



wwPDB X-ray Structure Validation Summary Report

Oct 7, 2023 – 04:19 PM EDT

PDB ID : 4Q0N
Title : Crystal Structure of the fifth bromodomain of Human Poly-bromodomain containing protein 1 (PB1) in complex with a hydroxyphenyl-propenone ligand
Authors : Filippakopoulos, P.; Picaud, S.; Felletar, I.; Martin, S.; Monteiro, O.; Fedorov, O.; Chaikuad, A.; Yue, W.; von Delft, F.; Arrowsmith, C.H.; Edwards, A.M.; Bountra, C.; Knapp, S.; Structural Genomics Consortium (SGC)
Deposited on : 2014-04-02
Resolution : 1.78 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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](#) ①) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

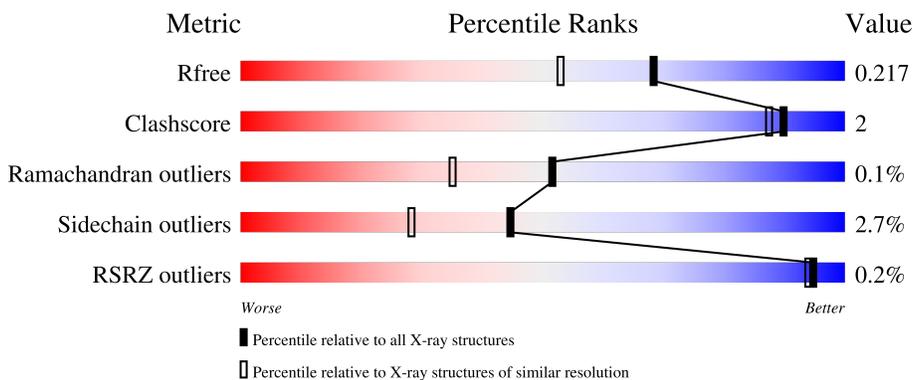
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.78 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}	130704	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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	145	
1	B	145	
1	C	145	
1	D	145	
1	E	145	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	145	 74% • 23%
1	G	145	 72% 5% • 23%
1	H	145	 73% • 23%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8215 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 Protein polybromo-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	113	948	603	162	173	10	0	2	0
1	B	111	915	579	155	171	10	0	0	0
1	C	111	919	585	153	170	11	0	2	0
1	D	112	931	593	155	172	11	0	2	0
1	E	113	943	601	157	175	10	0	3	0
1	F	112	927	588	157	172	10	0	0	0
1	G	112	923	587	154	172	10	0	1	0
1	H	111	928	589	157	172	10	0	1	0

There are 184 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	590	MET	-	expression tag	UNP Q86U86
A	591	HIS	-	expression tag	UNP Q86U86
A	592	HIS	-	expression tag	UNP Q86U86
A	593	HIS	-	expression tag	UNP Q86U86
A	594	HIS	-	expression tag	UNP Q86U86
A	595	HIS	-	expression tag	UNP Q86U86
A	596	HIS	-	expression tag	UNP Q86U86
A	597	SER	-	expression tag	UNP Q86U86
A	598	SER	-	expression tag	UNP Q86U86
A	599	GLY	-	expression tag	UNP Q86U86
A	600	VAL	-	expression tag	UNP Q86U86
A	601	ASP	-	expression tag	UNP Q86U86
A	602	LEU	-	expression tag	UNP Q86U86

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	603	GLY	-	expression tag	UNP Q86U86
A	604	THR	-	expression tag	UNP Q86U86
A	605	GLU	-	expression tag	UNP Q86U86
A	606	ASN	-	expression tag	UNP Q86U86
A	607	LEU	-	expression tag	UNP Q86U86
A	608	TYR	-	expression tag	UNP Q86U86
A	609	PHE	-	expression tag	UNP Q86U86
A	610	GLN	-	expression tag	UNP Q86U86
A	611	SER	-	expression tag	UNP Q86U86
A	612	MET	-	expression tag	UNP Q86U86
B	590	MET	-	expression tag	UNP Q86U86
B	591	HIS	-	expression tag	UNP Q86U86
B	592	HIS	-	expression tag	UNP Q86U86
B	593	HIS	-	expression tag	UNP Q86U86
B	594	HIS	-	expression tag	UNP Q86U86
B	595	HIS	-	expression tag	UNP Q86U86
B	596	HIS	-	expression tag	UNP Q86U86
B	597	SER	-	expression tag	UNP Q86U86
B	598	SER	-	expression tag	UNP Q86U86
B	599	GLY	-	expression tag	UNP Q86U86
B	600	VAL	-	expression tag	UNP Q86U86
B	601	ASP	-	expression tag	UNP Q86U86
B	602	LEU	-	expression tag	UNP Q86U86
B	603	GLY	-	expression tag	UNP Q86U86
B	604	THR	-	expression tag	UNP Q86U86
B	605	GLU	-	expression tag	UNP Q86U86
B	606	ASN	-	expression tag	UNP Q86U86
B	607	LEU	-	expression tag	UNP Q86U86
B	608	TYR	-	expression tag	UNP Q86U86
B	609	PHE	-	expression tag	UNP Q86U86
B	610	GLN	-	expression tag	UNP Q86U86
B	611	SER	-	expression tag	UNP Q86U86
B	612	MET	-	expression tag	UNP Q86U86
C	590	MET	-	expression tag	UNP Q86U86
C	591	HIS	-	expression tag	UNP Q86U86
C	592	HIS	-	expression tag	UNP Q86U86
C	593	HIS	-	expression tag	UNP Q86U86
C	594	HIS	-	expression tag	UNP Q86U86
C	595	HIS	-	expression tag	UNP Q86U86
C	596	HIS	-	expression tag	UNP Q86U86
C	597	SER	-	expression tag	UNP Q86U86
C	598	SER	-	expression tag	UNP Q86U86

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	599	GLY	-	expression tag	UNP Q86U86
C	600	VAL	-	expression tag	UNP Q86U86
C	601	ASP	-	expression tag	UNP Q86U86
C	602	LEU	-	expression tag	UNP Q86U86
C	603	GLY	-	expression tag	UNP Q86U86
C	604	THR	-	expression tag	UNP Q86U86
C	605	GLU	-	expression tag	UNP Q86U86
C	606	ASN	-	expression tag	UNP Q86U86
C	607	LEU	-	expression tag	UNP Q86U86
C	608	TYR	-	expression tag	UNP Q86U86
C	609	PHE	-	expression tag	UNP Q86U86
C	610	GLN	-	expression tag	UNP Q86U86
C	611	SER	-	expression tag	UNP Q86U86
C	612	MET	-	expression tag	UNP Q86U86
D	590	MET	-	expression tag	UNP Q86U86
D	591	HIS	-	expression tag	UNP Q86U86
D	592	HIS	-	expression tag	UNP Q86U86
D	593	HIS	-	expression tag	UNP Q86U86
D	594	HIS	-	expression tag	UNP Q86U86
D	595	HIS	-	expression tag	UNP Q86U86
D	596	HIS	-	expression tag	UNP Q86U86
D	597	SER	-	expression tag	UNP Q86U86
D	598	SER	-	expression tag	UNP Q86U86
D	599	GLY	-	expression tag	UNP Q86U86
D	600	VAL	-	expression tag	UNP Q86U86
D	601	ASP	-	expression tag	UNP Q86U86
D	602	LEU	-	expression tag	UNP Q86U86
D	603	GLY	-	expression tag	UNP Q86U86
D	604	THR	-	expression tag	UNP Q86U86
D	605	GLU	-	expression tag	UNP Q86U86
D	606	ASN	-	expression tag	UNP Q86U86
D	607	LEU	-	expression tag	UNP Q86U86
D	608	TYR	-	expression tag	UNP Q86U86
D	609	PHE	-	expression tag	UNP Q86U86
D	610	GLN	-	expression tag	UNP Q86U86
D	611	SER	-	expression tag	UNP Q86U86
D	612	MET	-	expression tag	UNP Q86U86
E	590	MET	-	expression tag	UNP Q86U86
E	591	HIS	-	expression tag	UNP Q86U86
E	592	HIS	-	expression tag	UNP Q86U86
E	593	HIS	-	expression tag	UNP Q86U86
E	594	HIS	-	expression tag	UNP Q86U86

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	595	HIS	-	expression tag	UNP Q86U86
E	596	HIS	-	expression tag	UNP Q86U86
E	597	SER	-	expression tag	UNP Q86U86
E	598	SER	-	expression tag	UNP Q86U86
E	599	GLY	-	expression tag	UNP Q86U86
E	600	VAL	-	expression tag	UNP Q86U86
E	601	ASP	-	expression tag	UNP Q86U86
E	602	LEU	-	expression tag	UNP Q86U86
E	603	GLY	-	expression tag	UNP Q86U86
E	604	THR	-	expression tag	UNP Q86U86
E	605	GLU	-	expression tag	UNP Q86U86
E	606	ASN	-	expression tag	UNP Q86U86
E	607	LEU	-	expression tag	UNP Q86U86
E	608	TYR	-	expression tag	UNP Q86U86
E	609	PHE	-	expression tag	UNP Q86U86
E	610	GLN	-	expression tag	UNP Q86U86
E	611	SER	-	expression tag	UNP Q86U86
E	612	MET	-	expression tag	UNP Q86U86
F	590	MET	-	expression tag	UNP Q86U86
F	591	HIS	-	expression tag	UNP Q86U86
F	592	HIS	-	expression tag	UNP Q86U86
F	593	HIS	-	expression tag	UNP Q86U86
F	594	HIS	-	expression tag	UNP Q86U86
F	595	HIS	-	expression tag	UNP Q86U86
F	596	HIS	-	expression tag	UNP Q86U86
F	597	SER	-	expression tag	UNP Q86U86
F	598	SER	-	expression tag	UNP Q86U86
F	599	GLY	-	expression tag	UNP Q86U86
F	600	VAL	-	expression tag	UNP Q86U86
F	601	ASP	-	expression tag	UNP Q86U86
F	602	LEU	-	expression tag	UNP Q86U86
F	603	GLY	-	expression tag	UNP Q86U86
F	604	THR	-	expression tag	UNP Q86U86
F	605	GLU	-	expression tag	UNP Q86U86
F	606	ASN	-	expression tag	UNP Q86U86
F	607	LEU	-	expression tag	UNP Q86U86
F	608	TYR	-	expression tag	UNP Q86U86
F	609	PHE	-	expression tag	UNP Q86U86
F	610	GLN	-	expression tag	UNP Q86U86
F	611	SER	-	expression tag	UNP Q86U86
F	612	MET	-	expression tag	UNP Q86U86
G	590	MET	-	expression tag	UNP Q86U86

Continued on next page...

