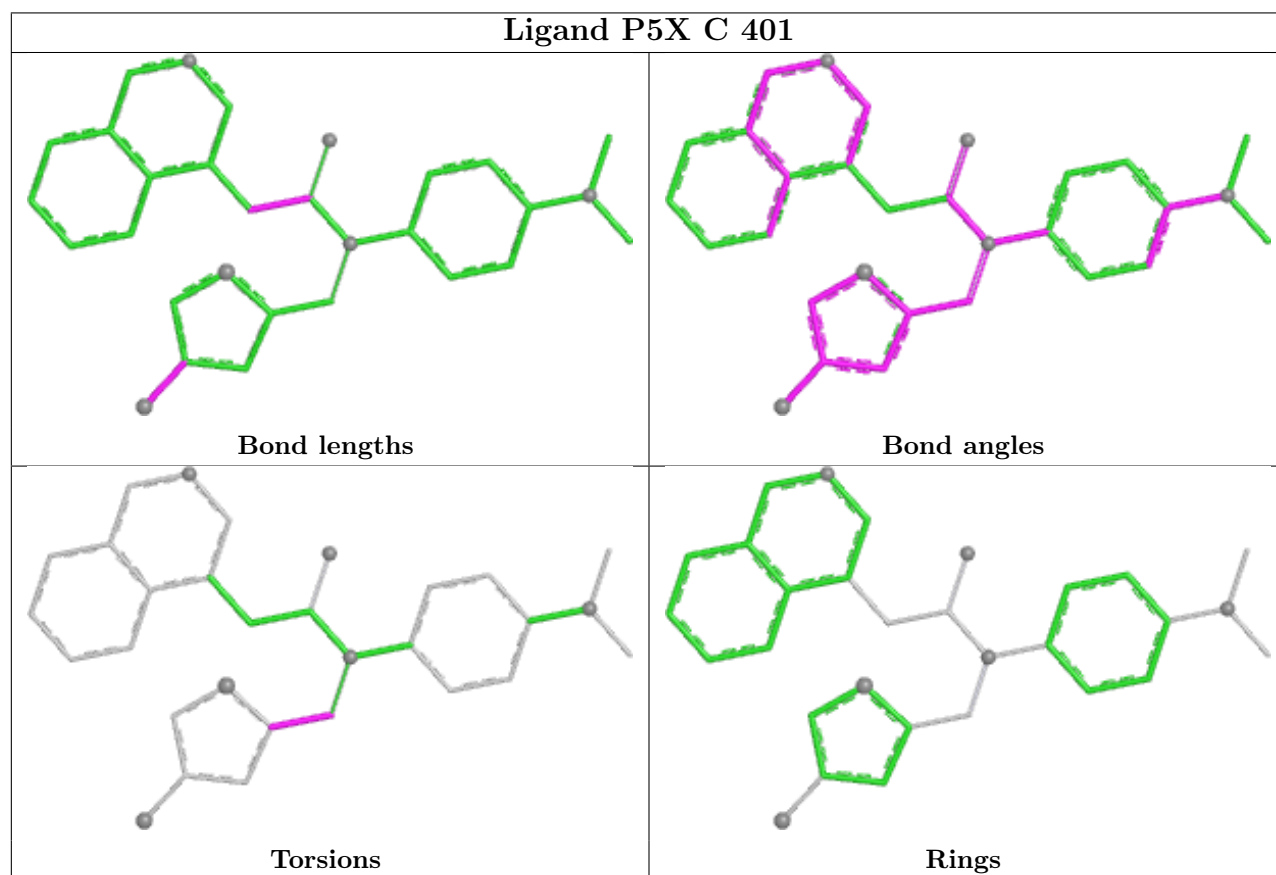
Mol	Chain	Res	Type	Atoms
2	B	401	P5X	S01-C19-C23-N02
2	C	401	P5X	S01-C19-C23-N02
2	A	401	P5X	C09-C19-C23-N02
2	B	401	P5X	C09-C19-C23-N02
2	C	401	P5X	C09-C19-C23-N02
2	D	401	P5X	C09-C19-C23-N02
2	A	401	P5X	S01-C19-C23-N02
2	D	401	P5X	S01-C19-C23-N02

There are no ring outliers.

