



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:23 AM UTC

PDB ID : 2XH1 / pdb_00002xh1
Title : Crystal structure of human KAT II-inhibitor complex
Authors : Rossi, F.; Casazza, V.; Garavaglia, S.; Sathyaikumar, K.V.; Schwarcz, R.;
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Deposited on : 2010-06-08
Resolution : 2.10 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

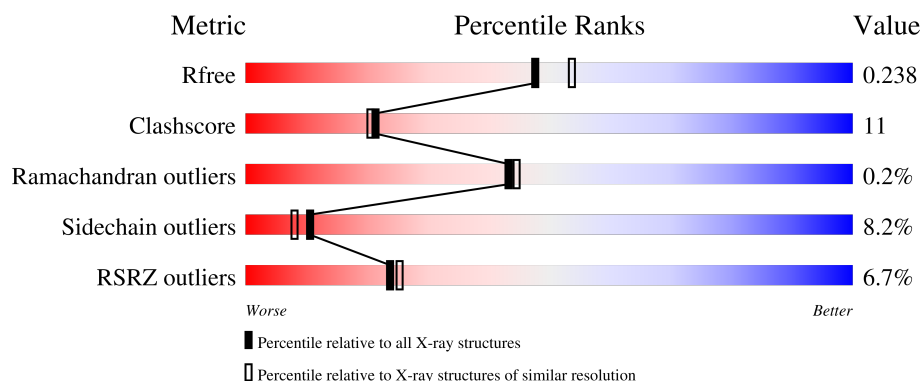
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	425	4% 77% 17% • •
1	B	425	8% 71% 21% • •

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6972 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.

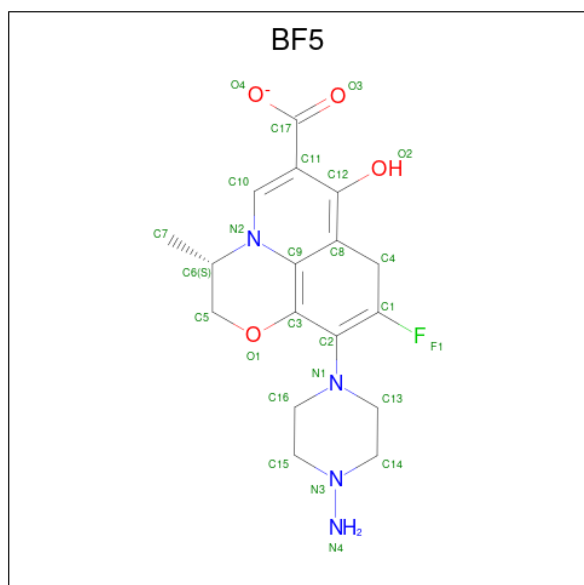
- Molecule 1 is a protein called KYNURENINE/ α -AMINOADIPATE AMINOTRANSFERASE, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	412	Total	C	N	O	S	0	0	0
			3233	2079	539	599	16			
1	B	412	Total	C	N	O	S	0	0	0
			3234	2079	539	600	16			

There are 2 discrepancies between the modelled and reference sequences:

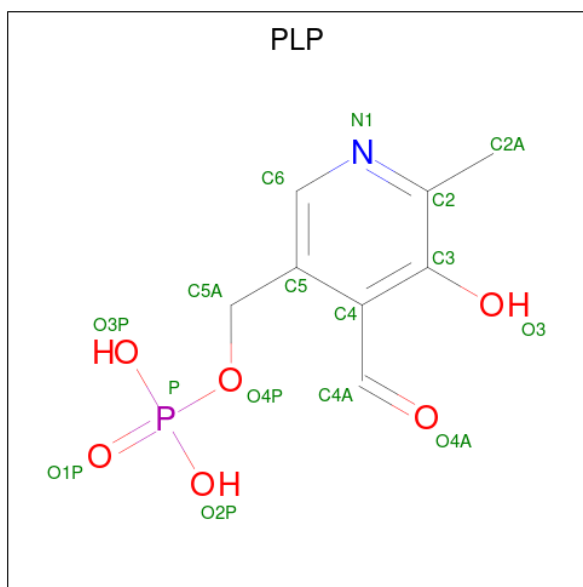
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLU	MET	engineered mutation	UNP Q8N5Z0
B	-1	GLU	MET	engineered mutation	UNP Q8N5Z0

- Molecule 2 is (3S)-10-(4-AMINOPIPERAZIN-1-YL)-9-FLUORO-7-HYDROXY-3-METHY L-2,3-DIHYDRO-8H-[1,4]OXAZINO[2,3,4-IJ]QUINOLINE-6-CARBOXYLATE (CCD ID: BF5) (formula: C₁₇H₁₉FN₄O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			26	17	1	4	4		
2	B	1	Total	C	F	N	O	0	0
			26	17	1	4	4		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).

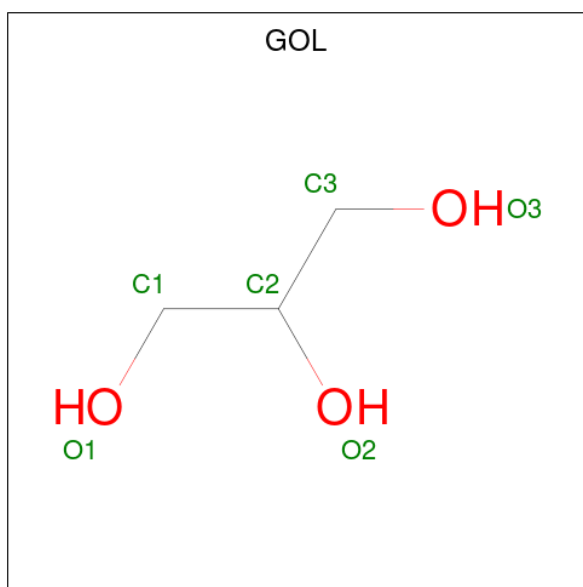


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 4 is IODIDE ION (CCD ID: IOD) (formula: I).

