



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:31 PM UTC

PDB ID : 8BDT / pdb_00008bdt
Title : Ternary complex between VCB, BRD4-BD2 and PROTAC 51
Authors : Sorrell, F.J.; Mueller, J.E.; Lehmann, M.; Wegener, A.A.
Deposited on : 2022-10-20
Resolution : 2.70 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

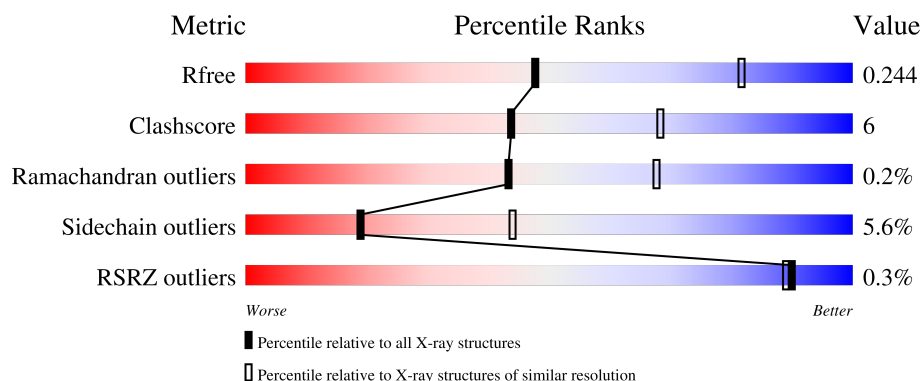
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	129	69% 13% • 16%
1	E	129	67% 18% 16%
2	B	104	78% 18% •
2	F	104	80% 17% •
3	C	97	87% 10% ••

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Mol	Chain	Length	Quality of chain
3	G	97	% 70%16%11%
4	D	162	% 77%12%8%
4	H	162	74%17%8%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7335 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 Bromodomain-containing protein 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	108	Total	C	N	O	S	0	0	0
			841	541	137	152	11			
1	E	109	Total	C	N	O	S	0	0	0
			851	544	139	157	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	332	GLY	-	expression tag	UNP O60885
E	332	GLY	-	expression tag	UNP O60885

- Molecule 2 is a protein called Elongin-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	104	Total	C	N	O	S	0	0	0
			814	515	137	158	4			
2	F	104	Total	C	N	O	S	0	0	0
			753	478	121	149	5			

- Molecule 3 is a protein called Elongin-C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	97	Total	C	N	O	S	0	0	0
			762	489	120	146	7			
3	G	86	Total	C	N	O	S	0	0	0
			665	429	104	126	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	16	MET	-	initiating methionine	UNP Q15369
G	16	MET	-	initiating methionine	UNP Q15369

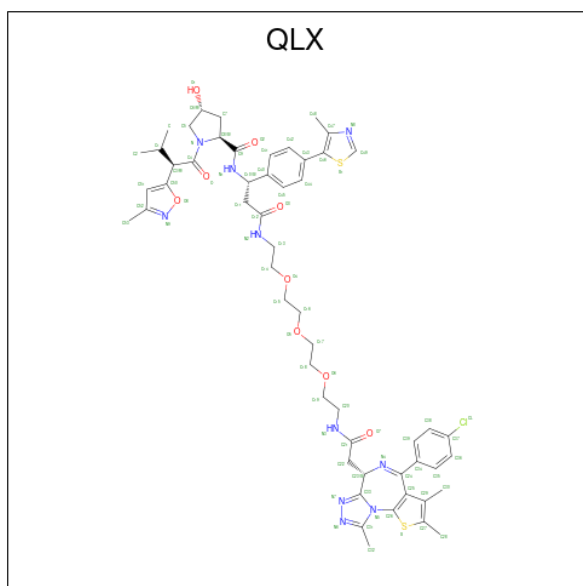
- Molecule 4 is a protein called von Hippel-Lindau disease tumor suppressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	149	Total	C	N	O	S	0	1	0
			1222	775	225	220	2			
4	H	149	Total	C	N	O	S	0	0	0
			1202	764	218	218	2			

There are 4 discrepancies between the modelled and reference sequences:

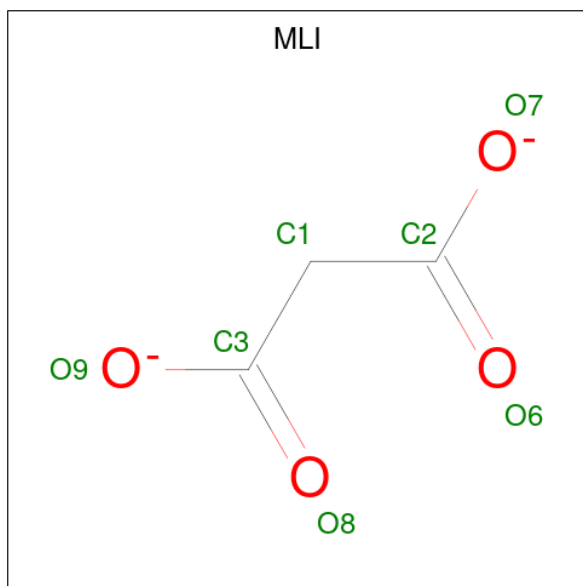
Chain	Residue	Modelled	Actual	Comment	Reference
D	52	GLY	-	expression tag	UNP P40337
D	53	SER	-	expression tag	UNP P40337
H	52	GLY	-	expression tag	UNP P40337
H	53	SER	-	expression tag	UNP P40337