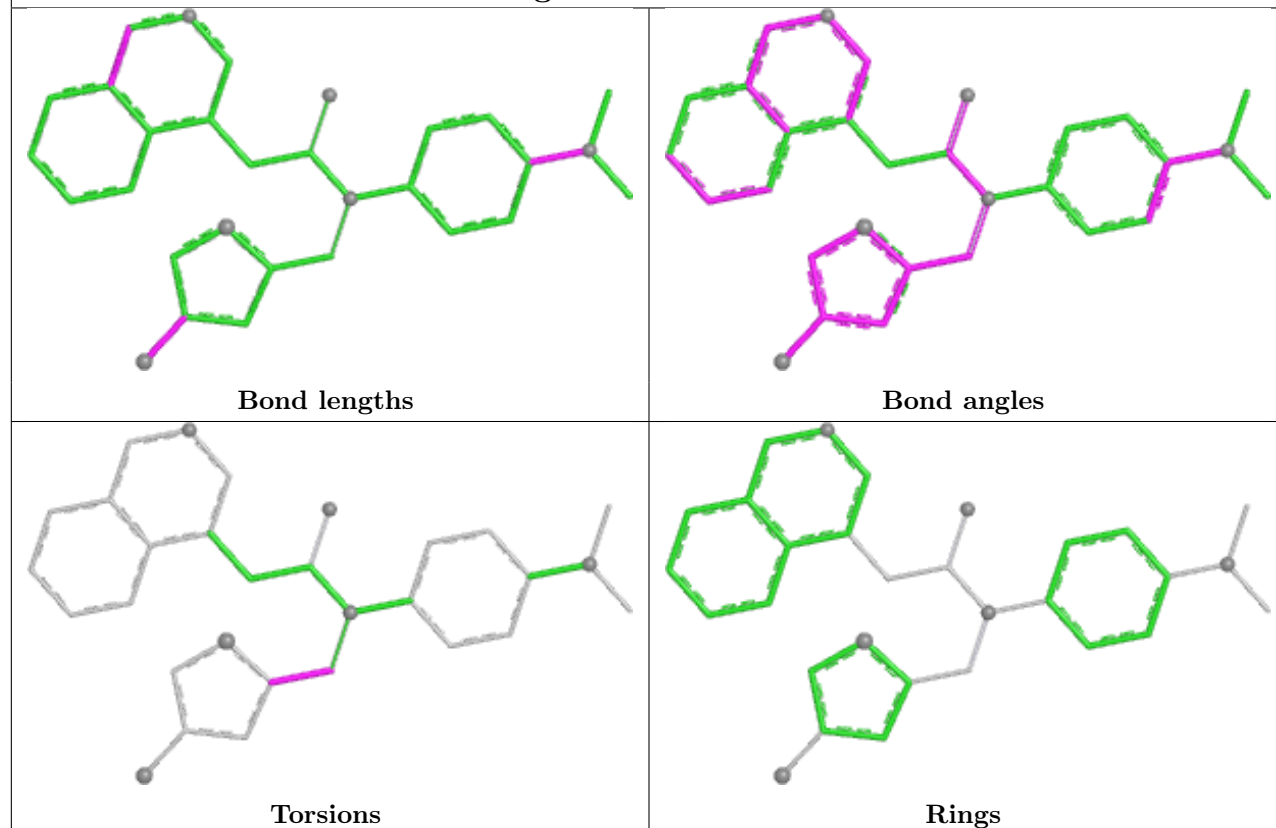
4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	P5X	1	0
2	D	401	P5X	1	0
2	B	401	P5X	1	0
2	A	401	P5X	3	0