Continued from previous page...

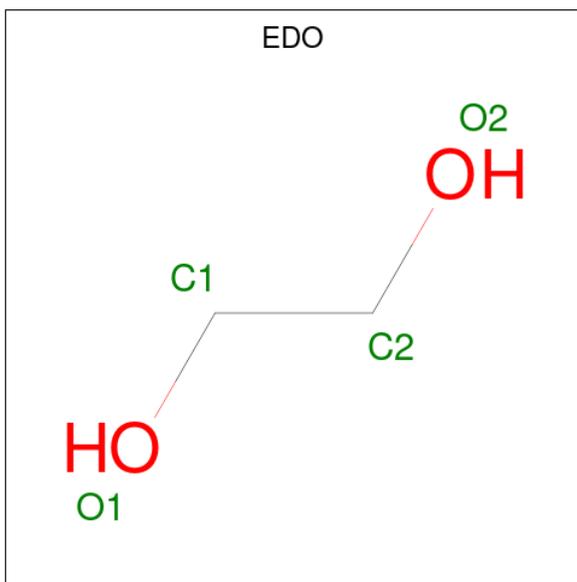
Chain	Residue	Modelled	Actual	Comment	Reference
G	591	HIS	-	expression tag	UNP Q86U86
G	592	HIS	-	expression tag	UNP Q86U86
G	593	HIS	-	expression tag	UNP Q86U86
G	594	HIS	-	expression tag	UNP Q86U86
G	595	HIS	-	expression tag	UNP Q86U86
G	596	HIS	-	expression tag	UNP Q86U86
G	597	SER	-	expression tag	UNP Q86U86
G	598	SER	-	expression tag	UNP Q86U86
G	599	GLY	-	expression tag	UNP Q86U86
G	600	VAL	-	expression tag	UNP Q86U86
G	601	ASP	-	expression tag	UNP Q86U86
G	602	LEU	-	expression tag	UNP Q86U86
G	603	GLY	-	expression tag	UNP Q86U86
G	604	THR	-	expression tag	UNP Q86U86
G	605	GLU	-	expression tag	UNP Q86U86
G	606	ASN	-	expression tag	UNP Q86U86
G	607	LEU	-	expression tag	UNP Q86U86
G	608	TYR	-	expression tag	UNP Q86U86
G	609	PHE	-	expression tag	UNP Q86U86
G	610	GLN	-	expression tag	UNP Q86U86
G	611	SER	-	expression tag	UNP Q86U86
G	612	MET	-	expression tag	UNP Q86U86
H	590	MET	-	expression tag	UNP Q86U86
H	591	HIS	-	expression tag	UNP Q86U86
H	592	HIS	-	expression tag	UNP Q86U86
H	593	HIS	-	expression tag	UNP Q86U86
H	594	HIS	-	expression tag	UNP Q86U86
H	595	HIS	-	expression tag	UNP Q86U86
H	596	HIS	-	expression tag	UNP Q86U86
H	597	SER	-	expression tag	UNP Q86U86
H	598	SER	-	expression tag	UNP Q86U86
H	599	GLY	-	expression tag	UNP Q86U86
H	600	VAL	-	expression tag	UNP Q86U86
H	601	ASP	-	expression tag	UNP Q86U86
H	602	LEU	-	expression tag	UNP Q86U86
H	603	GLY	-	expression tag	UNP Q86U86
H	604	THR	-	expression tag	UNP Q86U86
H	605	GLU	-	expression tag	UNP Q86U86
H	606	ASN	-	expression tag	UNP Q86U86
H	607	LEU	-	expression tag	UNP Q86U86
H	608	TYR	-	expression tag	UNP Q86U86
H	609	PHE	-	expression tag	UNP Q86U86

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
H	610	GLN	-	expression tag	UNP Q86U86
H	611	SER	-	expression tag	UNP Q86U86
H	612	MET	-	expression tag	UNP Q86U86

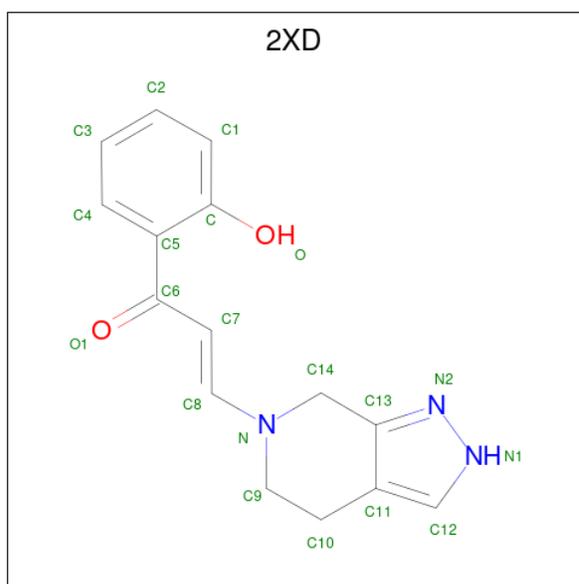
- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	E	1	Total C O 4 2 2	0	0
2	F	1	Total C O 4 2 2	0	0
2	G	1	Total C O 4 2 2	0	0
2	G	1	Total C O 4 2 2	0	0
2	H	1	Total C O 4 2 2	0	0

- Molecule 3 is (2E)-1-(2-hydroxyphenyl)-3-(2,4,5,7-tetrahydro-6H-pyrazolo[3,4-c]pyridin-6-yl

)prop-2-en-1-one (three-letter code: 2XD) (formula: C₁₅H₁₅N₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	1
			30	22	6	2		
3	B	1	Total	C	N	O	0	0
			20	15	3	2		
3	C	1	Total	C	N	O	0	1
			30	22	6	2		
3	D	1	Total	C	N	O	0	1
			30	22	6	2		
3	E	1	Total	C	N	O	0	1
			30	22	6	2		
3	F	1	Total	C	N	O	0	0
			20	15	3	2		
3	G	1	Total	C	N	O	0	0
			20	15	3	2		
3	H	1	Total	C	N	O	0	0
			20	15	3	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	64	Total	O	0	0
			64	64		
4	B	65	Total	O	0	0
			65	65		
4	C	61	Total	O	0	0
			61	61		

Continued on next page...

Continued from previous page...

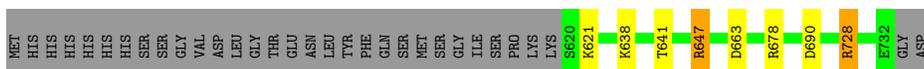
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	69	Total O 69 69	0	0
4	E	76	Total O 76 76	0	0
4	F	58	Total O 58 58	0	0
4	G	72	Total O 72 72	0	0
4	H	80	Total O 80 80	0	0

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: Protein polybromo-1

Chain A:  72% 22%



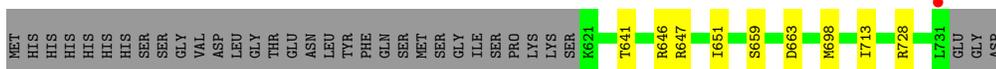
- Molecule 1: Protein polybromo-1

Chain B:  71% 23%



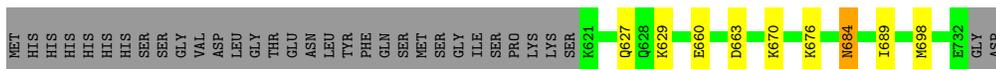
- Molecule 1: Protein polybromo-1

Chain C:  70% 23%



- Molecule 1: Protein polybromo-1

Chain D:  71% 23%



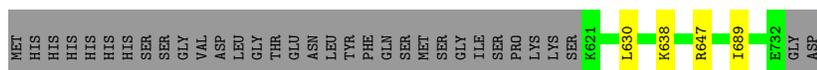
- Molecule 1: Protein polybromo-1

Chain E:  74% 22%



- Molecule 1: Protein polybromo-1

Chain F:  74% 23%



• Molecule 1: Protein polybromo-1

Chain G:  72% 5% 23%



• Molecule 1: Protein polybromo-1

Chain H:  73% 23%



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.02Å 136.72Å 114.41Å 90.00° 91.40° 90.00°	Depositor
Resolution (Å)	26.69 – 1.78 26.59 – 1.78	Depositor EDS
% Data completeness (in resolution range)	96.9 (26.69-1.78) 96.9 (26.59-1.78)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 1.78Å)	Xtrriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.176 , 0.214 0.186 , 0.217	Depositor DCC
R_{free} test set	1938 reflections (1.62%)	wwPDB-VP
Wilson B-factor (Å ²)	22.1	Xtrriage
Anisotropy	0.090	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 31.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.079 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8215	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.15 % 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 7.4557e-03.*

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: 2XD, EDO

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	0.77	0/971	0.84	2/1304 (0.2%)
1	B	0.66	0/932	0.84	2/1255 (0.2%)
1	C	0.67	0/942	0.85	2/1269 (0.2%)
1	D	0.71	0/954	0.76	0/1283
1	E	0.80	0/969	0.82	1/1303 (0.1%)
1	F	0.68	0/944	0.77	0/1270
1	G	0.76	0/944	0.81	1/1273 (0.1%)
1	H	0.69	0/948	0.81	0/1274
All	All	0.72	0/7604	0.81	8/10231 (0.1%)

There are no bond length outliers.

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	698	MET	CG-SD-CE	-11.18	82.32	100.20
1	B	626	MET	CG-SD-CE	8.31	113.49	100.20
1	G	692	MET	CG-SD-CE	-8.21	87.06	100.20
1	A	647[A]	ARG	NE-CZ-NH2	-5.79	117.40	120.30
1	A	647[B]	ARG	NE-CZ-NH2	-5.79	117.40	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	948	0	955	4	0
1	B	915	0	901	4	0
1	C	919	0	908	2	0
1	D	931	0	928	4	0
1	E	943	0	950	3	0
1	F	927	0	923	2	0
1	G	923	0	903	3	0
1	H	928	0	935	3	0
2	A	4	0	6	0	0
2	C	4	0	6	0	0
2	D	4	0	6	0	0
2	E	8	0	12	1	0
2	F	4	0	6	0	0
2	G	8	0	12	0	0
2	H	4	0	6	0	0
3	A	30	0	20	1	0
3	B	20	0	14	0	0
3	C	30	0	20	0	0
3	D	30	0	20	1	0
3	E	30	0	20	0	0
3	F	20	0	15	0	0
3	G	20	0	15	0	0
3	H	20	0	15	0	0
4	A	64	0	0	1	0
4	B	65	0	0	0	0
4	C	61	0	0	0	2
4	D	69	0	0	0	0
4	E	76	0	0	0	0
4	F	58	0	0	3	2
4	G	72	0	0	0	0
4	H	80	0	0	2	0
All	All	8215	0	7596	26	2

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 2.