- Molecule 5 is (2 {S},4 {R})- {N}-[(1 {S})-3-[2-[2-[2-[2-[2-[(9 {S})-7-(4-chlorophenyl)-4,5,1 3-trimethyl-3-thia-1,8,11,12-tetrazatricyclo[8.3.0.0[^]{2,6}}]trideca-2(6),4,7,10,12-pentaen-9-yl]ethanoylamino]ethoxy]ethoxy]ethoxy]ethylamino]-1-[4-(4-methyl-1,3-thiazol-5-yl)phenyl]-3-oxidanylidene-propyl]-1-[(2 {R})-3-methyl-2-(3-methyl-1,2-oxazol-5-yl)butanoyl]-4-oxi danyl-pyrrolidine-2-carboxamide (CCD ID: QLX) (formula: C₅₄H₆₅ClN₁₀O₉S₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	D	1	Total	C	Cl	N	O	S	0	0
			76	54	1	10	9	2		
5	H	1	Total	C	Cl	N	O	S	0	0
			76	54	1	10	9	2		

- Molecule 6 is MALONATE ION (CCD ID: MLI) (formula: $\text{C}_3\text{H}_2\text{O}_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	G	1	Total	C	O	0	0
			7	3	4		

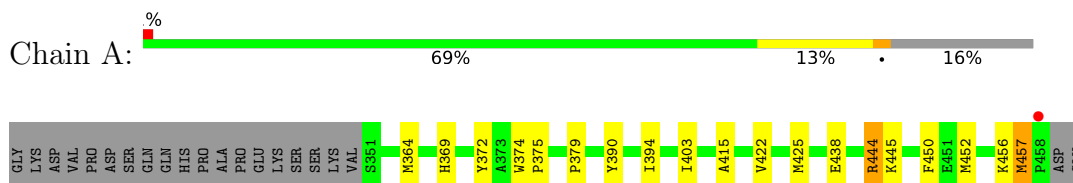
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	7	Total	O	0	0
			7	7		
7	B	8	Total	O	0	0
			8	8		
7	C	15	Total	O	0	0
			15	15		
7	D	9	Total	O	0	0
			9	9		
7	E	9	Total	O	0	0
			9	9		
7	F	3	Total	O	0	0
			3	3		
7	G	10	Total	O	0	0
			10	10		
7	H	5	Total	O	0	0
			5	5		

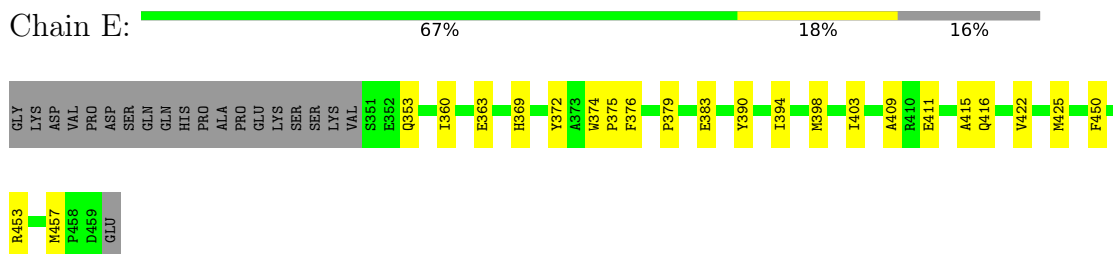
3 Residue-property plots [i](#)

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.

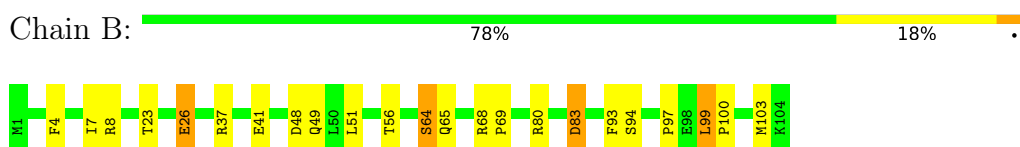
- Molecule 1: Bromodomain-containing protein 4



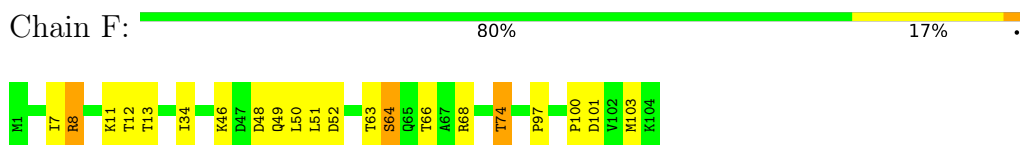
- Molecule 1: Bromodomain-containing protein 4