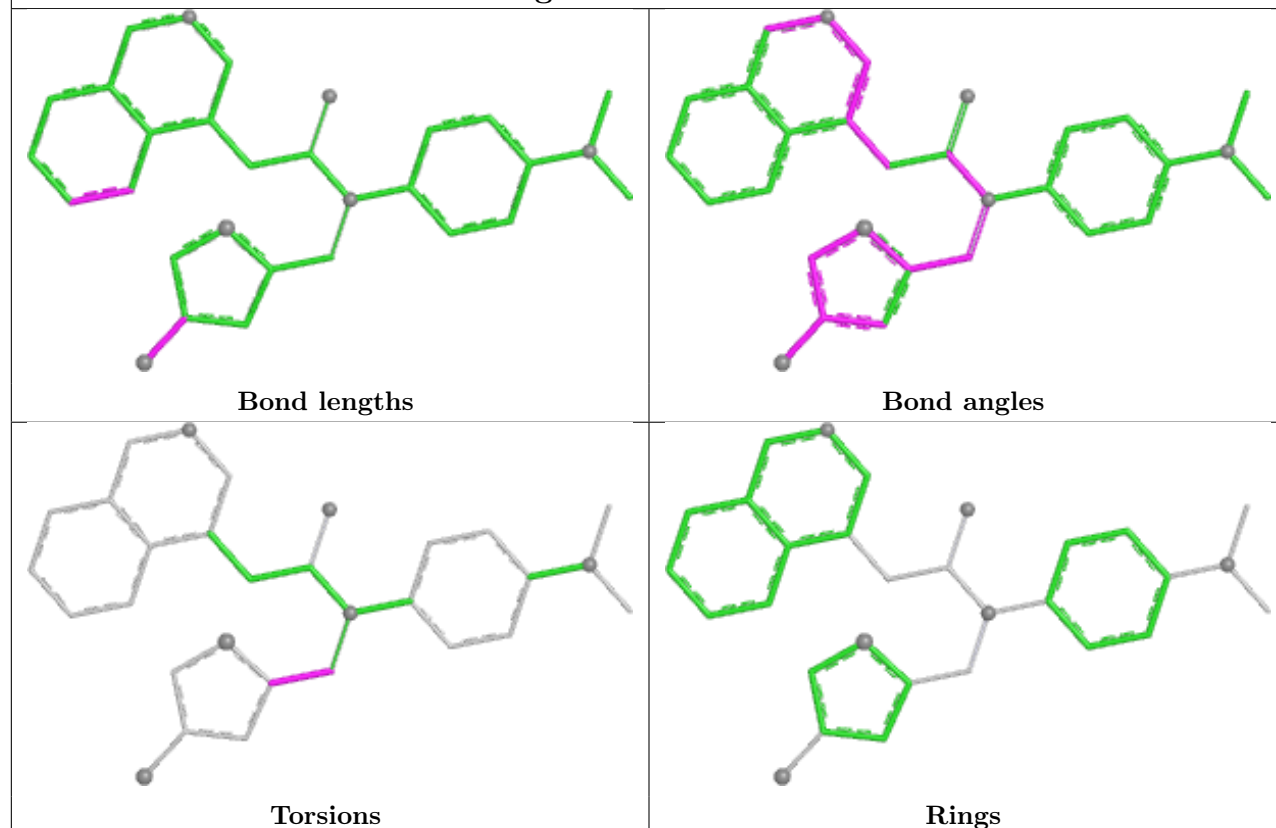
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