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

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

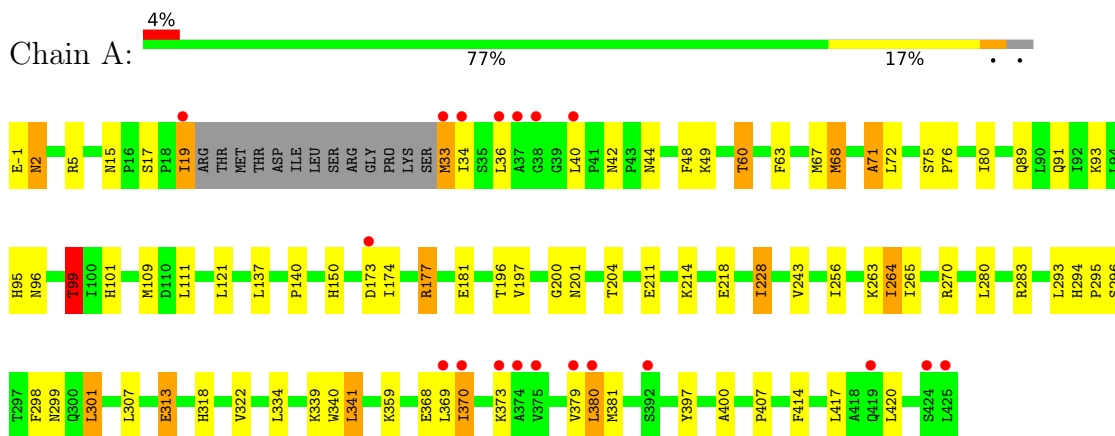
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	241	Total	O	0	0
			241	241		
6	B	175	Total	O	0	0
			175	175		

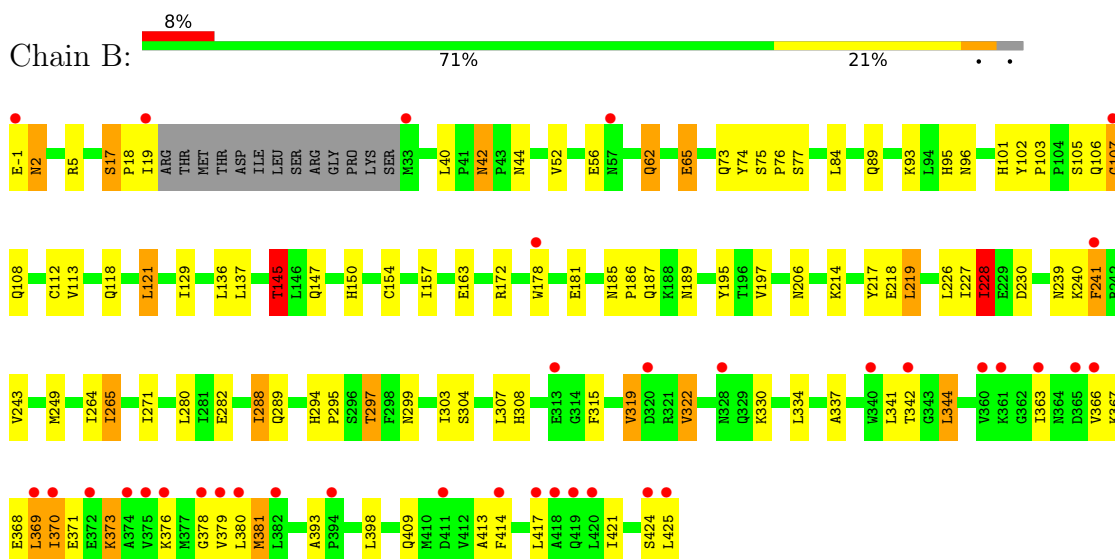
3 Residue-property plots

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: KYNURENINE/ALPHA-AMINOADIPATE AMINOTRANSFERASE, MITOCHONDRIAL



- Molecule 1: KYNURENINE/ALPHA-AMINOADIPATE AMINOTRANSFERASE, MITOCHONDRIAL



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	98.32Å 152.86Å 60.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.04 – 2.10 32.04 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (32.04-2.10) 99.9 (32.04-2.10)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.39 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.188 , 0.241 0.191 , 0.238	Depositor DCC
R_{free} test set	1111 reflections (2.05%)	wwPDB-VP
Wilson B-factor (Å ²)	24.0	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6972	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% 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: GOL, BF5, PLP, IOD

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	1.07	1/3313 (0.0%)	1.07	4/4498 (0.1%)
1	B	1.01	4/3314 (0.1%)	1.11	15/4498 (0.3%)
All	All	1.04	5/6627 (0.1%)	1.09	19/8996 (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

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	228	ILE	CA-CB	8.26	1.64	1.53
1	B	103	PRO	CA-C	6.54	1.58	1.52
1	B	107	GLY	N-CA	6.42	1.54	1.45
1	B	322	VAL	CA-CB	5.76	1.62	1.54
1	A	71	ALA	CA-CB	-5.72	1.44	1.53