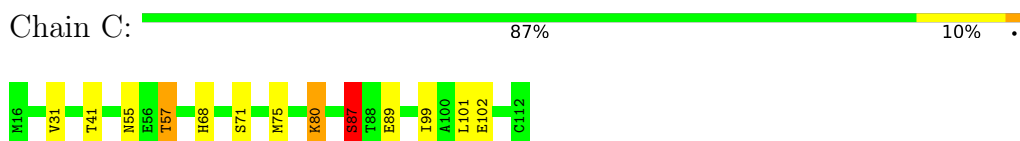
- Molecule 2: Elongin-B



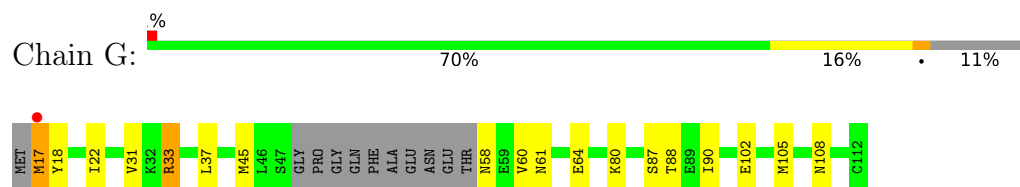
- Molecule 2: Elongin-B



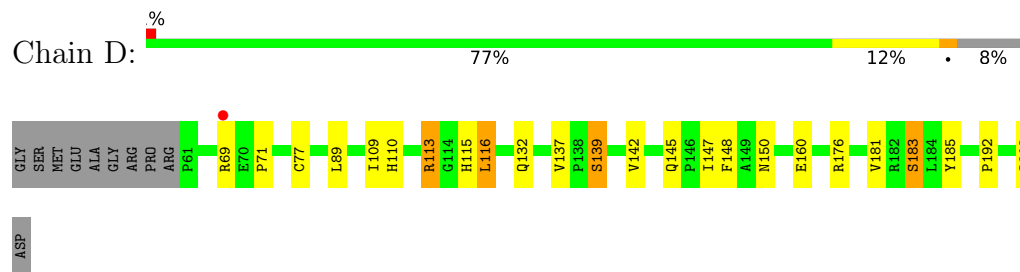
- Molecule 3: Elongin-C



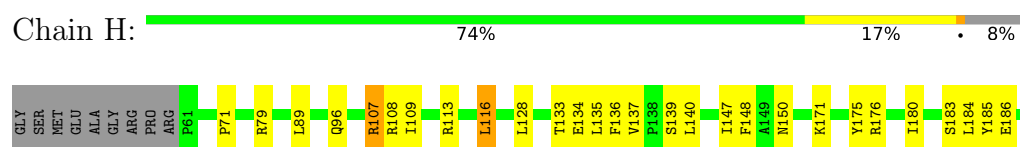
- Molecule 3: Elongin-C



- Molecule 4: von Hippel-Lindau disease tumor suppressor



- Molecule 4: von Hippel-Lindau disease tumor suppressor



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	82.48Å 82.48Å 169.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	65.80 – 2.70 65.80 – 2.70	Depositor EDS
% Data completeness (in resolution range)	82.5 (65.80-2.70) 82.6 (65.80-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.188 , 0.243 0.191 , 0.244	Depositor DCC
R_{free} test set	1513 reflections (4.30%)	wwPDB-VP
Wilson B-factor (Å ²)	81.1	Xtriage
Anisotropy	0.007	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l 0.035 for h,-h-k,-l 0.019 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7335	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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: MLI, QLX

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	1.15	1/864 (0.1%)	1.28	2/1169 (0.2%)
1	E	1.11	0/874	1.29	1/1185 (0.1%)
2	B	1.15	4/830 (0.5%)	1.27	2/1122 (0.2%)
2	F	1.11	2/769 (0.3%)	1.26	1/1049 (0.1%)
3	C	1.18	2/779 (0.3%)	1.29	1/1053 (0.1%)
3	G	1.09	0/679	1.19	1/921 (0.1%)
4	D	1.06	2/1254 (0.2%)	1.24	0/1712
4	H	0.99	1/1233 (0.1%)	1.24	1/1685 (0.1%)
All	All	1.10	12/7282 (0.2%)	1.26	9/9896 (0.1%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	57	THR	C-O	9.32	1.34	1.23
4	D	110	HIS	CG-CD2	6.39	1.42	1.35
2	B	49	GLN	C-O	-6.19	1.16	1.24
2	B	64	SER	C-O	6.15	1.31	1.24
2	F	49	GLN	C-O	-6.11	1.16	1.24
1	A	364	MET	C-O	6.03	1.31	1.24
2	B	93	PHE	C-O	5.84	1.31	1.23
4	H	175	TYR	C-O	5.57	1.30	1.24
2	F	64	SER	C-O	5.33	1.30	1.24
3	C	87	SER	C-O	5.30	1.30	1.24
2	B	8	ARG	C-O	-5.27	1.18	1.24
4	D	192	PRO	C-O	-5.16	1.17	1.23

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	379	PRO	CB-CA-C	-7.12	102.28	111.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	379	PRO	CB-CA-C	-6.51	103.05	111.39
1	A	444	ARG	CG-CD-NE	6.22	125.68	112.00
2	B	83	ASP	CA-CB-CG	5.53	118.13	112.60
2	F	12	THR	CB-CA-C	5.42	118.58	109.48
2	B	48	ASP	CA-CB-CG	5.19	117.79	112.60
4	H	171	LYS	CB-CA-C	5.05	117.50	109.52
3	C	41	THR	CA-CB-OG1	-5.03	102.06	109.60
3	G	90	ILE	N-CA-CB	-5.01	107.06	112.37