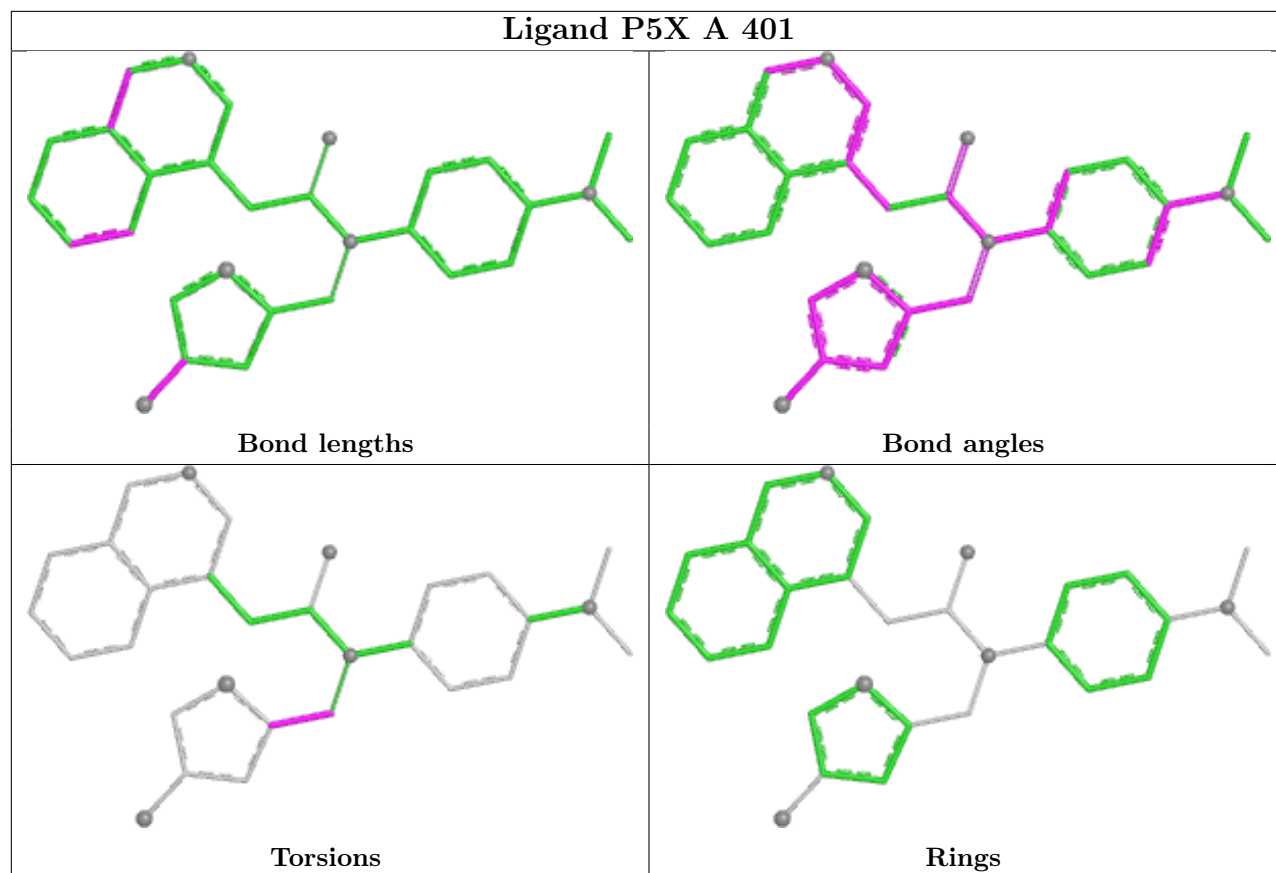


Ligand P5X D 401



Ligand P5X B 401





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	305/306 (99%)	0.22	15 (4%) 35 33	13, 28, 44, 65	3 (0%)
1	B	306/306 (100%)	0.48	21 (6%) 23 21	14, 32, 57, 88	3 (0%)
1	C	305/306 (99%)	0.72	38 (12%) 8 6	13, 34, 68, 122	4 (1%)
1	D	306/306 (100%)	0.26	18 (5%) 28 27	15, 28, 47, 68	2 (0%)
All	All	1222/1224 (99%)	0.42	92 (7%) 20 18	13, 30, 57, 122	12 (0%)