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	106	GLN	CA-C-N	-9.67	102.46	121.41
1	B	106	GLN	C-N-CA	-9.67	102.46	121.41
1	A	2	ASN	N-CA-C	-8.36	98.09	110.46
1	A	264	ILE	CB-CA-C	-7.69	104.29	111.06
1	B	101	HIS	N-CA-C	5.88	119.42	112.72
1	B	5	ARG	CB-CA-C	-5.78	100.86	110.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	106	GLN	O-C-N	-5.69	114.55	121.83
1	A	99	THR	N-CA-CB	-5.56	101.72	110.46
1	B	393	ALA	CA-C-N	5.36	125.30	119.78
1	B	393	ALA	C-N-CA	5.36	125.30	119.78
1	B	288	ILE	N-CA-C	5.32	116.05	110.62
1	B	228	ILE	CA-CB-CG2	5.27	119.45	110.50
1	B	2	ASN	CB-CA-C	5.26	118.51	109.51
1	B	145	THR	N-CA-CB	5.22	117.97	110.20
1	B	265	ILE	CB-CA-C	5.20	115.68	111.05
1	A	265	ILE	N-CA-C	5.12	115.67	111.62
1	B	228	ILE	CB-CG1-CD1	-5.04	103.21	113.80
1	B	227	ILE	CA-C-N	-5.01	116.29	123.11
1	B	227	ILE	C-N-CA	-5.01	116.29	123.11

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	-1	GLU	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	3233	0	3241	70	1
1	B	3234	0	3241	78	0
2	A	26	0	17	4	0
2	B	26	0	17	1	0
3	A	15	0	6	0	0
3	B	15	0	7	0	0
4	A	1	0	0	0	0
5	B	6	0	8	0	0
6	A	241	0	0	15	0
6	B	175	0	0	8	0
All	All	6972	0	6537	142	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 11.