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	841	0	785	13	0
1	E	851	0	776	14	0
2	B	814	0	806	11	0
2	F	753	0	689	11	0
3	C	762	0	747	6	0
3	G	665	0	640	10	0
4	D	1222	0	1210	19	0
4	H	1202	0	1187	19	0
5	D	76	0	0	3	0
5	H	76	0	0	0	0
6	G	7	0	2	0	0
7	A	7	0	0	0	0
7	B	8	0	0	0	0
7	C	15	0	0	0	0
7	D	9	0	0	0	0
7	E	9	0	0	0	0
7	F	3	0	0	1	0
7	G	10	0	0	0	0
7	H	5	0	0	0	0
All	All	7335	0	6842	86	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 (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:438:GLU:OE2	4:D:69[A]:ARG:NH2	1.92	1.03
1:A:369:HIS:HD1	1:A:372:TYR:HH	1.08	0.95
2:F:63:THR:H	2:F:66:THR:HG22	1.28	0.94
3:G:108:ASN:HD22	4:H:184:LEU:HD21	1.35	0.90
1:E:369:HIS:HD1	1:E:372:TYR:HH	0.86	0.85
2:B:37:ARG:NH1	2:B:41:GLU:OE1	2.14	0.81
4:H:107:ARG:HD3	4:H:109:ILE:HD11	1.64	0.80
2:F:63:THR:H	2:F:66:THR:CG2	1.96	0.78
3:G:33:ARG:O	3:G:37:LEU:HD23	1.88	0.73
2:B:99:LEU:HD23	2:B:100:PRO:HD2	1.72	0.72
4:D:113:ARG:HD3	4:D:113:ARG:C	2.15	0.71
2:B:80:ARG:HD2	2:B:83:ASP:HA	1.76	0.67
3:G:108:ASN:HD22	4:H:184:LEU:CD2	2.07	0.66
2:F:100:PRO:HD2	2:F:103:MET:HE2	1.77	0.65
1:A:374:TRP:CE3	4:D:71:PRO:HG3	2.34	0.61
1:E:374:TRP:CE3	4:H:71:PRO:HG3	2.36	0.61
2:F:63:THR:N	2:F:66:THR:HG22	2.10	0.61
4:H:89:LEU:HD12	4:H:116:LEU:HG	1.84	0.60
4:D:89:LEU:HD12	4:D:116:LEU:HG	1.85	0.59
4:H:107:ARG:HD3	4:H:109:ILE:CD1	2.33	0.58
4:H:133:THR:OG1	4:H:134:GLU:N	2.20	0.58
4:D:150:ASN:OD1	4:H:79:ARG:NH2	2.39	0.56
1:E:390:TYR:CZ	1:E:394:ILE:HG13	2.40	0.55
2:F:101:ASP:HB3	7:F:201:HOH:O	2.06	0.55
2:F:97:PRO:HD2	3:G:102:GLU:OE1	2.07	0.54
1:E:398:MET:HE2	1:E:403:ILE:HG12	1.87	0.54
4:D:113:ARG:C	4:D:113:ARG:CD	2.80	0.54
3:C:80:LYS:HD2	3:C:80:LYS:C	2.33	0.53
3:C:71:SER:O	3:C:75:MET:HG3	2.09	0.52
1:A:369:HIS:ND1	1:A:372:TYR:OH	2.12	0.52
2:B:97:PRO:HD2	3:C:102:GLU:OE1	2.09	0.52
1:E:409:ALA:HB3	1:E:411:GLU:HG3	1.92	0.52
3:G:22:ILE:HB	3:G:61:ASN:HD22	1.74	0.51
4:H:176:ARG:HA	4:H:185:TYR:CE1	2.46	0.51
2:F:8:ARG:HG3	2:F:74:THR:HG22	1.93	0.51
4:D:142:VAL:O	4:D:145:GLN:HG3	2.11	0.51
4:D:148:PHE:CE2	4:H:150:ASN:CB	2.95	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:64:SER:O	2:B:68:ARG:HG3	2.11	0.50
2:B:94:SER:O	3:C:68:HIS:HB3	2.10	0.50
2:F:34:ILE:HD11	3:G:18:TYR:CZ	2.46	0.50
2:B:99:LEU:CD2	2:B:100:PRO:HD2	2.42	0.49
4:D:148:PHE:CD1	4:H:148:PHE:CD1	3.00	0.48
2:B:23:THR:HG22	2:B:26:GLU:CG	2.44	0.48
2:F:64:SER:O	2:F:68:ARG:HG3	2.14	0.48
4:D:109:ILE:HD12	5:D:301:QLX:C44	2.43	0.48
5:D:301:QLX:C51	5:D:301:QLX:C2	2.88	0.47
1:E:422:VAL:HG11	1:E:450:PHE:CD1	2.49	0.47
1:A:422:VAL:HG11	1:A:450:PHE:CD1	2.49	0.47
1:A:390:TYR:CZ	1:A:394:ILE:HG13	2.49	0.47
1:A:452:MET:HE3	1:A:456:LYS:CE	2.45	0.47
4:D:176:ARG:HA	4:D:185:TYR:CE1	2.50	0.47
1:E:353:GLN:HG2	1:E:415:ALA:HB2	1.96	0.46
4:D:150:ASN:CB	4:H:148:PHE:CE2	2.98	0.46
1:E:369:HIS:HA	1:E:372:TYR:CZ	2.51	0.46
1:E:376:PHE:HB3	1:E:425:MET:SD	2.56	0.45
2:F:50:LEU:HD22	2:F:51:LEU:H	1.82	0.45
4:D:77:CYS:SG	3:G:88:THR:HG21	2.56	0.45
4:D:139:SER:O	4:D:147:ILE:CD1	2.65	0.45
1:A:452:MET:HE3	1:A:456:LYS:NZ	2.33	0.44
4:D:181:VAL:HG12	4:D:183:SER:H	1.83	0.44
4:H:113:ARG:HB2	4:H:140:LEU:HD22	2.00	0.44
1:A:415:ALA:CB	1:A:457:MET:HE1	2.48	0.44
4:D:145:GLN:NE2	4:H:135:LEU:O	2.51	0.44
2:F:8:ARG:CD	2:F:13:THR:HG23	2.47	0.44
3:G:17:MET:HE1	3:G:58:ASN:HD21	1.83	0.43
3:C:87:SER:HA	4:D:132:GLN:HE21	1.84	0.43
1:E:374:TRP:N	1:E:375:PRO:CD	2.81	0.43
4:H:116:LEU:HD13	4:H:137:VAL:CG2	2.49	0.43
1:E:398:MET:HE3	1:E:403:ILE:HG13	2.00	0.43
1:A:452:MET:HE3	1:A:456:LYS:HE2	2.01	0.43
1:E:360:ILE:HG12	1:E:453:ARG:HB3	2.00	0.43
1:E:353:GLN:CG	1:E:415:ALA:HB2	2.49	0.42
3:G:45:MET:CE	3:G:60:VAL:HG11	2.49	0.42
2:B:23:THR:HA	2:B:56:THR:HA	2.02	0.42
4:D:116:LEU:HD13	4:D:137:VAL:CG2	2.49	0.41
4:H:116:LEU:HD13	4:H:137:VAL:HG22	2.01	0.41
2:B:4:PHE:CD2	2:B:69:PRO:HD3	2.55	0.41
4:D:115:HIS:ND1	5:D:301:QLX:O1	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:383:GLU:OE2	4:H:108:ARG:N	2.36	0.41
2:B:103:MET:HE2	2:B:103:MET:HB3	1.92	0.41
3:C:101:LEU:HD23	3:C:101:LEU:HA	1.83	0.41
1:A:369:HIS:HA	1:A:372:TYR:CZ	2.56	0.41
3:G:105:MET:CE	4:H:180:ILE:HA	2.51	0.41
1:A:403:ILE:HD12	1:A:425:MET:HE2	2.02	0.41
1:A:374:TRP:N	1:A:375:PRO:CD	2.84	0.40
4:H:136:PHE:HZ	4:H:147:ILE:HG22	1.86	0.40