The worst 5 of 26 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:F:638:LYS:NZ	4:F:942:HOH:O	1.97	0.89
1:G:680[A]:HIS:CE1	1:G:685:LYS:CB	2.70	0.73
1:B:626:MET:CE	1:B:626:MET:HA	2.32	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:630:LEU:HD23	1:F:689:ILE:HD13	1.82	0.60
1:A:647[A]:ARG:HD3	4:F:907:HOH:O	2.02	0.60

All (2) 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)
4:C:934:HOH:O	4:F:915:HOH:O[2_546]	2.04	0.16
4:C:956:HOH:O	4:F:928:HOH:O[2_546]	2.16	0.04

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	113/145 (78%)	112 (99%)	1 (1%)	0	100	100
1	B	109/145 (75%)	107 (98%)	1 (1%)	1 (1%)	17	5
1	C	111/145 (77%)	111 (100%)	0	0	100	100
1	D	112/145 (77%)	112 (100%)	0	0	100	100
1	E	114/145 (79%)	114 (100%)	0	0	100	100
1	F	110/145 (76%)	110 (100%)	0	0	100	100
1	G	111/145 (77%)	111 (100%)	0	0	100	100
1	H	110/145 (76%)	110 (100%)	0	0	100	100
All	All	890/1160 (77%)	887 (100%)	2 (0%)	1 (0%)	51	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	684	ASN

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	105/135 (78%)	99 (94%)	6 (6%)	20	7
1	B	101/135 (75%)	97 (96%)	4 (4%)	31	14
1	C	101/135 (75%)	96 (95%)	5 (5%)	24	9
1	D	103/135 (76%)	100 (97%)	3 (3%)	42	25
1	E	106/135 (78%)	104 (98%)	2 (2%)	57	43
1	F	103/135 (76%)	102 (99%)	1 (1%)	76	68
1	G	101/135 (75%)	99 (98%)	2 (2%)	55	40
1	H	105/135 (78%)	104 (99%)	1 (1%)	76	68
All	All	825/1080 (76%)	801 (97%)	24 (3%)	44	25

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	728	ARG
1	D	684	ASN
1	D	676	LYS
1	E	663	ASP
1	B	626	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	684	ASN
1	D	684	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

21 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
3	2XD	E	803[A]	-	22,22,22	2.45	8 (36%)	22,30,30	1.41	4 (18%)
3	2XD	E	803[B]	-	22,22,22	2.46	8 (36%)	22,30,30	1.27	3 (13%)
3	2XD	D	802[A]	-	22,22,22	2.18	9 (40%)	22,30,30	1.34	2 (9%)
3	2XD	D	802[B]	-	22,22,22	2.14	10 (45%)	22,30,30	1.29	2 (9%)
2	EDO	D	801	-	3,3,3	0.50	0	2,2,2	0.26	0
2	EDO	H	801	-	3,3,3	0.60	0	2,2,2	0.15	0
2	EDO	E	802	-	3,3,3	0.20	0	2,2,2	0.39	0
3	2XD	A	802[A]	-	22,22,22	2.03	7 (31%)	22,30,30	1.19	1 (4%)
2	EDO	E	801	-	3,3,3	0.76	0	2,2,2	0.20	0
3	2XD	A	802[B]	-	22,22,22	2.18	9 (40%)	22,30,30	1.08	1 (4%)
2	EDO	A	801	-	3,3,3	0.29	0	2,2,2	1.18	0
3	2XD	G	803	-	22,22,22	1.80	7 (31%)	22,30,30	1.29	3 (13%)
3	2XD	H	802	-	22,22,22	2.11	9 (40%)	22,30,30	1.15	1 (4%)
3	2XD	F	802	-	22,22,22	1.94	7 (31%)	22,30,30	1.19	1 (4%)
2	EDO	G	802	-	3,3,3	0.32	0	2,2,2	0.83	0
2	EDO	G	801	-	3,3,3	0.58	0	2,2,2	0.50	0
2	EDO	C	801	-	3,3,3	0.41	0	2,2,2	0.54	0
3	2XD	B	801	-	22,22,22	1.77	7 (31%)	22,30,30	0.90	1 (4%)
2	EDO	F	801	-	3,3,3	0.36	0	2,2,2	0.50	0
3	2XD	C	802[A]	-	22,22,22	2.02	9 (40%)	22,30,30	1.20	2 (9%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	2XD	C	802[B]	-	22,22,22	2.05	8 (36%)	22,30,30	1.08	2 (9%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	2XD	E	803[A]	-	-	0/9/18/18	0/3/3/3
3	2XD	E	803[B]	-	-	0/9/18/18	0/3/3/3
3	2XD	D	802[A]	-	-	0/9/18/18	0/3/3/3
3	2XD	D	802[B]	-	-	0/9/18/18	0/3/3/3
2	EDO	D	801	-	-	0/1/1/1	-
2	EDO	H	801	-	-	1/1/1/1	-
2	EDO	E	802	-	-	1/1/1/1	-
3	2XD	A	802[A]	-	-	0/9/18/18	0/3/3/3
2	EDO	E	801	-	-	0/1/1/1	-
3	2XD	A	802[B]	-	-	0/9/18/18	0/3/3/3
2	EDO	A	801	-	-	1/1/1/1	-
3	2XD	G	803	-	-	0/9/18/18	0/3/3/3
3	2XD	H	802	-	-	0/9/18/18	0/3/3/3
3	2XD	F	802	-	-	0/9/18/18	0/3/3/3
2	EDO	G	802	-	-	0/1/1/1	-
2	EDO	G	801	-	-	0/1/1/1	-
2	EDO	C	801	-	-	0/1/1/1	-
3	2XD	B	801	-	-	0/9/18/18	0/3/3/3
2	EDO	F	801	-	-	1/1/1/1	-
3	2XD	C	802[A]	-	-	1/9/18/18	0/3/3/3
3	2XD	C	802[B]	-	-	0/9/18/18	0/3/3/3

The worst 5 of 98 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	803[A]	2XD	O-C	5.21	1.47	1.36
3	E	803[B]	2XD	O-C	5.21	1.47	1.36
3	D	802[A]	2XD	O-C	4.98	1.46	1.36
3	D	802[B]	2XD	O-C	4.98	1.46	1.36
3	H	802	2XD	C4-C5	4.64	1.47	1.39

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	802[A]	2XD	C-C5-C6	-3.39	117.63	119.82
3	D	802[B]	2XD	C-C5-C6	-3.39	117.63	119.82
3	A	802[A]	2XD	C8-C7-C6	-3.37	116.60	119.75
3	G	803	2XD	C8-C7-C6	-3.21	116.75	119.75
3	F	802	2XD	C9-C10-C11	-2.62	106.67	111.35

There are no chirality outliers.

All (5) torsion outliers are listed below:

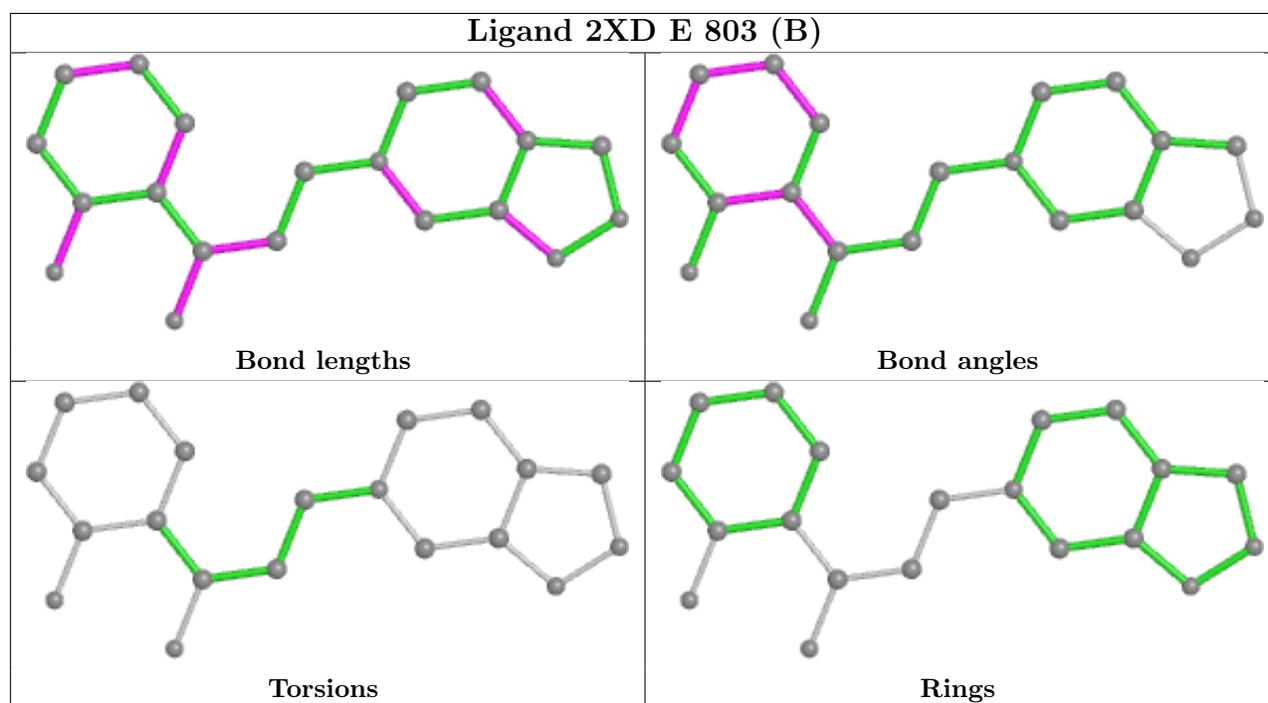
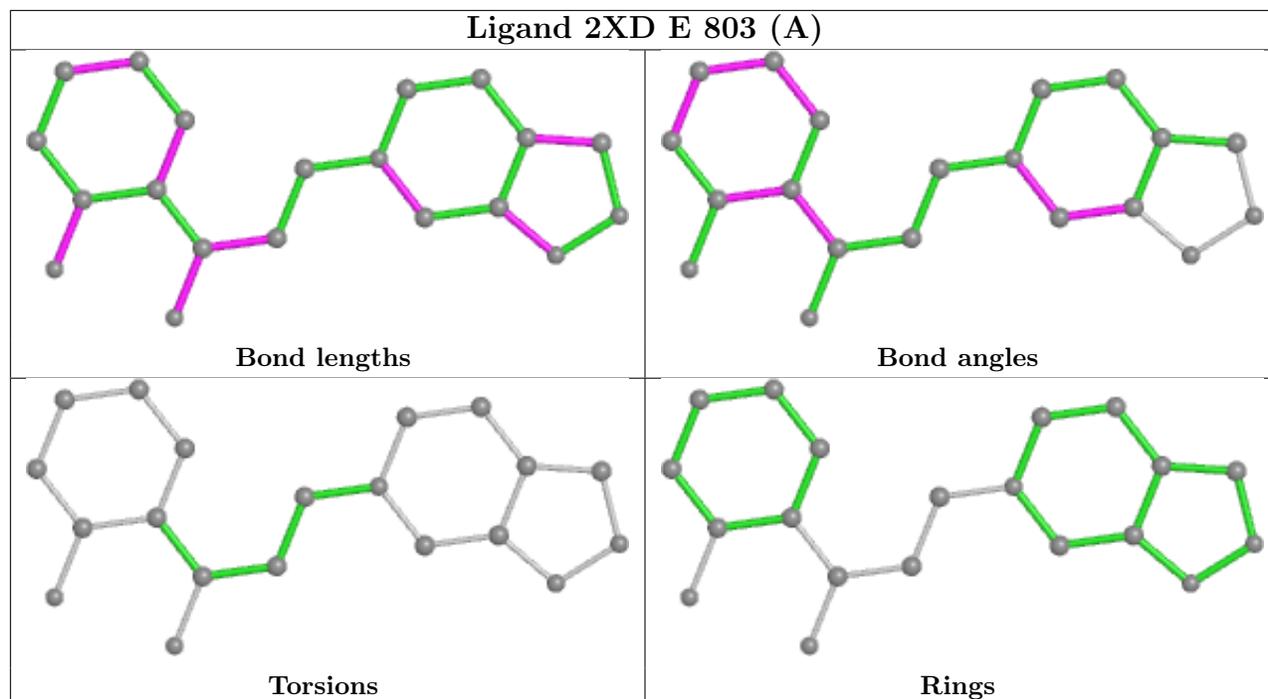
Mol	Chain	Res	Type	Atoms
3	C	802[A]	2XD	C7-C8-N-C14
2	E	802	EDO	O1-C1-C2-O2
2	A	801	EDO	O1-C1-C2-O2
2	F	801	EDO	O1-C1-C2-O2
2	H	801	EDO	O1-C1-C2-O2