All (142) 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:B:89:GLN:HG3	1:B:93:LYS:HE2	1.42	0.99
1:B:112:CYS:HB2	6:B:2076:HOH:O	1.66	0.94
1:A:99:THR:HG21	6:A:2073:HOH:O	1.71	0.89
1:B:112:CYS:SG	6:B:2076:HOH:O	2.36	0.82
1:A:196:THR:HG22	6:A:2170:HOH:O	1.80	0.81
1:A:121:LEU:CD2	1:A:228:ILE:HD11	2.15	0.76
1:A:368:GLU:HG3	6:A:2232:HOH:O	1.87	0.74
1:A:63:PHE:HD2	1:A:68:MET:HE1	1.52	0.73
1:B:121:LEU:CD1	1:B:228:ILE:HD11	2.20	0.72
1:B:112:CYS:CB	6:B:2076:HOH:O	2.28	0.71
1:B:172:ARG:HG2	1:B:219:LEU:HD11	1.71	0.71
1:B:379:VAL:HG11	1:B:413:ALA:HA	1.71	0.71
1:A:264:ILE:O	1:A:318:HIS:HE1	1.75	0.69
1:B:145:THR:HG21	1:B:195:TYR:CZ	2.29	0.67
1:A:5:ARG:HD3	6:A:2004:HOH:O	1.94	0.67
1:B:294:HIS:HD2	1:B:295:PRO:O	1.78	0.66
1:A:63:PHE:HD2	1:A:68:MET:CE	2.09	0.65
1:A:34:ILE:HD13	1:B:409:GLN:HB3	1.79	0.65
1:B:271:ILE:HG13	1:B:303:ILE:HD12	1.78	0.65
1:B:42:ASN:HD22	1:B:44:ASN:H	1.44	0.65
1:B:73:GLN:O	1:B:297:THR:HG21	1.97	0.64
1:A:67:MET:HE3	6:A:2201:HOH:O	1.98	0.63
1:A:370:ILE:HD11	1:A:380:LEU:HA	1.78	0.63
1:A:299:ASN:HD21	1:B:299:ASN:HD21	1.46	0.63
1:A:68:MET:CE	1:A:68:MET:HA	2.28	0.63
1:B:185:ASN:C	1:B:185:ASN:HD22	2.07	0.62
1:A:381:MET:HE3	1:A:400:ALA:HB2	1.81	0.62
1:B:121:LEU:HD11	1:B:228:ILE:HD11	1.80	0.62
1:A:173:ASP:OD1	1:A:173:ASP:C	2.41	0.62
1:A:80:ILE:CD1	1:A:301:LEU:HD13	2.30	0.62
1:B:344:LEU:HD23	1:B:344:LEU:H	1.65	0.62
1:A:283:ARG:HD3	6:A:2192:HOH:O	1.99	0.61
1:A:174:ILE:O	1:A:177:ARG:HD3	2.01	0.61
2:A:1426:BF5:H4	1:B:40:LEU:CD2	2.31	0.61
1:B:369:LEU:HA	1:B:373:LYS:HG3	1.82	0.61
1:A:17:SER:C	1:A:19:ILE:H	2.09	0.60
1:B:84:LEU:HD11	1:B:113:VAL:HG23	1.84	0.60
1:B:89:GLN:CG	1:B:93:LYS:HE2	2.26	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:LEU:HD12	1:B:381:MET:HE3	1.83	0.60
1:B:214:LYS:O	1:B:218:GLU:HG3	2.02	0.59
1:A:99:THR:HG22	1:A:109:MET:HB2	1.83	0.59
1:A:101:HIS:HD2	6:A:2068:HOH:O	1.84	0.59
1:B:304:SER:O	1:B:308:HIS:HD2	1.85	0.58
1:B:218:GLU:CD	6:B:2130:HOH:O	2.46	0.58
1:B:195:TYR:OH	1:B:230:ASP:OD2	2.17	0.58
1:A:341:LEU:HD22	1:A:414:PHE:HD2	1.69	0.58
1:B:370:ILE:HG13	1:B:398:LEU:HD21	1.85	0.58
1:A:2:ASN:HA	6:A:2071:HOH:O	2.03	0.58
1:B:344:LEU:HD23	1:B:344:LEU:N	2.19	0.58
1:A:75:SER:HB2	1:A:76:PRO:CD	2.33	0.57
1:B:421:ILE:O	1:B:425:LEU:HG	2.04	0.57
1:B:217:TYR:HB2	1:B:249:MET:CE	2.35	0.57
2:A:1426:BF5:H4	1:B:40:LEU:HD21	1.85	0.56
1:B:84:LEU:CD1	1:B:113:VAL:HG23	2.35	0.55
1:A:89:GLN:HE21	1:A:93:LYS:HG3	1.70	0.55
1:B:42:ASN:ND2	1:B:44:ASN:H	2.04	0.55
1:B:239:ASN:O	1:B:241:PHE:N	2.40	0.54
1:B:52:VAL:HG22	1:B:62:GLN:NE2	2.21	0.54
2:A:1426:BF5:H16A	2:A:1426:BF5:F1	1.97	0.54
1:A:67:MET:CE	6:A:2201:HOH:O	2.57	0.53
1:B:18:PRO:O	1:B:19:ILE:HB	2.08	0.53
1:B:136:LEU:HD23	1:B:157:ILE:HB	1.91	0.53
1:A:294:HIS:HD2	1:A:295:PRO:O	1.90	0.52
1:B:121:LEU:HD21	1:B:230:ASP:CG	2.33	0.52
1:A:121:LEU:HD22	1:A:228:ILE:HD11	1.90	0.52
1:B:17:SER:HB2	1:B:18:PRO:O	2.10	0.52
1:A:256:ILE:HD11	1:A:280:LEU:HD13	1.91	0.51
1:A:341:LEU:HD22	1:A:414:PHE:CD2	2.44	0.51
1:A:68:MET:HA	1:A:68:MET:HE2	1.93	0.51
1:A:15:ASN:O	1:B:147:GLN:NE2	2.43	0.51
1:A:369:LEU:HA	1:A:373:LYS:HE2	1.92	0.51
1:A:293:LEU:HD21	1:B:118:GLN:HE21	1.75	0.50
1:A:80:ILE:HD13	1:A:301:LEU:HD13	1.93	0.50
1:A:283:ARG:CD	6:A:2192:HOH:O	2.56	0.50
1:B:367:LYS:HE3	1:B:371:GLU:OE1	2.12	0.50
1:A:63:PHE:CD2	1:A:68:MET:CE	2.93	0.50
1:A:173:ASP:OD1	1:A:174:ILE:N	2.45	0.49
1:A:295:PRO:O	1:A:296:SER:C	2.55	0.49
1:B:2:ASN:ND2	6:B:2173:HOH:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121:LEU:HD12	1:B:228:ILE:HD11	1.93	0.49
1:B:145:THR:HG21	1:B:195:TYR:OH	2.12	0.49
1:A:121:LEU:HD23	1:A:228:ILE:HD11	1.94	0.49
2:A:1426:BF5:F1	2:A:1426:BF5:C16	2.51	0.49
1:B:185:ASN:ND2	1:B:187:GLN:H	2.11	0.48
1:A:68:MET:HA	1:A:68:MET:HE3	1.95	0.48
1:A:313:GLU:H	1:A:313:GLU:CD	2.20	0.48
1:B:73:GLN:O	1:B:297:THR:CG2	2.60	0.48
1:B:74:TYR:HA	1:B:294:HIS:CE1	2.48	0.48
1:A:197:VAL:HB	1:A:201:ASN:HD22	1.79	0.48
1:A:95:HIS:O	1:A:96:ASN:C	2.57	0.48
1:B:337:ALA:O	1:B:341:LEU:HB2	2.14	0.47
1:A:150:HIS:HE1	6:B:2012:HOH:O	1.96	0.47
1:A:299:ASN:HD21	1:B:299:ASN:ND2	2.12	0.47
1:B:102:TYR:O	1:B:108:GLN:HB2	2.14	0.47
1:B:264:ILE:HG13	1:B:265:ILE:HG13	1.97	0.46
1:A:42:ASN:OD1	1:A:44:ASN:HB2	2.16	0.46
1:A:99:THR:HG23	1:A:109:MET:N	2.31	0.45
1:A:214:LYS:O	1:A:218:GLU:HG3	2.15	0.45
1:B:2:ASN:N	6:B:2003:HOH:O	2.47	0.45
1:A:95:HIS:HE1	6:A:2185:HOH:O	2.00	0.45
1:B:-1:GLU:O	1:B:105:SER:O	2.34	0.45
1:A:68:MET:HE1	1:A:298:PHE:CE2	2.52	0.45
1:B:107:GLY:O	1:B:108:GLN:C	2.59	0.45
1:A:48:PHE:HB2	1:A:72:LEU:HD11	1.99	0.45
1:B:304:SER:O	1:B:308:HIS:CD2	2.69	0.45
1:B:181:GLU:CD	1:B:181:GLU:H	2.24	0.45
1:B:42:ASN:HD21	1:B:44:ASN:HD22	1.65	0.45
1:A:264:ILE:O	1:A:318:HIS:CE1	2.65	0.44
1:B:77:SER:OG	1:B:289:GLN:HG2	2.17	0.44
1:A:283:ARG:HD2	1:A:283:ARG:N	2.32	0.44
1:A:200:GLY:O	1:A:201:ASN:C	2.61	0.44
1:B:42:ASN:HD22	1:B:42:ASN:C	2.24	0.44
1:B:95:HIS:O	1:B:96:ASN:C	2.60	0.44
1:A:60:THR:HG22	6:A:2025:HOH:O	2.17	0.44
1:A:68:MET:HE2	1:A:71:ALA:HB3	2.00	0.44
1:A:270:ARG:HD3	1:B:294:HIS:CE1	2.52	0.44
1:B:217:TYR:HB2	1:B:249:MET:HE2	2.00	0.44
1:A:359:LYS:HB2	1:A:397:TYR:CE2	2.53	0.44
6:A:2007:HOH:O	1:B:150:HIS:HE1	2.01	0.44
1:A:95:HIS:HD2	6:A:2180:HOH:O	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:SER:HB2	1:B:76:PRO:CD	2.48	0.43
1:B:315:PHE:O	1:B:319:VAL:HG12	2.18	0.43
1:A:33:MET:HG3	1:A:34:ILE:N	2.34	0.42
1:A:49:LYS:HE2	1:B:56:GLU:HB2	2.00	0.42
1:A:339:LYS:HE2	1:A:340:TRP:CZ2	2.54	0.42
1:B:17:SER:C	1:B:18:PRO:O	2.61	0.42
1:B:206:ASN:ND2	6:B:2125:HOH:O	2.40	0.42
1:B:294:HIS:NE2	1:B:297:THR:HG22	2.35	0.42
1:B:239:ASN:C	1:B:241:PHE:H	2.28	0.41
1:B:65:GLU:H	1:B:65:GLU:HG2	1.60	0.41
1:B:185:ASN:HD22	1:B:186:PRO:N	2.17	0.41
1:A:80:ILE:HD12	1:A:301:LEU:HD13	2.01	0.41
1:A:137:LEU:C	1:A:137:LEU:HD12	2.45	0.41
1:B:366:VAL:HG21	1:B:398:LEU:HG	2.03	0.41
1:A:140:PRO:HG2	1:A:204:THR:HG23	2.03	0.41
1:A:368:GLU:CG	6:A:2232:HOH:O	2.59	0.41
1:B:197:VAL:HA	1:B:230:ASP:O	2.20	0.41
1:B:341:LEU:HD22	1:B:414:PHE:CD2	2.56	0.41
1:A:40:LEU:HD21	2:B:1426:BF5:H4	2.02	0.40
1:B:129:ILE:CG2	1:B:154:CYS:HB3	2.51	0.40
1:B:341:LEU:HD22	1:B:414:PHE:HD2	1.85	0.40
1:A:91:GLN:HA	1:A:91:GLN:NE2	2.36	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)
1:A:211:GLU:OE1	1:A:211:GLU:OE1[2_565]	1.73	0.47