There are no symmetry-related clashes.

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	106/129 (82%)	105 (99%)	1 (1%)	0	100	100
1	E	107/129 (83%)	104 (97%)	3 (3%)	0	100	100
2	B	102/104 (98%)	99 (97%)	3 (3%)	0	100	100
2	F	102/104 (98%)	99 (97%)	3 (3%)	0	100	100
3	C	95/97 (98%)	87 (92%)	6 (6%)	2 (2%)	5	15
3	G	82/97 (84%)	80 (98%)	2 (2%)	0	100	100
4	D	148/162 (91%)	143 (97%)	5 (3%)	0	100	100
4	H	147/162 (91%)	141 (96%)	6 (4%)	0	100	100
All	All	889/984 (90%)	858 (96%)	29 (3%)	2 (0%)	43	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	89	GLU
3	C	55	ASN

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	85/113 (75%)	82 (96%)	3 (4%)	32	61
1	E	85/113 (75%)	82 (96%)	3 (4%)	32	61
2	B	89/92 (97%)	84 (94%)	5 (6%)	19	44
2	F	75/92 (82%)	68 (91%)	7 (9%)	8	21
3	C	85/86 (99%)	80 (94%)	5 (6%)	18	42
3	G	73/86 (85%)	67 (92%)	6 (8%)	10	27
4	D	138/148 (93%)	132 (96%)	6 (4%)	26	54
4	H	135/148 (91%)	127 (94%)	8 (6%)	18	42
All	All	765/878 (87%)	722 (94%)	43 (6%)	19	44

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	444	ARG
1	A	445	LYS
1	A	457	MET
2	B	7	ILE
2	B	26	GLU
2	B	51	LEU
2	B	65	GLN
2	B	99	LEU
3	C	31	VAL
3	C	57	THR
3	C	80	LYS
3	C	87	SER
3	C	99	ILE
4	D	113	ARG
4	D	116	LEU
4	D	139	SER
4	D	160	GLU
4	D	183	SER
4	D	203	GLN

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Mol	Chain	Res	Type
1	E	363	GLU
1	E	416	GLN
1	E	457	MET
2	F	7	ILE
2	F	8	ARG
2	F	11	LYS
2	F	46	LYS
2	F	48	ASP
2	F	52	ASP
2	F	74	THR
3	G	17	MET
3	G	31	VAL
3	G	33	ARG
3	G	64	GLU
3	G	80	LYS
3	G	87	SER
4	H	96	GLN
4	H	107	ARG
4	H	116	LEU
4	H	128	LEU
4	H	139	SER
4	H	183	SER
4	H	186	GLU
4	H	209	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	433	ASN
3	C	35	HIS
3	C	61	ASN
4	D	73	GLN
4	D	96	GLN
4	D	110	HIS
4	D	145	GLN
4	D	203	GLN
1	E	396	HIS
1	E	433	ASN
3	G	35	HIS
3	G	58	ASN
3	G	61	ASN
3	G	108	ASN

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Mol	Chain	Res	Type
4	H	96	GLN
4	H	132	GLN
4	H	191	HIS