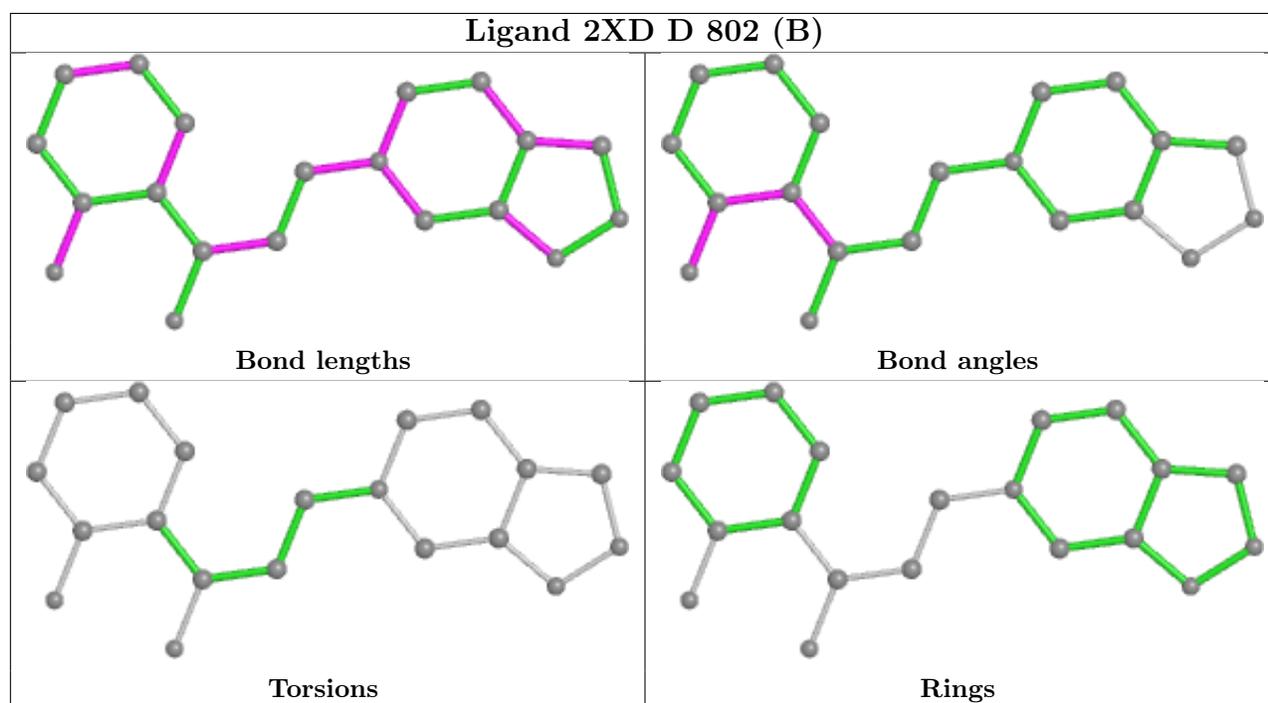
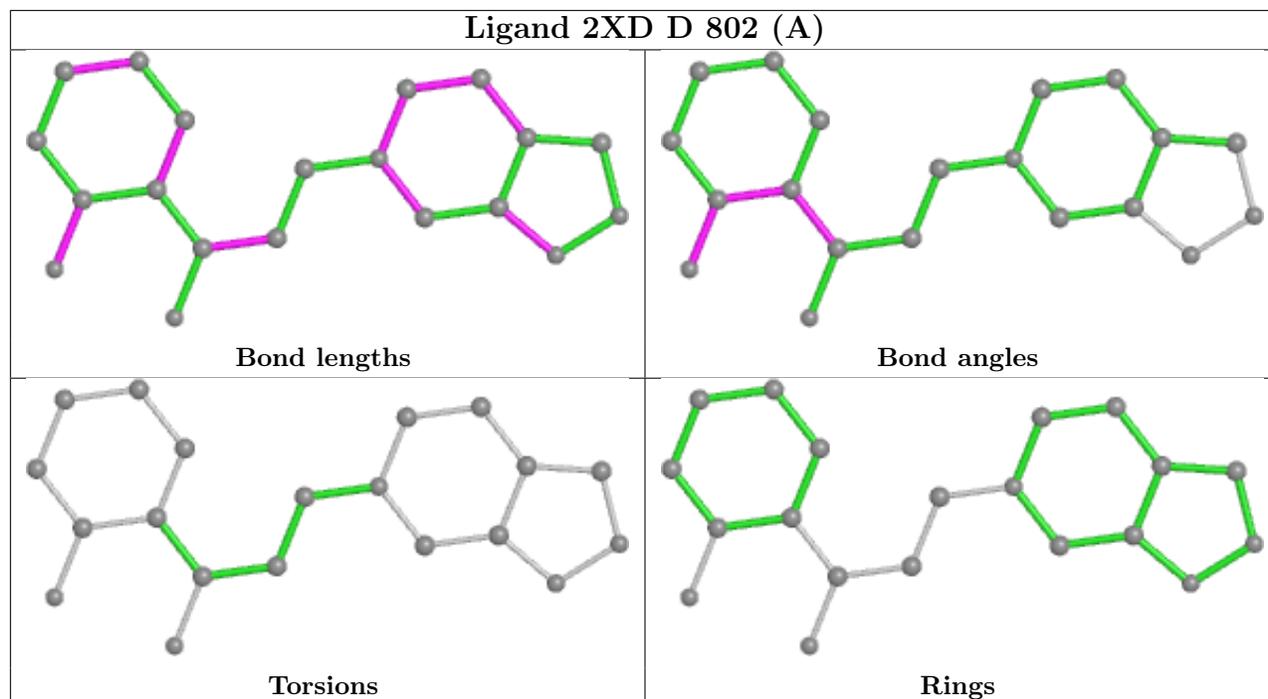
There are no ring outliers.