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	278	GLY	7.0
1	C	223	PHE	6.4
1	C	305	PHE	6.2
1	C	235	MET	5.7
1	B	154	TYR	5.6
1	C	302	GLY	5.2
1	C	276[A]	MET	4.9
1	B	301	SER	4.9
1	C	277	ASN	4.7
1	C	154	TYR	4.5
1	C	232	LEU	4.4
1	B	72	ASN	4.3
1	C	218	TRP	4.3
1	D	191	ALA	4.2
1	A	294	PHE	4.1
1	C	236	LYS	4.1
1	B	153	ASP	4.0
1	C	224	THR	4.0
1	D	306	GLN	4.0
1	C	72	ASN	3.9
1	C	222	ARG	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	279	ARG	3.9
1	C	220	LEU	3.8
1	C	303	VAL	3.7
1	D	305	PHE	3.5
1	B	306	GLN	3.4
1	A	305	PHE	3.4
1	B	224	THR	3.4
1	A	67	LEU	3.4
1	C	153	ASP	3.4
1	C	275	GLY	3.4
1	B	223	PHE	3.3
1	B	235	MET	3.3
1	D	294	PHE	3.3
1	B	283	GLY	3.3
1	D	46	SER	3.2
1	C	285	ALA	3.2
1	A	277	ASN	3.2
1	B	222	ARG	3.2
1	B	305	PHE	3.0
1	C	255	ALA	3.0
1	A	72	ASN	3.0
1	B	67	LEU	2.9
1	A	191	ALA	2.9
1	D	67	LEU	2.9
1	A	46	SER	2.9
1	D	154	TYR	2.9
1	D	72	ASN	2.8
1	C	301	SER	2.8
1	C	257	THR	2.8
1	C	219	PHE	2.7
1	D	195	GLY	2.7
1	B	294	PHE	2.7
1	D	235	MET	2.6
1	A	154	TYR	2.6
1	D	153	ASP	2.6
1	D	298	ARG	2.5
1	B	232	LEU	2.5
1	A	100	LYS	2.5
1	C	216	ASP	2.5
1	B	220	LEU	2.4
1	B	279	ARG	2.4
1	C	300	CYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	46	SER	2.4
1	C	225	THR	2.3
1	C	298	ARG	2.3
1	A	47	GLU	2.3
1	A	45	THR	2.3
1	C	260	ALA	2.3
1	D	78	ILE	2.3
1	A	73	VAL	2.2
1	C	283	GLY	2.2
1	C	294	PHE	2.2
1	C	281	ILE	2.2
1	B	234	ALA	2.2
1	A	303	VAL	2.1
1	C	304	THR	2.1
1	D	169	THR	2.1
1	D	190	THR	2.1
1	D	304	THR	2.1
1	A	298	ARG	2.1
1	D	168	PRO	2.1
1	D	277	ASN	2.1
1	C	76	ARG	2.1
1	C	274	ASN	2.1
1	A	285	ALA	2.1
1	C	155	ASP	2.1
1	C	259	ILE	2.0
1	B	227	LEU	2.0
1	B	1	SER	2.0
1	B	304	THR	2.0
1	C	233	VAL	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

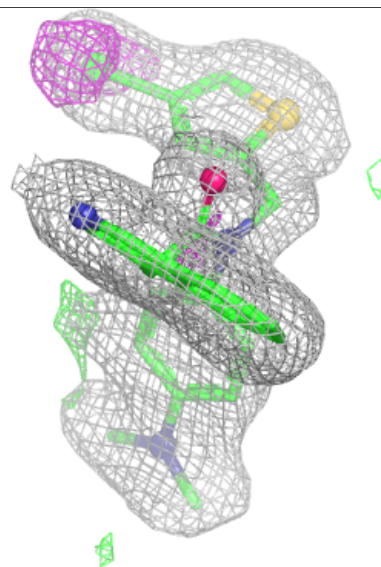
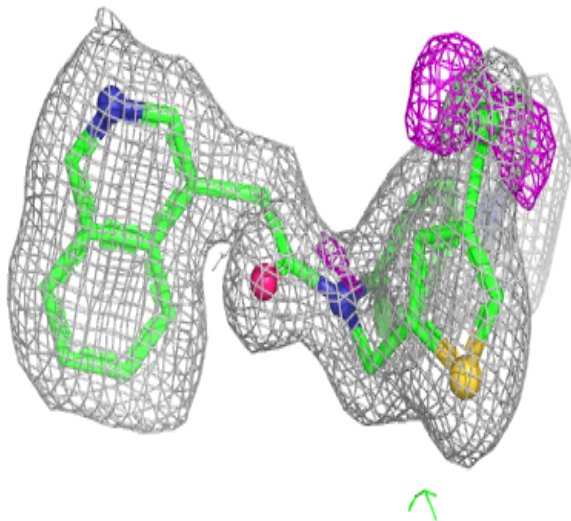
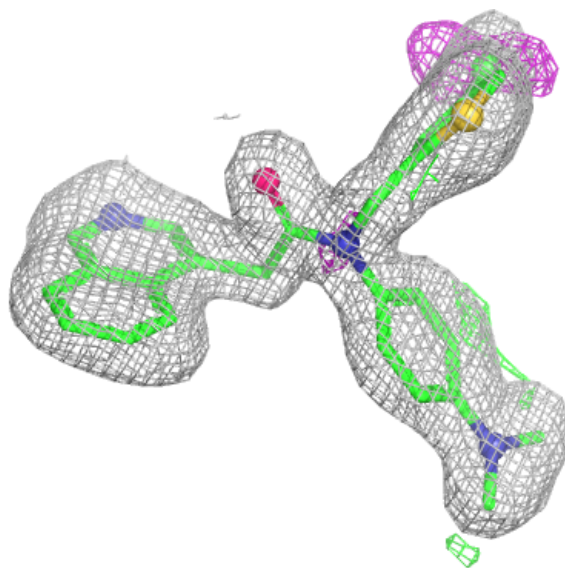
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	P5X	D	401	30/30	0.90	0.10	22,29,36,39	0
2	P5X	A	401	30/30	0.92	0.10	22,29,33,37	0
2	P5X	C	401	30/30	0.94	0.09	23,28,33,35	0
2	P5X	B	401	30/30	0.95	0.08	21,25,33,38	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.