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](#)

3 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	$\# Z > 2$
5	QLX	H	301	-	77,83,83	0.78	2 (2%)	98,117,117	0.65	1 (1%)
6	MLI	G	201	-	6,6,6	1.70	2 (33%)	7,7,7	1.09	0
5	QLX	D	301	-	77,83,83	0.66	2 (2%)	98,117,117	0.74	3 (3%)

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
5	QLX	H	301	-	-	17/60/88/88	0/7/8/8
6	MLI	G	201	-	-	2/4/4/4	-
5	QLX	D	301	-	-	13/60/88/88	0/7/8/8

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	301	QLX	C52-N9	4.37	1.38	1.30
5	H	301	QLX	C26-S	-3.86	1.66	1.74
5	D	301	QLX	C52-N9	2.70	1.35	1.30
6	G	201	MLI	C1-C3	2.68	1.55	1.51
6	G	201	MLI	C1-C2	2.09	1.54	1.51
5	D	301	QLX	C46-S1	2.07	1.78	1.73

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	301	QLX	C53-C52-N9	3.03	122.45	119.97
5	D	301	QLX	C11-C10-N1	-2.21	104.94	108.94
5	D	301	QLX	C10-C11-C12	-2.18	107.58	114.49
5	D	301	QLX	C-C1-C3	2.00	116.42	111.68

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	301	QLX	C2-C1-C3-C4
5	D	301	QLX	C2-C1-C3-C50
5	D	301	QLX	C-C1-C3-C4
5	D	301	QLX	C-C1-C3-C50
5	H	301	QLX	C11-C12-N2-C13
5	H	301	QLX	C2-C1-C3-C4
5	H	301	QLX	C2-C1-C3-C50
5	H	301	QLX	C-C1-C3-C4
5	H	301	QLX	C-C1-C3-C50
5	H	301	QLX	C4-C3-C50-O8
5	H	301	QLX	O3-C12-N2-C13
5	D	301	QLX	O6-C19-C20-N3
5	D	301	QLX	O7-C21-N3-C20
5	D	301	QLX	C22-C21-N3-C20
5	D	301	QLX	N2-C13-C14-O4

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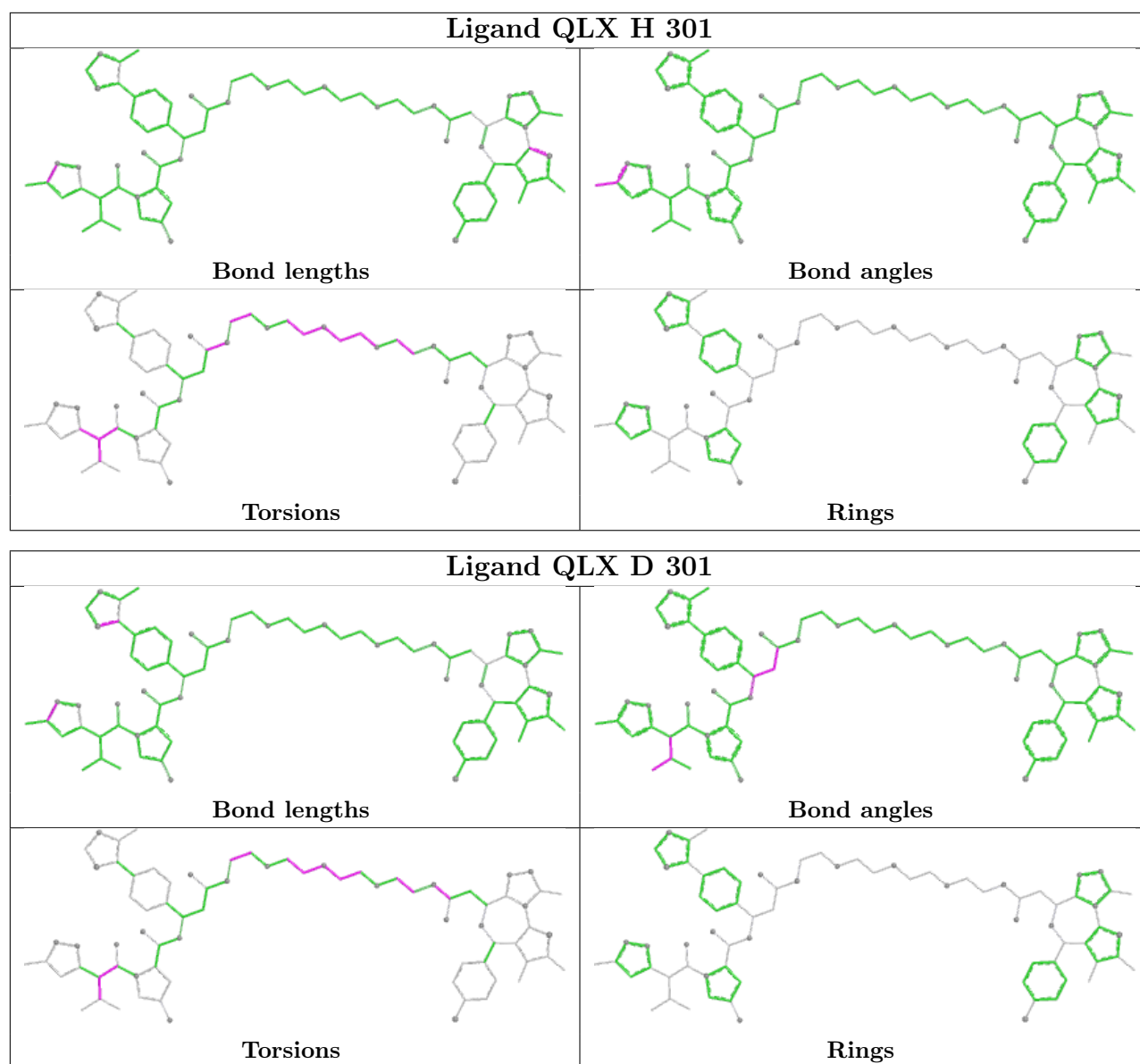
Mol	Chain	Res	Type	Atoms
5	H	301	QLX	O4-C15-C16-O5
5	H	301	QLX	O6-C19-C20-N3
5	H	301	QLX	N2-C13-C14-O4
5	H	301	QLX	O5-C17-C18-O6
5	D	301	QLX	O4-C15-C16-O5
5	D	301	QLX	C18-C17-O5-C16
5	H	301	QLX	C17-C18-O6-C19
5	D	301	QLX	C15-C16-O5-C17
5	D	301	QLX	C1-C3-C4-O
5	H	301	QLX	C1-C3-C4-O
5	H	301	QLX	C1-C3-C50-O8
5	H	301	QLX	C18-C17-O5-C16
5	D	301	QLX	O5-C17-C18-O6
5	H	301	QLX	C15-C16-O5-C17
6	G	201	MLI	C3-C1-C2-O7
6	G	201	MLI	C3-C1-C2-O6
5	H	301	QLX	C1-C3-C4-N