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	408/425 (96%)	395 (97%)	13 (3%)	0	100	100
1	B	408/425 (96%)	390 (96%)	16 (4%)	2 (0%)	24	22
All	All	816/850 (96%)	785 (96%)	29 (4%)	2 (0%)	43	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	240	LYS
1	B	378	GLY

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	358/370 (97%)	335 (94%)	23 (6%)	16	14
1	B	358/370 (97%)	322 (90%)	36 (10%)	7	5
All	All	716/740 (97%)	657 (92%)	59 (8%)	10	8

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	ILE
1	A	33	MET
1	A	60	THR
1	A	68	MET
1	A	99	THR
1	A	111	LEU
1	A	177	ARG
1	A	181	GLU
1	A	228	ILE
1	A	243	VAL
1	A	263	LYS
1	A	301	LEU
1	A	307	LEU
1	A	313	GLU

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Mol	Chain	Res	Type
1	A	322	VAL
1	A	334	LEU
1	A	341	LEU
1	A	370	ILE
1	A	379	VAL
1	A	380	LEU
1	A	407	PRO
1	A	417	LEU
1	A	420	LEU
1	B	17	SER
1	B	42	ASN
1	B	62	GLN
1	B	65	GLU
1	B	121	LEU
1	B	137	LEU
1	B	145	THR
1	B	163	GLU
1	B	178	TRP
1	B	189	ASN
1	B	219	LEU
1	B	226	LEU
1	B	228	ILE
1	B	241	PHE
1	B	243	VAL
1	B	280	LEU
1	B	282	GLU
1	B	288	ILE
1	B	297	THR
1	B	307	LEU
1	B	319	VAL
1	B	322	VAL
1	B	330	LYS
1	B	334	LEU
1	B	342	THR
1	B	344	LEU
1	B	363	ILE
1	B	368	GLU
1	B	369	LEU
1	B	370	ILE
1	B	373	LYS
1	B	376	LYS
1	B	380	LEU

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Mol	Chain	Res	Type
1	B	381	MET
1	B	417	LEU
1	B	424	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	GLN
1	A	91	GLN
1	A	95	HIS
1	A	150	HIS
1	A	201	ASN
1	A	206	ASN
1	A	294	HIS
1	A	305	GLN
1	A	318	HIS
1	A	328	ASN
1	A	348	HIS
1	B	2	ASN
1	B	42	ASN
1	B	57	ASN
1	B	62	GLN
1	B	106	GLN
1	B	118	GLN
1	B	150	HIS
1	B	185	ASN
1	B	201	ASN
1	B	237	GLN
1	B	294	HIS
1	B	299	ASN
1	B	308	HIS
1	B	328	ASN
1	B	348	HIS

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](#)

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

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	BF5	B	1426	3	25,29,29	2.29	6 (24%)	23,44,44	1.49	2 (8%)
3	PLP	B	1427	2	15,15,16	1.04	0	21,22,23	1.48	4 (19%)
3	PLP	A	1427	2	15,15,16	0.91	1 (6%)	21,22,23	1.67	1 (4%)
5	GOL	B	1428	-	5,5,5	0.27	0	5,5,5	0.95	0
2	BF5	A	1426	3	25,29,29	2.48	5 (20%)	23,44,44	1.58	5 (21%)

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	BF5	B	1426	3	-	5/8/43/43	0/4/4/4
3	PLP	B	1427	2	-	3/6/6/8	0/1/1/1
3	PLP	A	1427	2	-	3/6/6/8	0/1/1/1
5	GOL	B	1428	-	-	2/4/4/4	-
2	BF5	A	1426	3	-	5/8/43/43	0/4/4/4