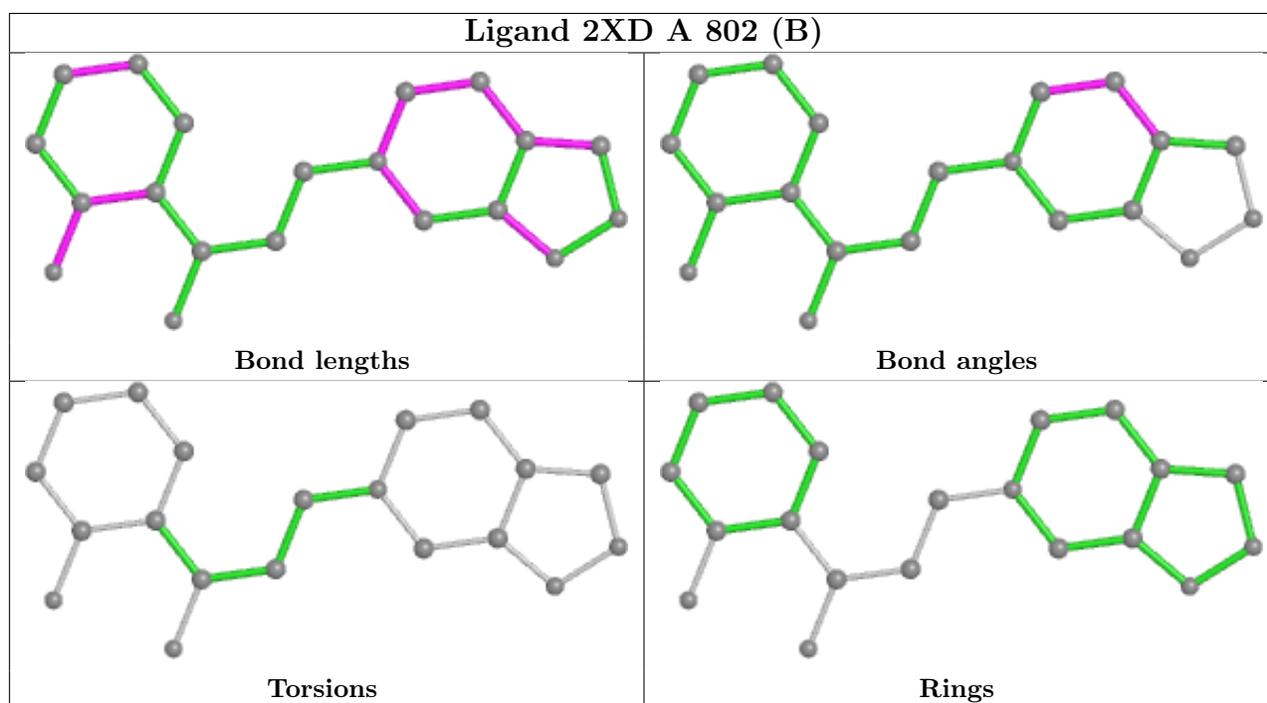
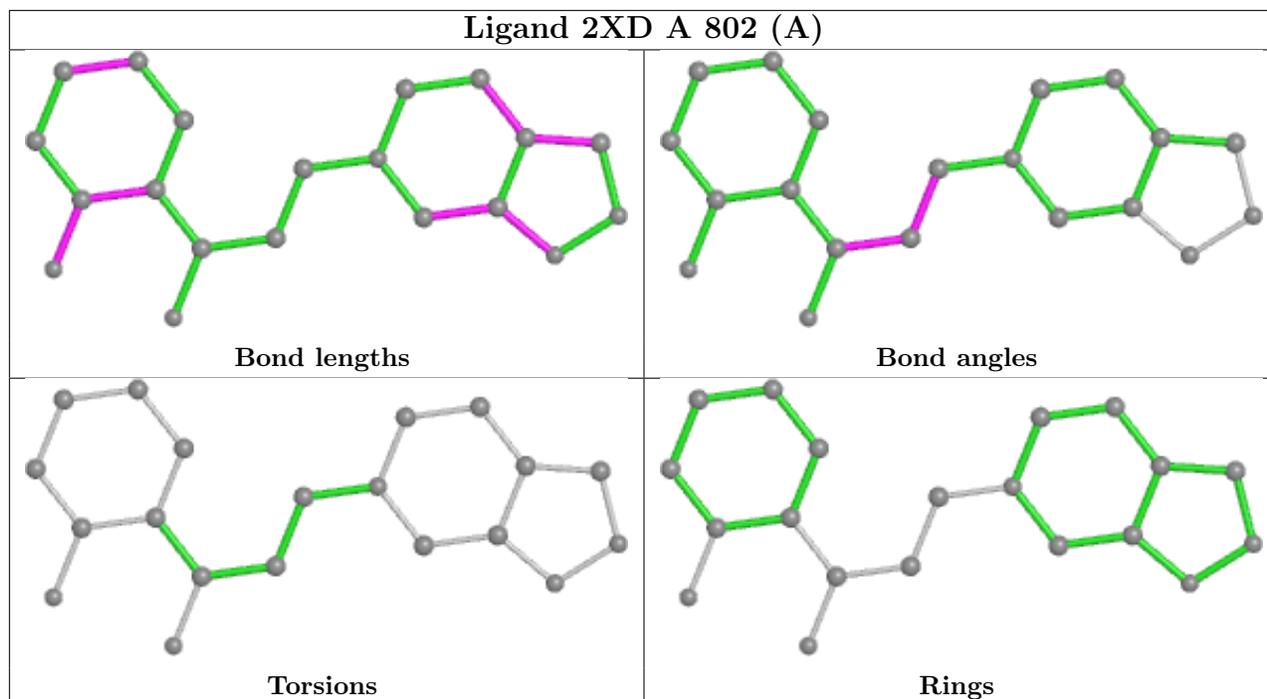
3 monomers are involved in 3 short contacts:

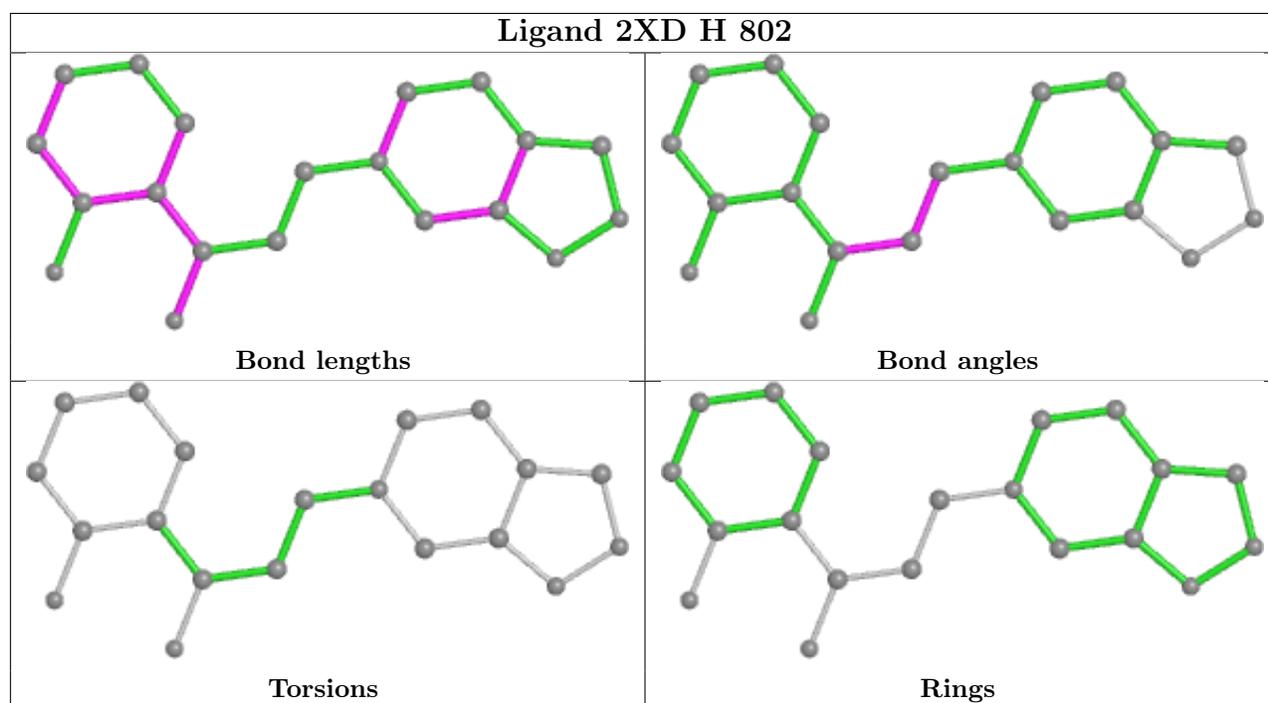
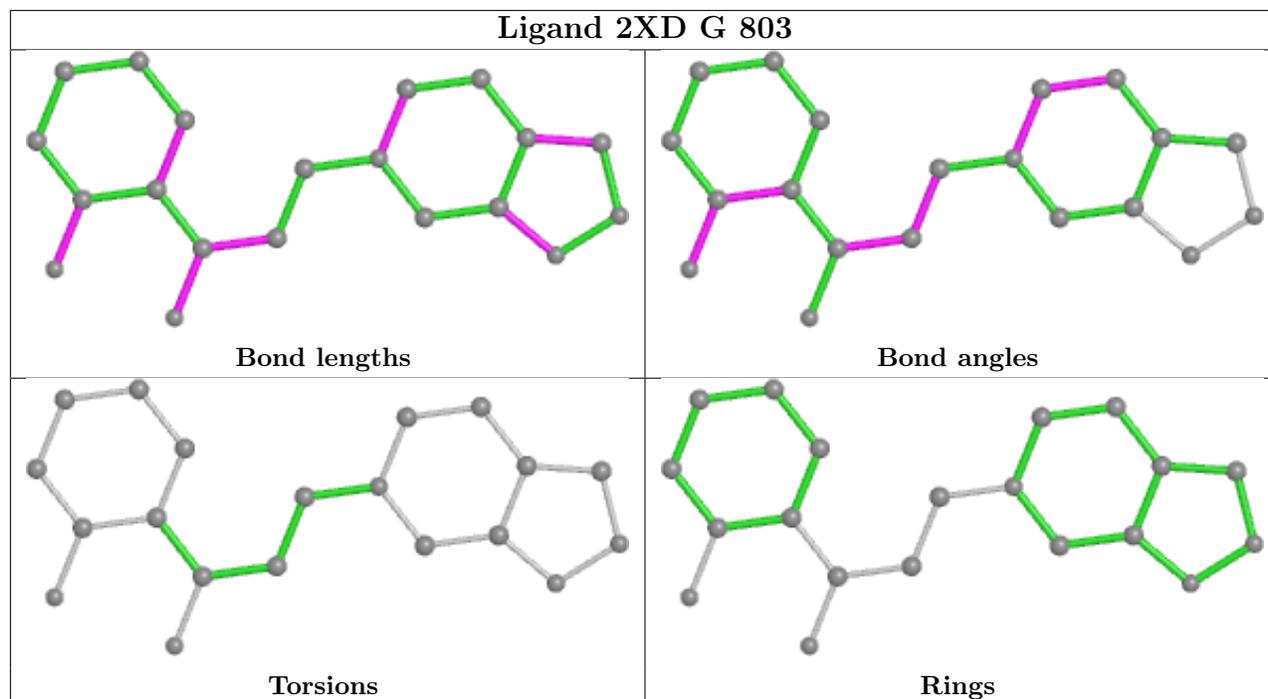
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	802[A]	2XD	1	0
2	E	802	EDO	1	0
3	A	802[B]	2XD	1	0