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	301	QLX	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 [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	108/129 (83%)	-0.14	1 (0%) 81 80	58, 74, 119, 137	0
1	E	109/129 (84%)	-0.27	0 100 100	54, 71, 113, 149	0
2	B	104/104 (100%)	-0.28	0 100 100	59, 89, 109, 127	0
2	F	104/104 (100%)	0.08	0 100 100	77, 111, 139, 180	0
3	C	97/97 (100%)	-0.25	0 100 100	52, 73, 139, 149	0
3	G	86/97 (88%)	-0.15	1 (1%) 76 75	69, 92, 112, 126	0
4	D	149/162 (91%)	-0.34	1 (0%) 84 83	41, 70, 106, 129	1 (0%)
4	H	149/162 (91%)	-0.19	0 100 100	57, 83, 126, 154	0
All	All	906/984 (92%)	-0.20	3 (0%) 90 89	41, 82, 127, 180	1 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	69[A]	ARG	3.6
1	A	458	PRO	2.6
3	G	17	MET	2.1

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

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

6.3 Carbohydrates [i](#)

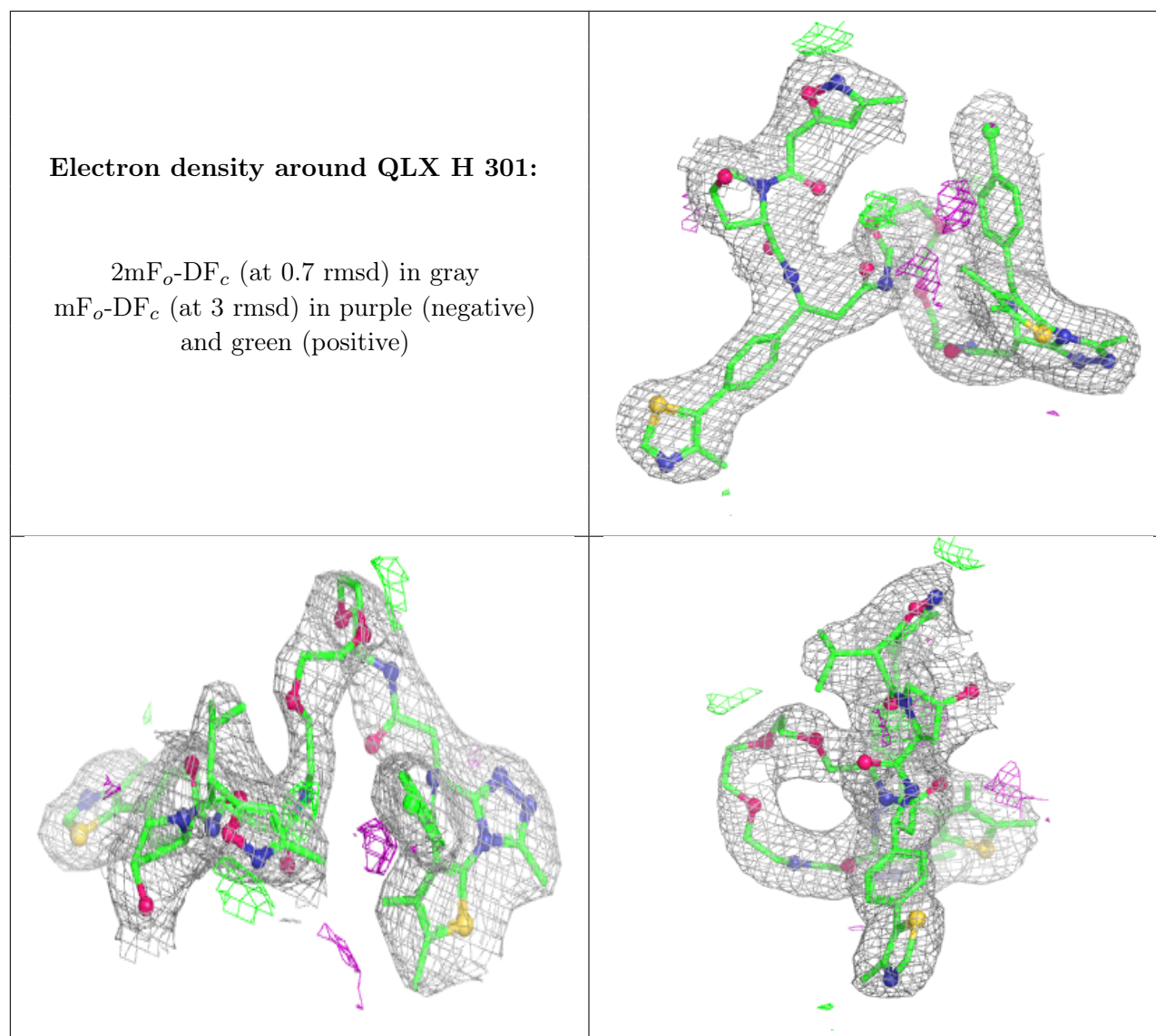
There are no oligosaccharides in this entry.