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1426	BF5	C12-C8	9.17	1.50	1.36
2	B	1426	BF5	C12-C8	8.17	1.49	1.36
2	A	1426	BF5	C4-C8	-5.31	1.37	1.50
2	B	1426	BF5	C4-C8	-4.64	1.39	1.50
2	B	1426	BF5	C9-N2	3.34	1.43	1.38
2	A	1426	BF5	C9-N2	2.74	1.42	1.38
2	B	1426	BF5	O2-C12	-2.44	1.25	1.33
2	A	1426	BF5	O4-C17	-2.38	1.24	1.30
2	B	1426	BF5	O3-C17	2.17	1.27	1.22
2	B	1426	BF5	C11-C17	2.10	1.52	1.48
3	A	1427	PLP	C5A-C5	2.07	1.56	1.50
2	A	1426	BF5	C2-N1	2.04	1.43	1.36

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1427	PLP	O4P-C5A-C5	5.14	118.99	109.36
2	B	1426	BF5	C13-N1-C2	-4.02	108.41	121.03
2	A	1426	BF5	C13-N1-C2	-3.98	108.55	121.03
2	A	1426	BF5	C14-C13-N1	3.08	116.53	110.42
3	B	1427	PLP	O4P-C5A-C5	3.02	115.01	109.36
3	B	1427	PLP	C3-C4-C5	2.66	121.78	118.59
2	A	1426	BF5	C10-C11-C17	2.61	122.43	118.27
2	B	1426	BF5	C16-C15-N3	-2.56	106.70	109.53
2	A	1426	BF5	C16-N1-C2	-2.48	113.26	121.03
3	B	1427	PLP	C6-C5-C4	-2.10	116.37	118.10
2	A	1426	BF5	C13-C14-N3	2.09	111.84	109.53
3	B	1427	PLP	O3P-P-O4P	-2.01	101.44	106.67

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1426	BF5	C3-C2-N1-C16
2	B	1426	BF5	C10-C11-C17-O3
2	B	1426	BF5	C10-C11-C17-O4
2	B	1426	BF5	C12-C11-C17-O3
2	B	1426	BF5	C12-C11-C17-O4
3	A	1427	PLP	C4-C5-C5A-O4P
3	A	1427	PLP	C6-C5-C5A-O4P
3	A	1427	PLP	C5A-O4P-P-O2P
3	B	1427	PLP	C5A-O4P-P-O1P