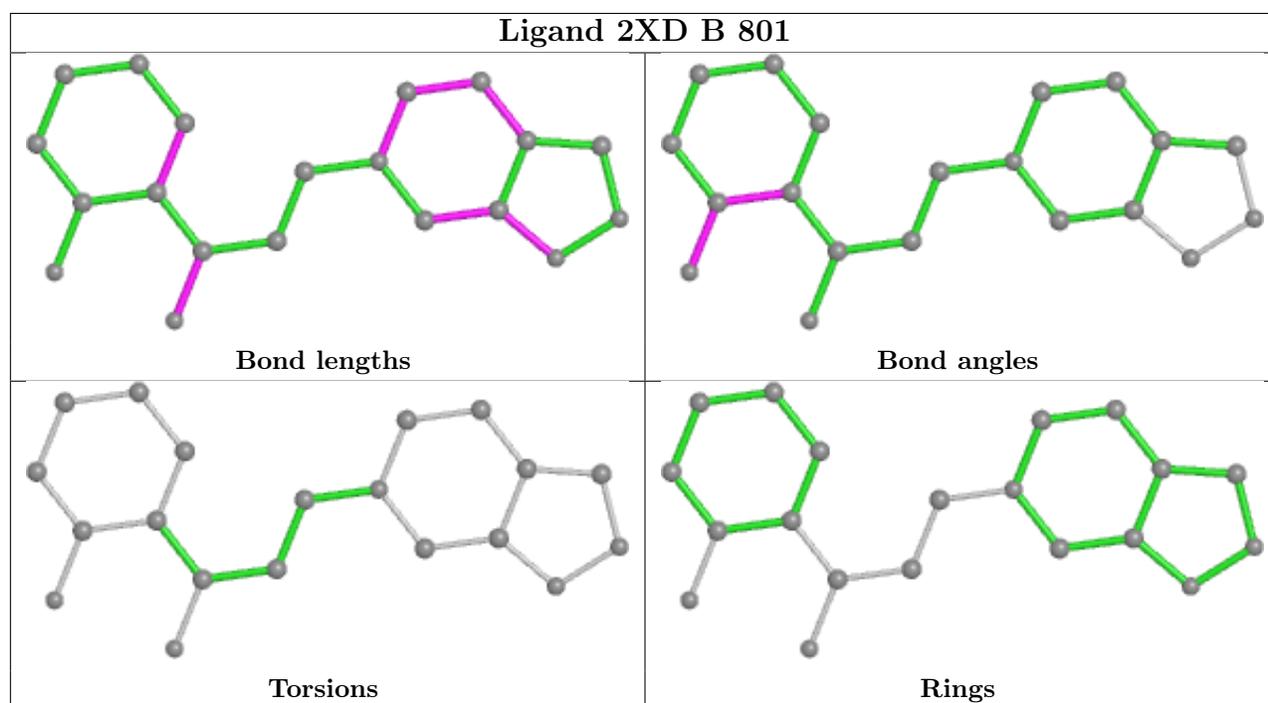
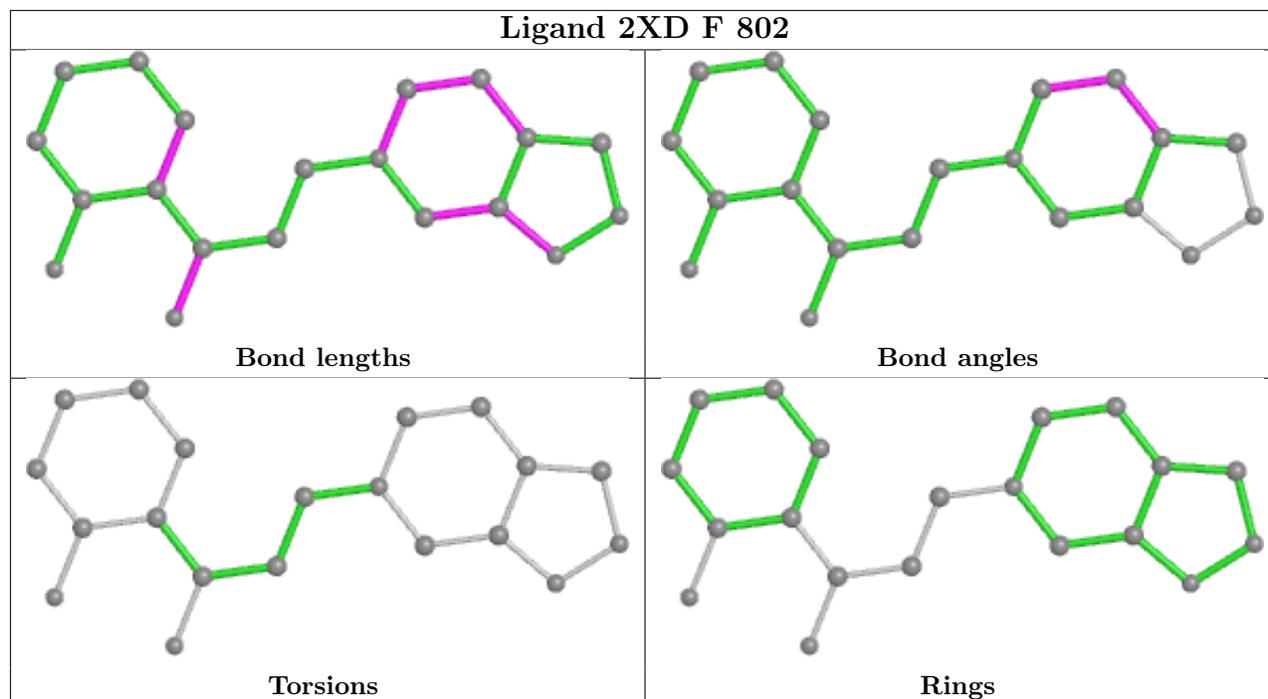
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

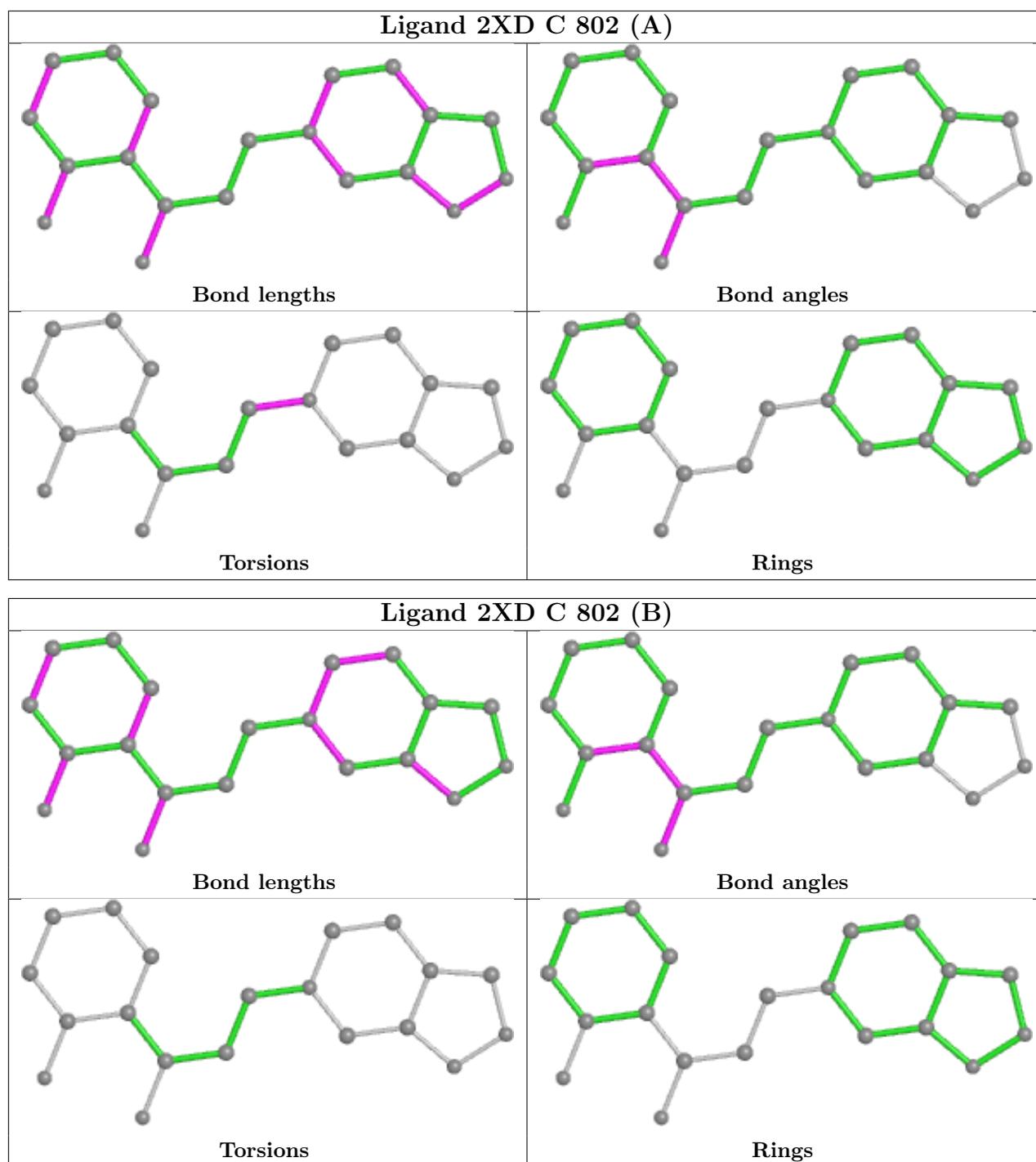












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	113/145 (77%)	-0.45	0 100 100	12, 27, 54, 72	0
1	B	111/145 (76%)	-0.44	1 (0%) 84 84	16, 29, 64, 72	0
1	C	111/145 (76%)	-0.43	1 (0%) 84 84	16, 29, 57, 75	0
1	D	112/145 (77%)	-0.47	0 100 100	14, 28, 60, 75	0
1	E	113/145 (77%)	-0.51	0 100 100	12, 25, 53, 70	0
1	F	112/145 (77%)	-0.51	0 100 100	15, 28, 56, 72	0
1	G	112/145 (77%)	-0.56	0 100 100	12, 24, 49, 66	0
1	H	111/145 (76%)	-0.57	0 100 100	13, 25, 47, 64	0
All	All	895/1160 (77%)	-0.49	2 (0%) 95 94	12, 27, 56, 75	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	622	TYR	2.3
1	C	731	LEU	2.2

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 monosaccharides in this entry.