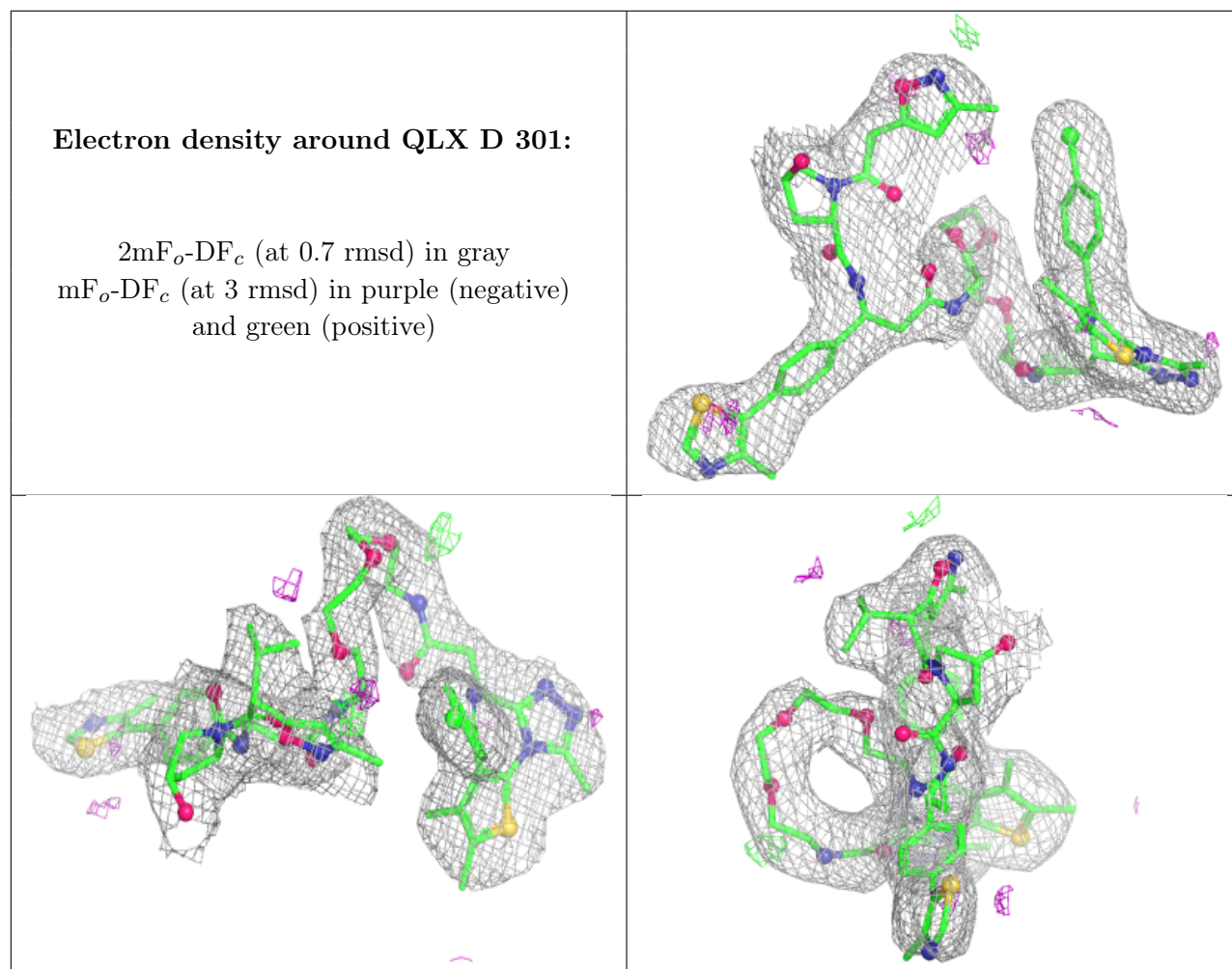
6.4 Ligands ⓘ

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	MLI	G	201	7/7	0.86	0.08	91,94,99,101	0
5	QLX	H	301	76/76	0.95	0.09	51,71,103,113	0
5	QLX	D	301	76/76	0.95	0.09	56,70,101,108	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 ⓘ

There are no such residues in this entry.