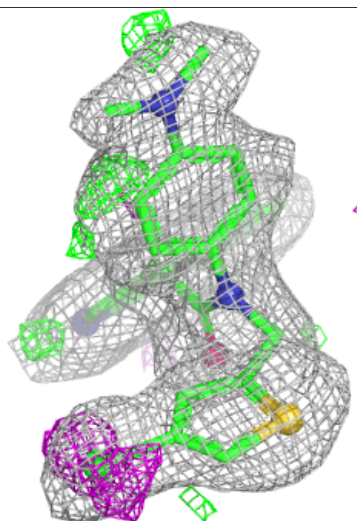
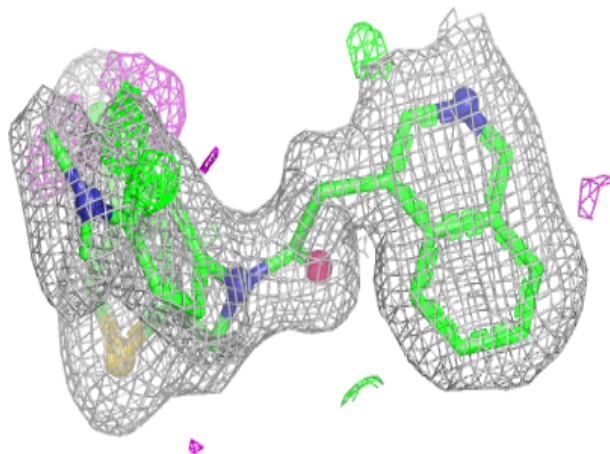
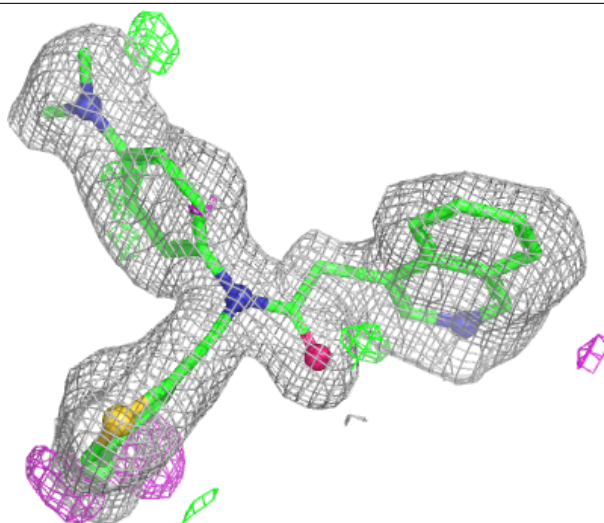
Electron density around P5X D 401:

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



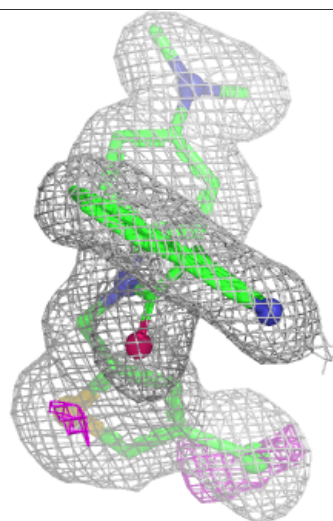
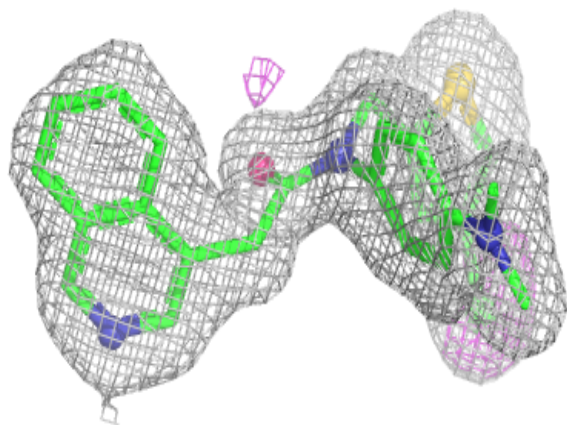
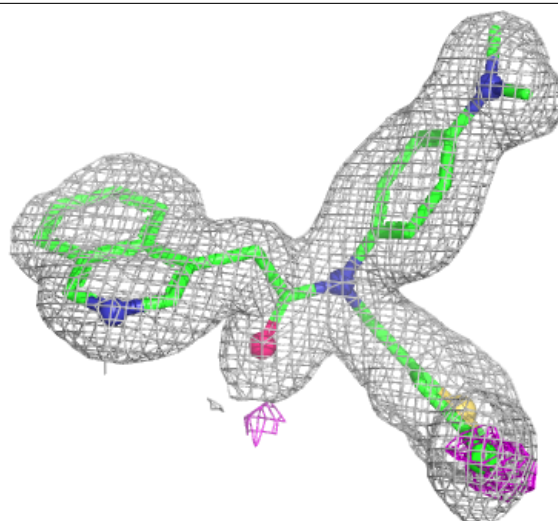
Electron density around P5X A 401:

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



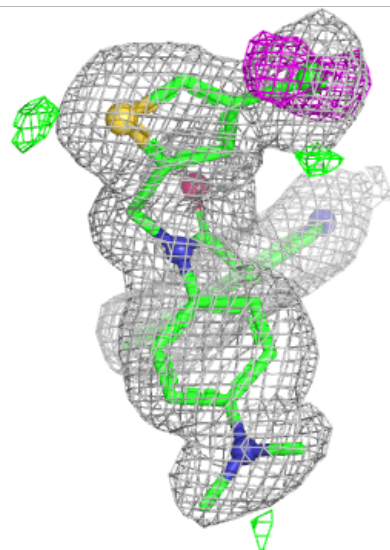
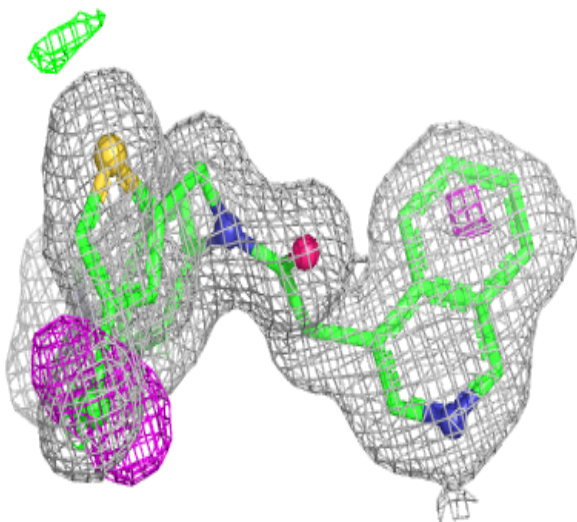
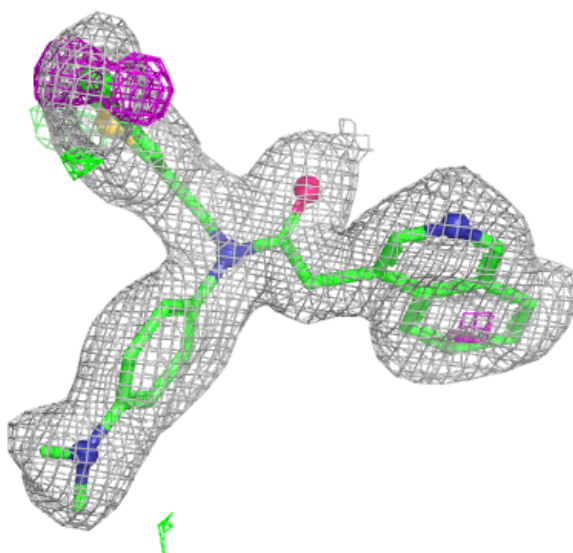
Electron density around P5X C 401:

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



Electron density around P5X B 401:

$2mF_o-DF_c$ (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.