6.4 Ligands

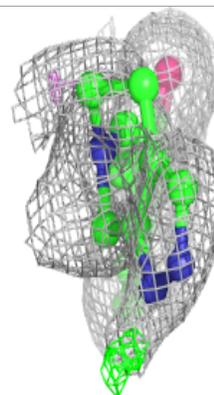
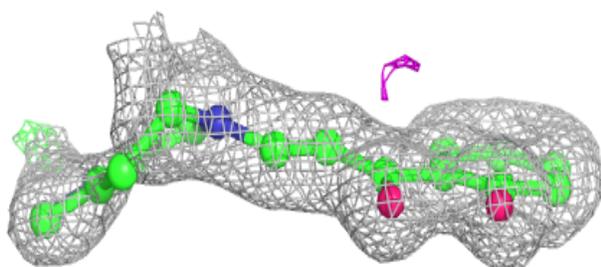
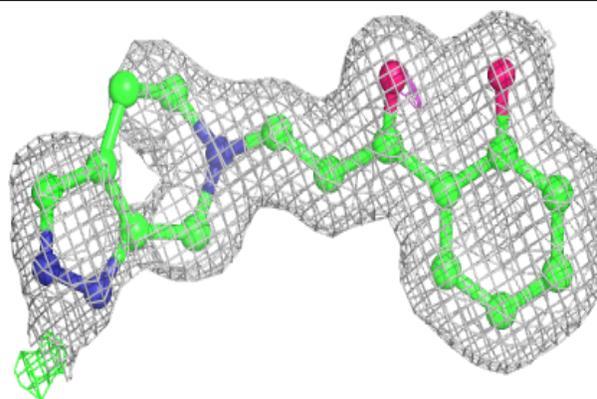
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	EDO	C	801	4/4	0.96	0.07	32,35,35,39	0
2	EDO	D	801	4/4	0.96	0.06	22,24,26,28	0
3	2XD	C	802[A]	20/20	0.96	0.09	19,28,44,50	10
3	2XD	C	802[B]	20/20	0.96	0.09	19,25,34,41	10
2	EDO	G	802	4/4	0.97	0.06	20,21,32,35	0
2	EDO	E	802	4/4	0.97	0.11	29,30,35,59	0
2	EDO	F	801	4/4	0.97	0.10	22,33,39,41	0
3	2XD	D	802[A]	20/20	0.97	0.08	13,20,30,32	10
3	2XD	D	802[B]	20/20	0.97	0.08	13,20,38,39	10
3	2XD	A	802[A]	20/20	0.98	0.08	13,17,24,34	10
3	2XD	A	802[B]	20/20	0.98	0.08	13,19,35,41	10
3	2XD	B	801	20/20	0.98	0.06	13,17,22,23	0
2	EDO	E	801	4/4	0.98	0.06	19,21,21,24	0
2	EDO	G	801	4/4	0.98	0.06	19,25,26,29	0
2	EDO	A	801	4/4	0.98	0.05	13,17,17,18	0
2	EDO	H	801	4/4	0.98	0.08	23,29,30,33	0
3	2XD	E	803[A]	20/20	0.98	0.08	12,16,24,24	10
3	2XD	E	803[B]	20/20	0.98	0.08	12,17,34,49	10
3	2XD	F	802	20/20	0.98	0.06	14,18,21,22	0
3	2XD	G	803	20/20	0.98	0.06	11,17,28,30	0
3	2XD	H	802	20/20	0.98	0.06	10,16,19,19	0

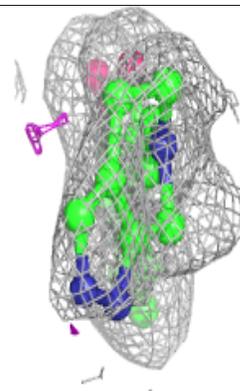
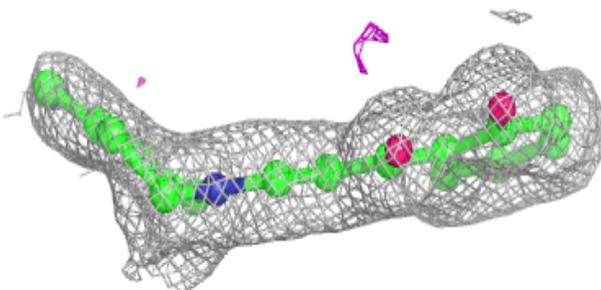
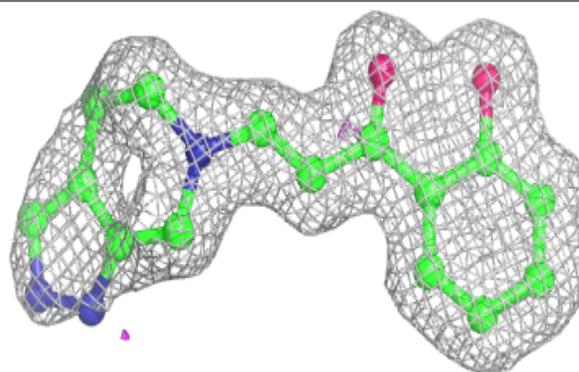
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 2XD C 802 (A):

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

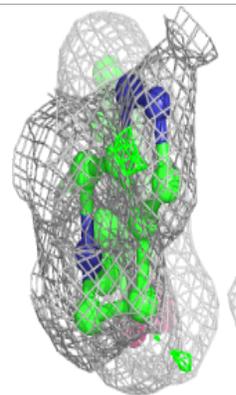
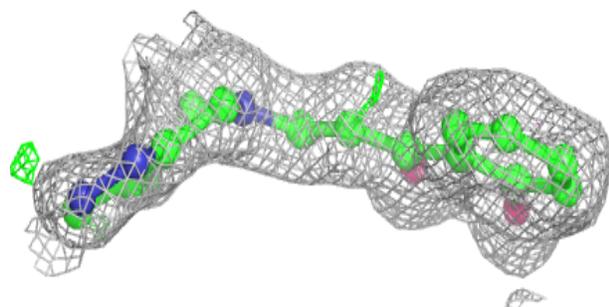
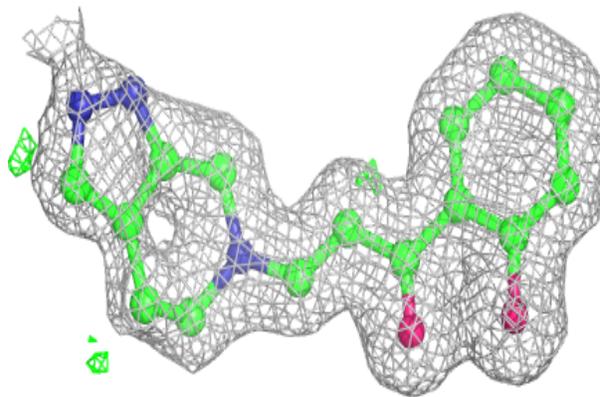
**Electron density around 2XD C 802 (B):**

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

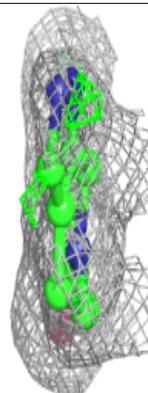
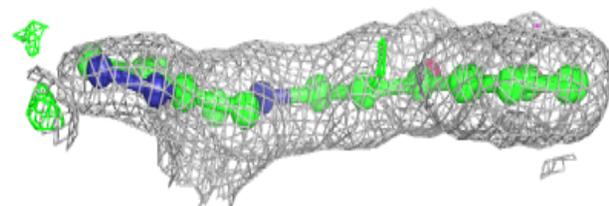
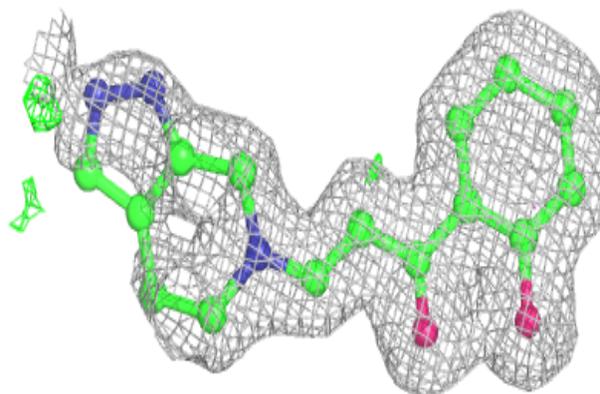


Electron density around 2XD D 802 (A):

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

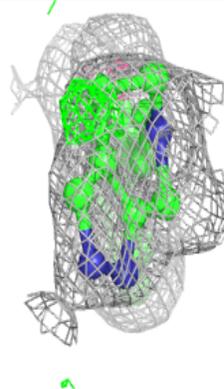
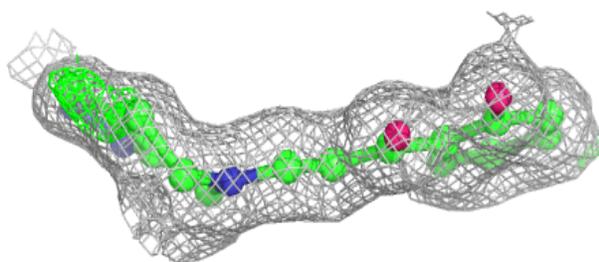
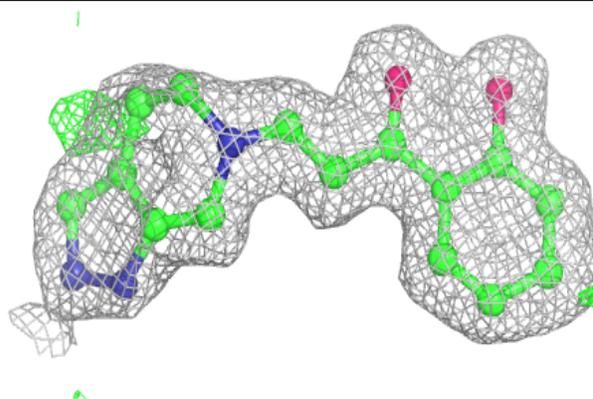
**Electron density around 2XD D 802 (B):**

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

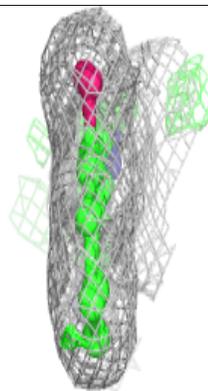
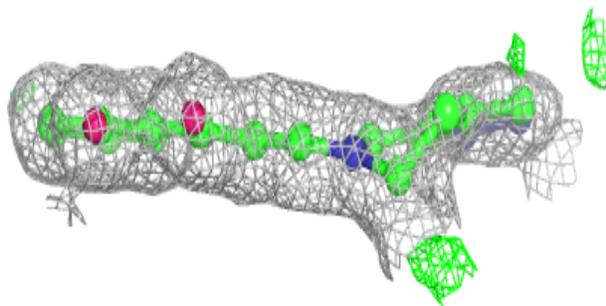
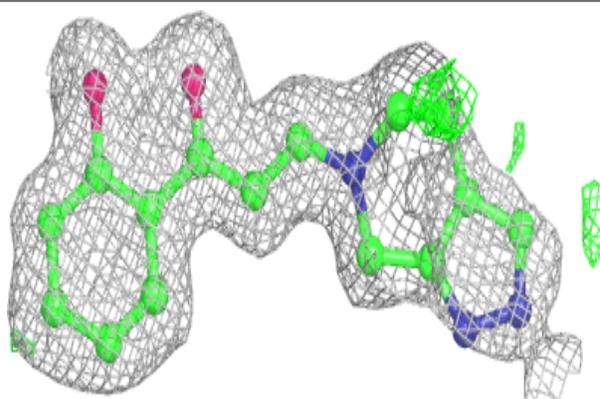