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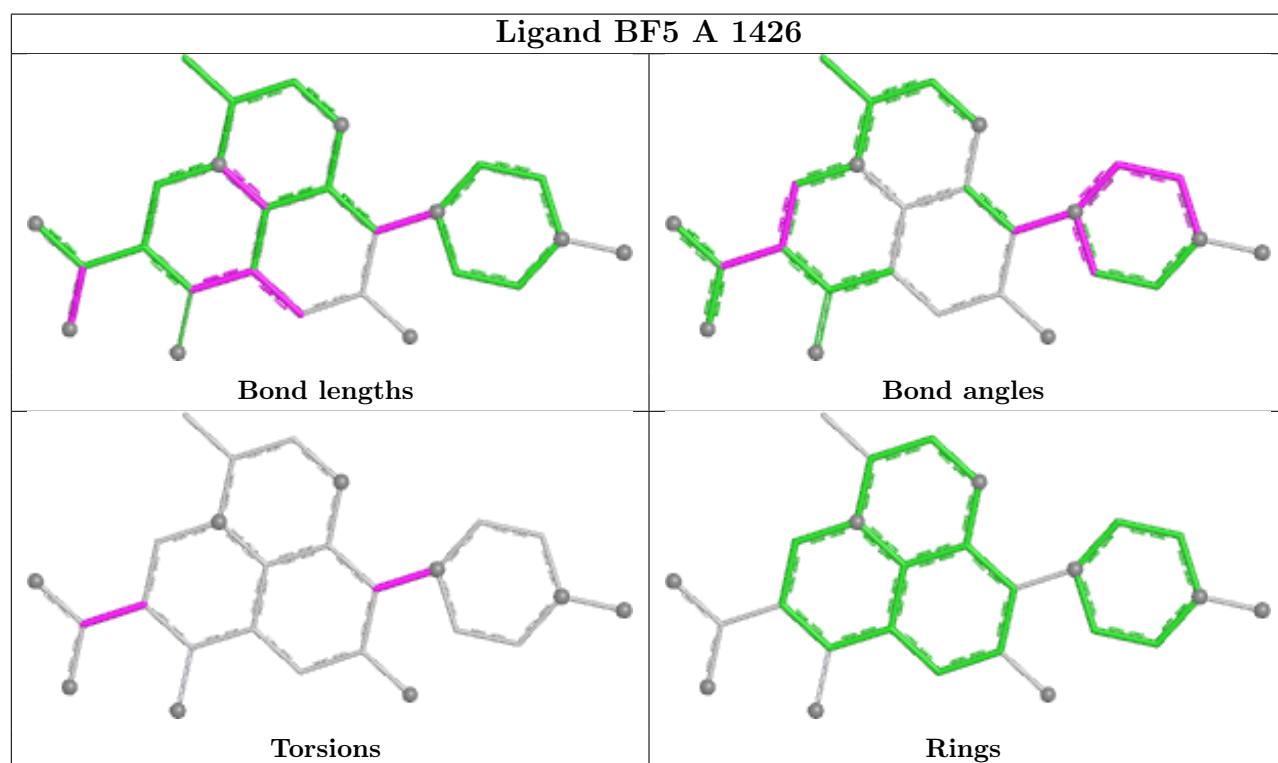
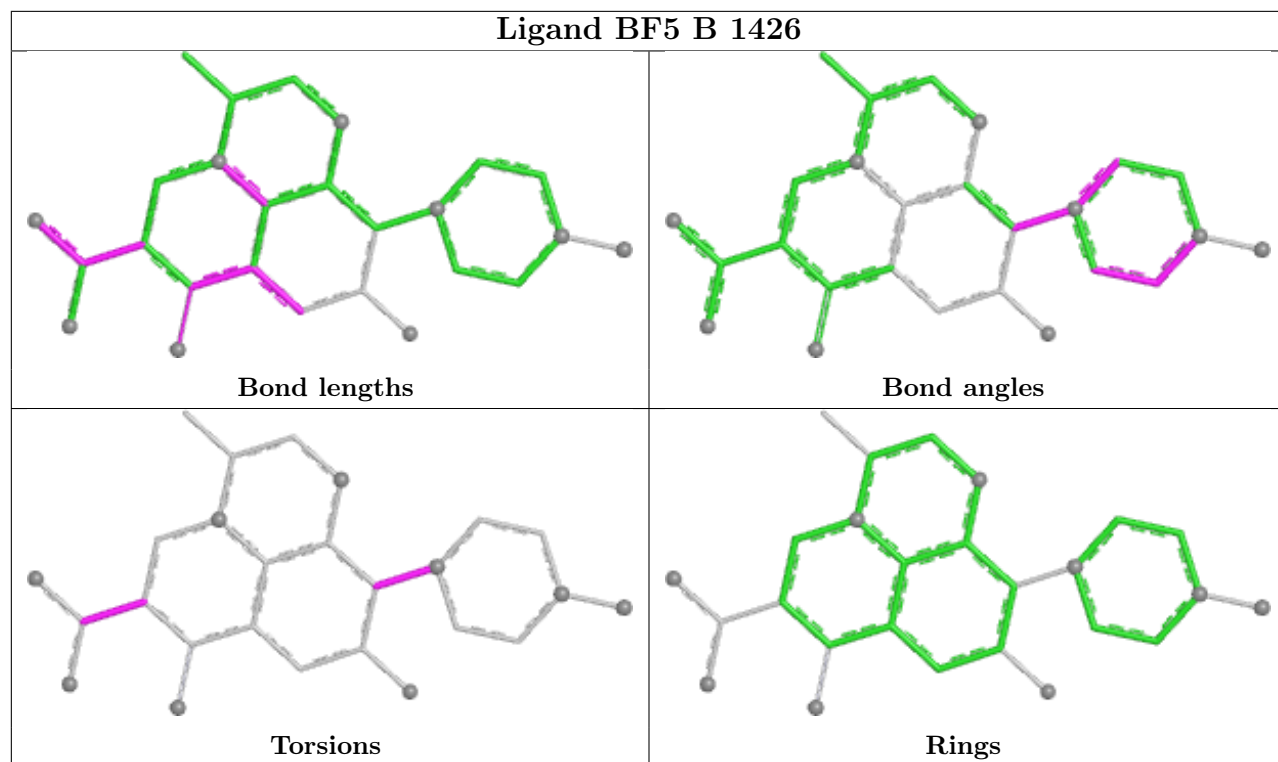
Mol	Chain	Res	Type	Atoms
3	B	1427	PLP	C5A-O4P-P-O2P
3	B	1427	PLP	C5A-O4P-P-O3P
5	B	1428	GOL	O1-C1-C2-C3
5	B	1428	GOL	O1-C1-C2-O2
2	A	1426	BF5	C12-C11-C17-O3
2	A	1426	BF5	C12-C11-C17-O4
2	A	1426	BF5	C10-C11-C17-O3
2	A	1426	BF5	C10-C11-C17-O4
2	B	1426	BF5	C3-C2-N1-C16

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1426	BF5	1	0
2	A	1426	BF5	4	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	412/425 (96%)	0.07	19 (4%) 37 39	11, 22, 52, 69	0
1	B	412/425 (96%)	0.38	36 (8%) 16 16	13, 27, 73, 91	0
All	All	824/850 (96%)	0.23	55 (6%) 24 25	11, 24, 64, 91	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	379	VAL	5.0
1	B	365	ASP	4.0
1	B	380	LEU	3.8
1	B	370	ILE	3.7
1	B	-1	GLU	3.6
1	A	19	ILE	3.6
1	A	425	LEU	3.5
1	B	411	ASP	3.3
1	B	394	PRO	3.3
1	B	379	VAL	3.3
1	B	369	LEU	3.2
1	B	425	LEU	3.2
1	B	342	THR	3.2
1	A	370	ILE	3.1
1	A	38	GLY	3.1
1	B	241	PHE	3.1
1	B	363	ILE	3.0
1	B	340	TRP	3.0
1	A	374	ALA	2.9
1	B	19	ILE	2.8
1	B	376	LYS	2.8
1	B	424	SER	2.8
1	A	40	LEU	2.7
1	B	313	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	34	ILE	2.7
1	A	36	LEU	2.7
1	B	57	ASN	2.7
1	A	392	SER	2.6
1	A	380	LEU	2.6
1	B	417	LEU	2.6
1	A	375	VAL	2.5
1	B	360	VAL	2.5
1	B	418	ALA	2.5
1	B	366	VAL	2.5
1	B	375	VAL	2.5
1	B	320	ASP	2.5
1	B	33	MET	2.5
1	A	33	MET	2.4
1	A	173	ASP	2.4
1	B	378	GLY	2.3
1	A	373	LYS	2.3
1	B	372	GLU	2.3
1	B	361	LYS	2.3
1	B	107	GLY	2.2
1	B	374	ALA	2.2
1	B	419	GLN	2.2
1	B	420	LEU	2.1
1	A	424	SER	2.1
1	B	328	ASN	2.1
1	A	369	LEU	2.1
1	B	382	LEU	2.1
1	B	414	PHE	2.1
1	B	178	TRP	2.1
1	A	37	ALA	2.1
1	A	419	GLN	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](#)