Electron density around 2XD A 802 (A):

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

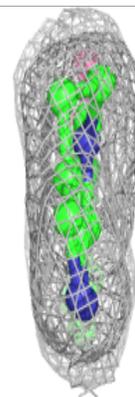
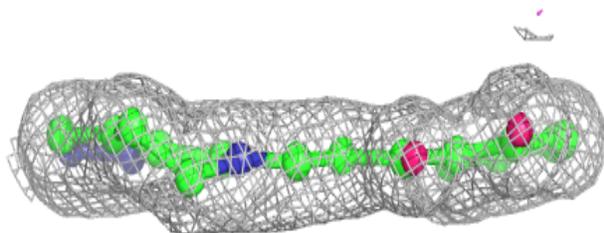
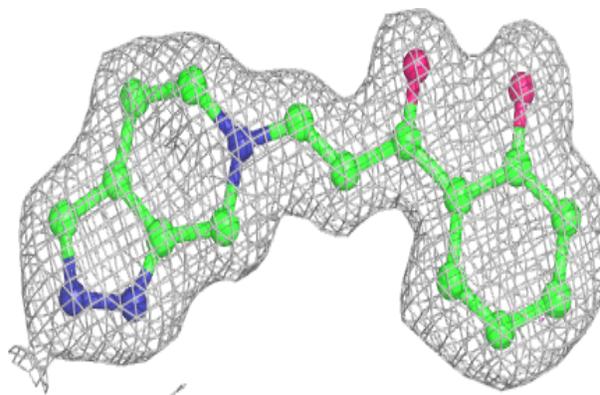
**Electron density around 2XD A 802 (B):**

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

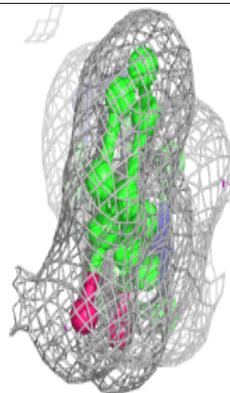
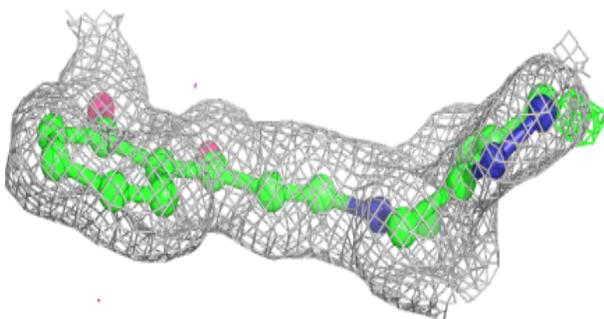
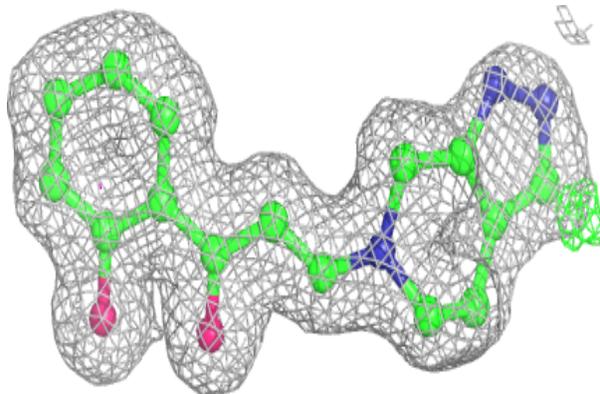


Electron density around 2XD B 801:

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

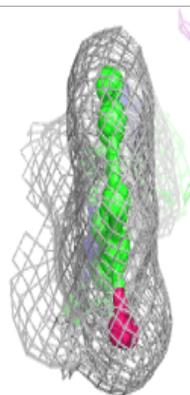
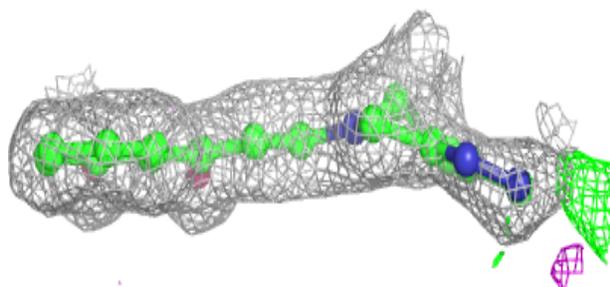
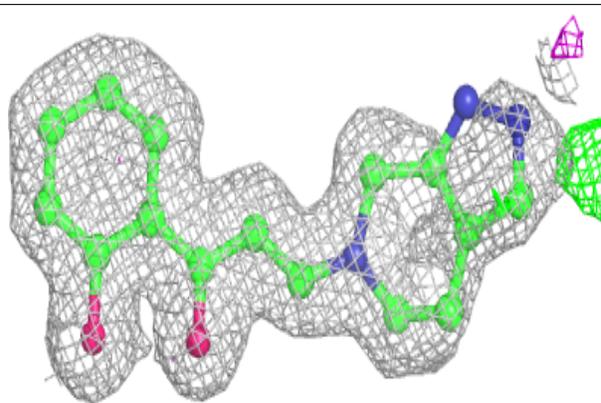
**Electron density around 2XD E 803 (A):**

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

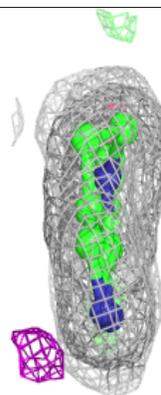
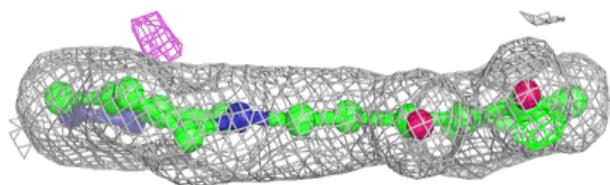
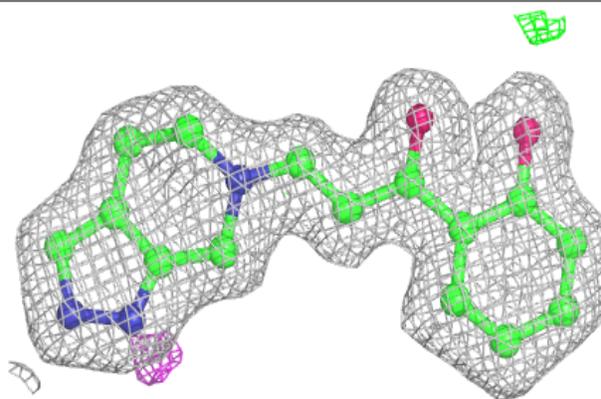


Electron density around 2XD E 803 (B):

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

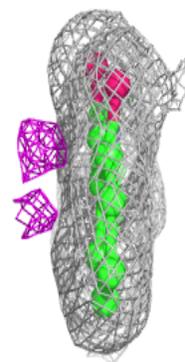
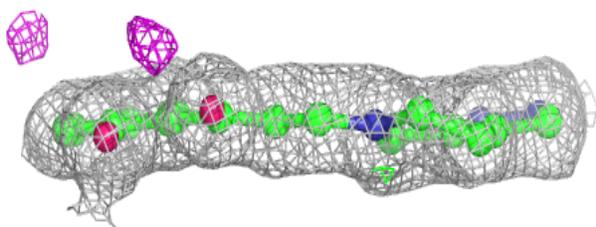
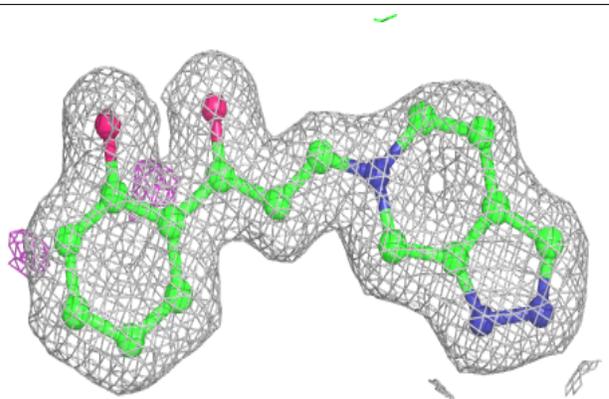
**Electron density around 2XD F 802:**

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

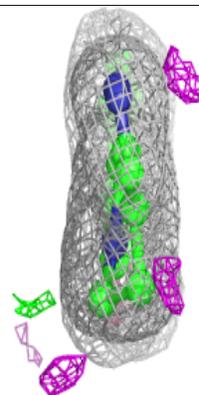
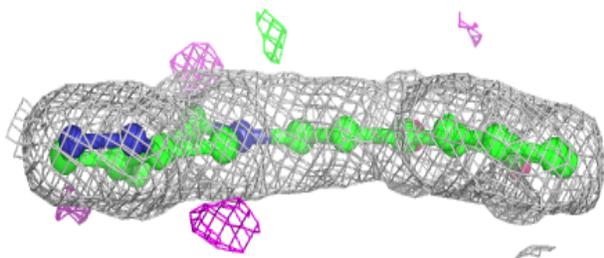
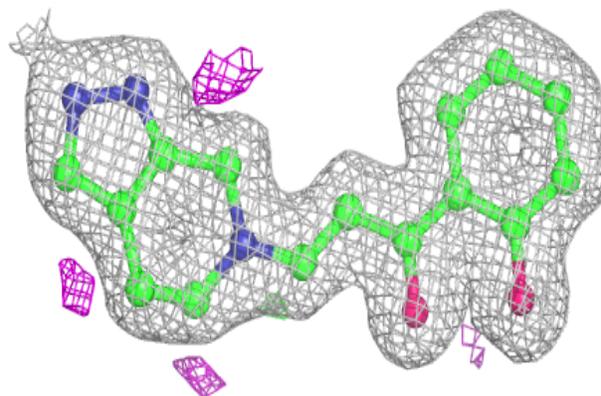


Electron density around 2XD G 803:

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

**Electron density around 2XD H 802:**

$2mF_o-DF_c$ (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.