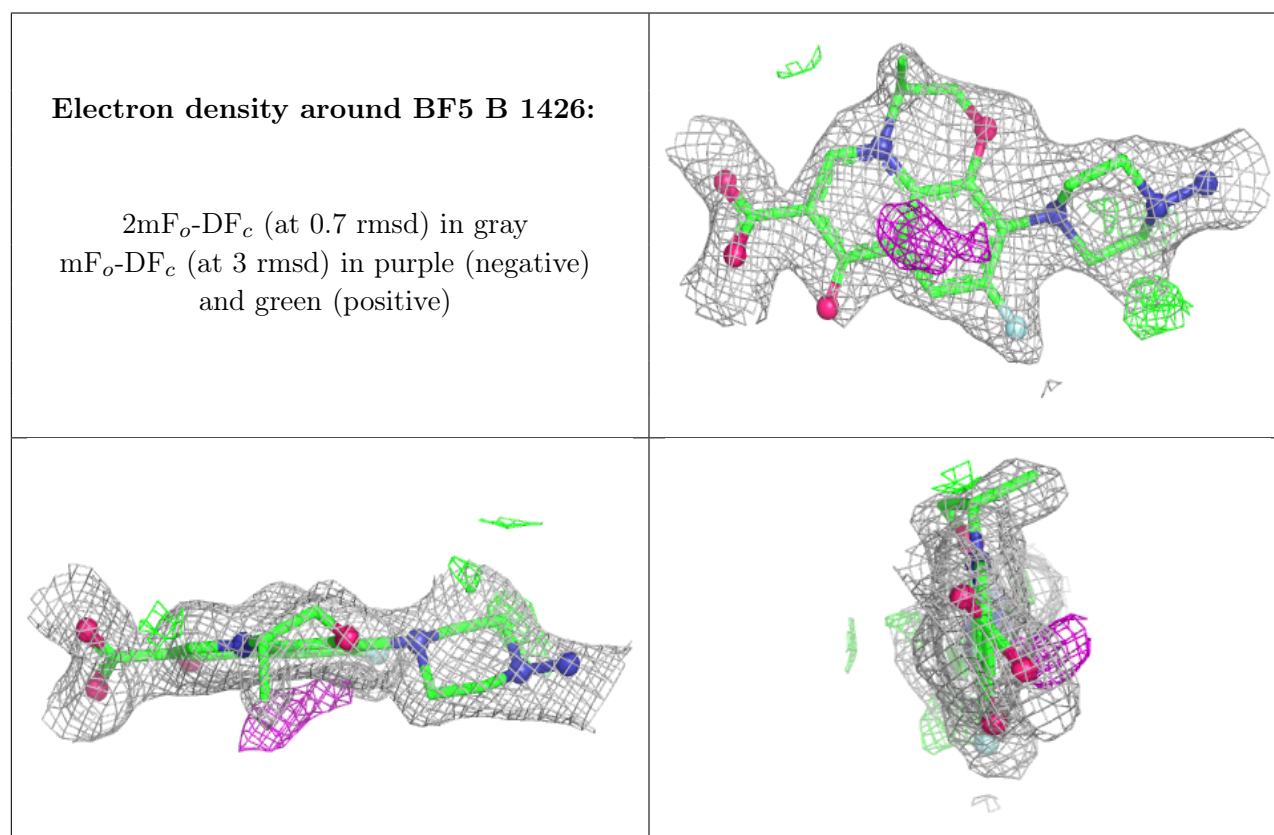
There are no oligosaccharides in this entry.

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.

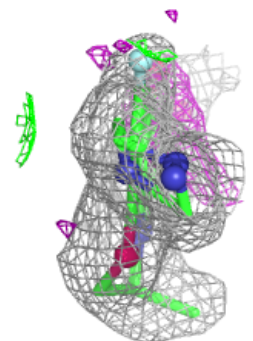
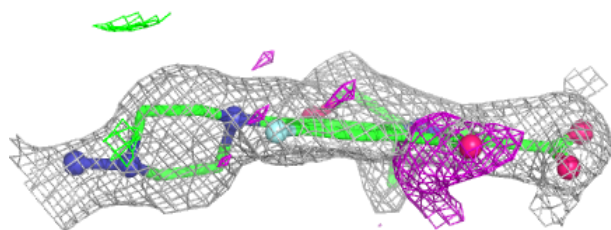
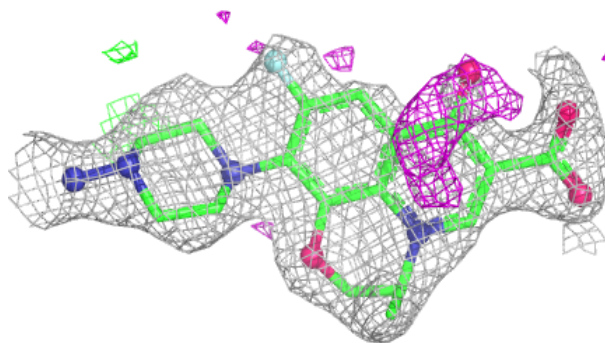
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	GOL	B	1428	6/6	0.77	0.18	47,49,50,52	0
2	BF5	B	1426	26/26	0.79	0.13	40,51,52,52	0
2	BF5	A	1426	26/26	0.84	0.12	35,40,41,43	0
3	PLP	B	1427	15/16	0.98	0.06	20,31,34,34	0
3	PLP	A	1427	15/16	0.98	0.06	18,26,29,29	0
4	IOD	A	1428	1/1	1.00	0.03	20,20,20,20	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 BF5 A 1426:

$2mF_o-DF_c$ (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.